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Final Report

Hot Mix Asphalt Plants Truck Loading and Silo Filling Manual Methods Testing

Asphalt Plant C Los Angeles, California

Volume 5 of 8



FINAL REPORT

**HOT MIX ASPHALT PLANTS
TRUCK LOADING AND SILO FILLING
MANUAL METHODS TESTING
ASPHALT PLANT C, LOS ANGELES, CALIFORNIA**

**VOLUME 5 OF 8
APPENDIX G.3**

**EPA Contract No. 68-D-98-004
Work Assignment No. 3-02**

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May 2000

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GLOSSARY OF TERMS

ASTM – American Society for Testing and Materials
 CEMS – Continuous Emissions Monitoring System
 CTS – Calibration Transfer Standard
 EMC – Emissions Measurement Center
 EMAD – Emission Monitoring and Analysis Division
 ESP – Electrostatic Precipitator
 FID – Flame Ionization Detector
 FTIR – Fourier Transform Infrared Spectroscopy
 HAP – Hazardous Air Pollutant
 MCEM – Methylene Chloride Extractable Matter
 MRI – Midwest Research Institute
 PES – Pacific Environmental Services
 PM – Particulate Matter
 PTE – Permanent Total Enclosure
 RAP – Recycled Asphalt
 RTFOT – Rolling Thin Film Oven Test
 SED – Silo Exhaust Duct

GLOSSARY OF TERMS (CONTINUED)

SMTG – Source Measurement Technology Group
SVOHAP – Semi-Volatile Organic Hazardous Air Pollutant
TED – Tunnel Emissions Duct
TFOT – Thin Film Oven Test
THC – Total Hydrocarbons
VOHAP – Volatile Organic Hazardous Air Pollutant
VOST – Volatile Organic Sampling Train

VOLUME 5A

APPENDIX G

ANALYTICAL DATA (CONTINUED)

G.3 SVOHAPS DATA

Organics Raw Data Package

Sample Data

Semivolatile Organics
Method 0010/8270

Client Name: Pacific Environmental Services
Client ID: S-MM5-2-F, FH, XAD, COND, BH
LAB ID: 300681-0001-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 25 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 100

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
Phenol	ND	ug/Sample	3000	GR
bis(2-Chloroethyl)ether	ND	ug/Sample	3000	
2-Chlorophenol	ND	ug/Sample	3000	
1,3-Dichlorobenzene	ND	ug/Sample	3000	
1,4-Dichlorobenzene	ND	ug/Sample	3000	
Benzyl alcohol	ND	ug/Sample	3000	
1,2-Dichlorobenzene	ND	ug/Sample	3000	
2-Methylphenol	ND	ug/Sample	3000	
2,2'-Oxybis(1-chloropropane)	ND	ug/Sample	3000	
3/4-Methylphenol	ND	ug/Sample	3000	
N-Nitroso-di-n-propylamine	ND	ug/Sample	3000	
Hexachloroethane	ND	ug/Sample	3000	
Nitrobenzene	ND	ug/Sample	3000	
Isophorone	ND	ug/Sample	3000	
2-Nitrophenol	ND	ug/Sample	3000	
2,4-Dimethylphenol	ND	ug/Sample	3000	
Benzoic acid	ND	ug/Sample	15000	
bis(2-Chloroethoxy)-methane	ND	ug/Sample	3000	
2,4-Dichlorophenol	ND	ug/Sample	3000	
1,2,4-Trichlorobenzene	ND	ug/Sample	3000	
Naphthalene	1600	ug/Sample	3000	J
4-Chloroaniline	ND	ug/Sample	3000	
Hexachlorobutadiene	ND	ug/Sample	3000	
4-Chloro-3-methylphenol	ND	ug/Sample	3000	
2-Methylnaphthalene	3600	ug/Sample	3000	
Hexachlorocyclopentadiene	ND	ug/Sample	3000	
2,4,6-Trichlorophenol	ND	ug/Sample	3000	
2,4,5-Trichlorophenol	ND	ug/Sample	15000	
2-Chloronaphthalene	ND	ug/Sample	3000	
2-Nitroaniline	ND	ug/Sample	3000	
Dimethyl phthalate	ND	ug/Sample	3000	
Acenaphthylene	ND	ug/Sample	3000	
3-Nitroaniline	ND	ug/Sample	15000	
Acenaphthene	ND	ug/Sample	3000	
2,4-Dinitrophenol	ND	ug/Sample	15000	

Note G = Reporting limit(s) raised due to matrix interference.

Note J = Result is detected below the reporting limit or is an estimated concentration.

Note R = Reporting limit(s) raised due to sample volume limitations.

ND = Not Detected

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Approved By: Mike Orbanosky

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Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: S-MM5-2-F, FH, XAD, COND, BH
LAB ID: 300681-0001-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 25 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 100

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
4-Nitrophenol	ND	ug/Sample	15000	
Dibenzofuran	ND	ug/Sample	3000	
2,4-Dinitrotoluene	ND	ug/Sample	3000	
2,6-Dinitrotoluene	ND	ug/Sample	3000	
Diethyl phthalate	ND	ug/Sample	3000	
4-Chlorophenyl phenyl ether	ND	ug/Sample	3000	
Fluorene	ND	ug/Sample	3000	
4-Nitroaniline	ND	ug/Sample	15000	
4,6-Dinitro-2-methylphenol	ND	ug/Sample	15000	
N-Nitrosodiphenylamine	ND	ug/Sample	3000	
4-Bromophenyl phenyl ether	ND	ug/Sample	3000	
Hexachlorobenzene	ND	ug/Sample	3000	
Pentachlorophenol	ND	ug/Sample	15000	
Phenanthrene	950	ug/Sample	3000	J
Anthracene	ND	ug/Sample	3000	
Di-n-butyl phthalate	ND	ug/Sample	3000	
Fluoranthene	ND	ug/Sample	3000	
Pyrene	ND	ug/Sample	3000	
Butyl benzyl phthalate	ND	ug/Sample	3000	
3,3'-Dichlorobenzidine	ND	ug/Sample	6000	
Benzo(a)anthracene	ND	ug/Sample	3000	
bis(2-Ethylhexyl)-phthalate	ND	ug/Sample	3000	
Chrysene	ND	ug/Sample	3000	
Di-n-octyl phthalate	ND	ug/Sample	3000	
Benzo(b)fluoranthene	ND	ug/Sample	3000	
Benzo(k)fluoranthene	ND	ug/Sample	3000	
Benzo(a)pyrene	ND	ug/Sample	3000	
Indeno(1,2,3-cd)pyrene	ND	ug/Sample	3000	
Dibenz(a,h)anthracene	ND	ug/Sample	3000	
Benzo(g,h,i)perylene	ND	ug/Sample	3000	
Acetophenone	ND	ug/Sample	3000	
4-Aminobiphenyl	ND	ug/Sample	15000	
Aniline	ND	ug/Sample	3000	
Benzdine	ND	ug/Sample	30000	
3,3'-Dimethylbenzidine	ND	ug/Sample	6000	
N-Nitrosodimethylamine	ND	ug/Sample	3000	
N-Nitrosomorpholine	ND	ug/Sample	3000	

Note J = Result is detected below the reporting limit or is an estimated concentration.
ND = Not Detected

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Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: S-MM5-2-F, FH, XAD, COND, BH
LAB ID: 300681-0001-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 25 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 100

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
Pentachloronitrobenzene (PCNB)	ND	ug/Sample	15000	
o-Toluidine	ND	ug/Sample	6000	
2-Methoxybenzenamine	ND	ug/Sample	--	
Biphenyl	ND	ug/Sample	--	
Chloroacetophenone	ND	ug/Sample	--	
Cumene	ND	ug/Sample	--	
DBCP (1,2-Dibromo-3-chloropropane)	ND	ug/Sample	--	
Benzo(e)pyrene	ND	ug/Sample	--	
N-N-Diethylaniline	ND	ug/Sample	--	
Dimethylaniline	ND	ug/Sample	--	
3,3'-Dimethoxybenzidine	ND	ug/Sample	--	
Hydroquinone	ND	ug/Sample	--	
4,4'-Methyl-bis(2-chloroaniline)	ND	ug/Sample	--	
4-Nitrodiphenyl	ND	ug/Sample	--	
Trifluralin	ND	ug/Sample	--	

Surrogate	Recovery	Acceptable Range
Nitrobenzene-d5	ND %	45 - 107
2-Fluorobiphenyl	ND %	62 - 110
Terphenyl-d14	ND %	58 - 135
Phenol-d5	ND %	43 - 130
2-Fluorophenol	ND %	36 - 111
2,4,6-Tribromophenol	ND %	58 - 131

ND = Not Detected

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Semivolatiles Library Search (20 Compound TID)
Method 8270

Client Name: Pacific Environmental Services
Client ID: S-MM5-2-F,FH,XAD,COND,BH
LAB ID: 300681-0001-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 25 JUL 98
Prepared: NA

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 100

Parameter	Result	Units	Reporting Limit	Qualifier
Unknown alkane	7900	ug/Sample	--	
Unknown	13000	ug/Sample	--	
Unknown PAH	15000	ug/Sample	--	
Unknown alkene	8900	ug/Sample	--	
Undecane	8600	ug/Sample	--	0
Unknown	8400	ug/Sample	--	
Dodecane, 2,6,10-trimethyl-	17000	ug/Sample	--	0
Tetradecane	22000	ug/Sample	--	0
Unknown alkene	12000	ug/Sample	--	
Heptadecane, 2,6,10,14 -tetramethyl-	18000	ug/Sample	--	0
Undecane, 2-methyl-	19000	ug/Sample	--	0
Unknown alkane	13000	ug/Sample	--	
Nonadecane	24000	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	15000	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	32000	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	10000	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	14000	ug/Sample	--	0
Nonadecane	7700	ug/Sample	--	0
Heptadecane, 2,6,10,14 -tetramethyl-	9700	ug/Sample	--	0
Nonadecane	7500	ug/Sample	--	0

Note 0 = Or structurally similar compound (isomer).
NA = Not Applicable

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QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068101
Std Id: ST16980831

Sample: S-MM5-2 1/3SA/100M INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 19:00
Analyst: DAT Instrument ID: F16 Run Factor: 300.
Quan List Threshold: 0.95 Surrogate Vol.: 100.

Surrogate Spike Recoveries
8270-G, LIMS 10/10/96

Surrogate	Surrogate Amount (ug)		% Recovery		QC limits
	Spiked	Measured	Measured		
CS20 NITROBENZENE-D5	50.0	BDL	0*	45	107
CS25 2-FLUOROBIPHENY	50.0	BDL	0*	62	110
CS30 TERPHENYL-D14	50.00	135.0	270.	*	58 135
CS45 PHENOL-D5	100.	BDL	0*	43	130
CS50 2-FLUOROPHENOL	100.	BDL	0*	36	111
CS55 2,4,6-TRIBROMOP	100.0	120.0	120.	58	131

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
HEXACHLOROBENZENE-C13		ND	UG/A	3000.0
C310 N-NITROSODIMETHYLAMINE		ND	UG/A	3000.0
PYRIDINE		ND	UG/A	6000.0
2-PICOLINE		ND	UG/A	3000.0
N-NITROSOMETHYLETHYLAMINE		ND	UG/A	3000.0
METHYLMETHANESULFONATE		ND	UG/A	3000.0
N-NITROSODIETHYLAMINE		ND	UG/A	3000.0
ETHYLMETHANE SULFONATE		ND	UG/A	3000.0
PENTACHLOROETHANE		ND	UG/A	3000.0
C320 ANILINE		ND	UG/A	3000.0
C315 PHENOL		ND	UG/A	3000.0
C325 BIS(2-CHLOROETHYL)ETHER		ND	UG/A	3000.0
C330 2-CHLOROPHENOL		ND	UG/A	3000.0
C335 1,3-DICHLOROBENZENE		ND	UG/A	3000.0
C340 1,4-DICHLOROBENZENE		ND	UG/A	3000.0
C345 BENZYL ALCOHOL		ND	UG/A	3000.0
C350 1,2-DICHLOROBENZENE		ND	UG/A	3000.0
C355 2-METHYLPHENOL		ND	UG/A	3000.0
C360 2,2'-OXYBIS(1-CLPROPAN)		ND	UG/A	3000.0
C361 ACETOPHENONE		ND	UG/A	3000.0
N-NITROSOPYRROLIDINE		ND	UG/A	3000.0
N-NITROSOMORPHOLINE		ND	UG/A	3000.0
3-METHYL PHENOL		ND	UG/A	3000.0
C365 4-METHYLPHENOL		ND	UG/A	3000.0
C370 N-NITROSO-DI-N-PROPYLAM		ND	UG/A	3000.0
O-TOLUIDINE		ND	UG/A	3000.0
C375 HEXACHLOROETHANE		ND	UG/A	3000.0
C410 NITROBENZENE		ND	UG/A	3000.0
N-NITROSOPIPERIDINE		ND	UG/A	3000.0

Reviewed by:

MM 9.18.98
DEG 10/6/98

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068101
Std Id: ST16980831

Sample: S-MM5-2 1/3SA/100M INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 19:00
Analyst: DAT Instrument ID: F16 Run Factor: 300.
Quan List Threshold: 0.95 Surrogate Vol.: 100.

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C415 ISOPHORONE		ND	UG/A	3000.0
C420 2-NITROPHENOL		ND	UG/A	3000.0
C425 2,4-DIMETHYLPHENOL		ND	UG/A	3000.0
C435 BIS(2-CHLOROETHOXY)METH		ND	UG/A	3000.0
C440 2,4-DICHLOROPHENOL		ND	UG/A	3000.0
C430 BENZOIC ACID		ND	UG/A	15000.0
AA-DIMETHYLPHENETHYLAMINE		ND	UG/A	3000.0
C445 1,2,4-TRICHLOROBENZENE		ND	UG/A	3000.0
C450 NAPHTHALENE		ND	UG/A	3000.0
C455 4-CHLOROANALINE		ND	UG/A	3000.0
2,6-DICHLOROPHENOL		ND	UG/A	3000.0
HEXACHLOROPROPENE		ND	UG/A	3000.0
C460 HEXACHLOROBUTADIENE		ND	UG/A	3000.0
P-PHENYLENE DIAMINE		ND	UG/A	3000.0
N-NITROSODI-N-BUTYLAMINE		ND	UG/A	3000.0
C465 4-CHLORO-3-METHYLPHENO		ND	UG/A	3000.0
SAFROLE		ND	UG/A	3000.0
C470 2-METHYLNAPHTHALENE	652	3600.	UG/A	3000.0
1,2,4,5-TETRACHLOROBENZENE		ND	UG/A	3000.0
ISOSAFROLE (#1)		ND	UG/A	6000.0
C510 HEXACHLOROCYCLOPENTADI		ND	UG/A	3000.0
C515 2,4,6-TRICHLOROPHENOL		ND	UG/A	3000.0
C520 2,4,5-TRICHLOROPHENOL		ND	UG/A	3000.0
ISOSAFROLE (#2)		ND	UG/A	6000.0
C525 2-CHLORONAPHTHALENE		ND	UG/A	3000.0
1-CHLORONAPHTHALENE		ND	UG/A	1500.0
C530 2-NITROANALINE		ND	UG/A	15000.0
1,4-NAPHTHOQUINONE		ND	UG/A	3000.0
C535 DIMETHYLPHTHALATE		ND	UG/A	3000.0
1,3-DINITROBENZENE		ND	UG/A	3000.0
C540 ACENAPHTHYLENE		ND	UG/A	3000.0
C543 2,6-DINITROTOLUENE		ND	UG/A	3000.0
C545 3-NITROANILINE		ND	UG/A	15000.0
C550 ACENAPHTHENE		ND	UG/A	3000.0
C555 2,4-DINITROPHENOL		ND	UG/A	15000.0
C565 DIBENZOFURAN		ND	UG/A	3000.0
C560 4-NITROPHENOL		ND	UG/A	15000.0
PENTACHLOROBENZENE		ND	UG/A	3000.0
C570 2,4-DINITROTOLUENE		ND	UG/A	3000.0
1-NAPHTHYLAMINE		ND	UG/A	3000.0
2-NAPHTHYLAMINE		ND	UG/A	3000.0
2,3,4,6-TETRACHLOROPHENOL		ND	UG/A	6000.0
C580 DIETHYLPHTHALATE		ND	UG/A	3000.0
C590 FLUORENE		ND	UG/A	3000.0

ND 470
< MOL

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068101
Std Id: ST16980831

Sample: S-MM5-2 1/3SA/100M INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 19:00
Analyst: DAT Instrument ID: F16 Run Factor: 300.
Quan List Threshold: 0.95 Surrogate Vol.: 100.

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C585 4-CHLOROPHENYL-PHENYLE		ND	UG/A	3000.0
5-NITRO-O-TOLUIDINE		ND	UG/A	3000.0
C595 4-NITROANALINE		ND	UG/A	15000.0
C610 4,6-DINITRO-2-METHYLPH		ND	UG/A	15000.0
C615 N-NITROSODIPHENYLAMINE		ND	UG/A	3000.0
C620 AZOBENZENE		ND	UG/A	3000.0
SYM-TRINITROBENZENE		ND	UG/A	3000.0
C625 4-BROMOPHENYL-PHENYLET		ND	UG/A	3000.0
PHENACETIN		ND	UG/A	3000.0
DIALATE	234	ND	UG/A	3000.0
C630 HEXACHLOROBENZENE		ND	UG/A	3000.0
4-AMINOBIIPHENYL		ND	UG/A	3000.0
C635 PENTACHLOROPHENOL		ND	UG/A	15000.0
PRONAMIDE		ND	UG/A	3000.0
PENTACHLORONITROBENZENE		ND	UG/A	15000.0
C640 PHENANTHRENE		ND 950	UG/A	3000.0
C645 ANTHRACENE		ND	UG/A	3000.0
2SECBUTYL-4,6-DINITROPHENOL		ND	UG/A	3000.0
C647 CARBAZOLE		ND	UG/A	3000.0
C650 DI-N-BUTYLPHTHALATE		ND	UG/A	3000.0
4-NITROQUINOLINE-1-OXIDE		ND	UG/A	3000.0
METHAPYRILENE		ND	UG/A	3000.0
ISODRIN		ND	UG/A	3000.0
C655 FLUORANTHENE		ND	UG/A	3000.0
CHLOROBENZILATE		ND	UG/A	3000.0
C710 BENZIDINE		ND	UG/A	30000.0
C715 PYRENE		ND	UG/A	3000.0
ARAMITE (#1)		ND	UG/A	3000.0
ARAMITE (#2)		ND	UG/A	3000.0
P-DIMETHYLAMINOAZOBENZENE		ND	UG/A	3000.0
3,3'-DIMETHYLBENZIDINE		ND	UG/A	3000.0
KEPONE		ND	UG/A	15000.0
C720 BUTYLBENZYLPHTHALATE		ND	UG/A	3000.0
2-ACETYLAMINOFLUORENE		ND	UG/A	3000.0
C730 BENZO(A)ANTHRACENE		ND	UG/A	3000.0
C725 3,3'-DICHLOBENZIDINE		ND	UG/A	6000.0
C740 CHRYSENE		ND	UG/A	3000.0
C745 BIS(2-ETHYLHEXYL)PHTHA		ND	UG/A	3000.0
3-METHYLCHOLANTHRENE		ND	UG/A	3000.0
C760 DI-N-OCTYL PHTHALATE		ND	UG/A	3000.0
C765 BENZO(B)FLUORANTHENE		ND	UG/A	3000.0
7,12-DIMETHYLBENZANTHRACENE		ND	UG/A	3000.0
C770 BENZO(K)FLUORANTHENE		ND	UG/A	3000.0
HEXACHLOROPHENE		ND	UG/A	3000.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

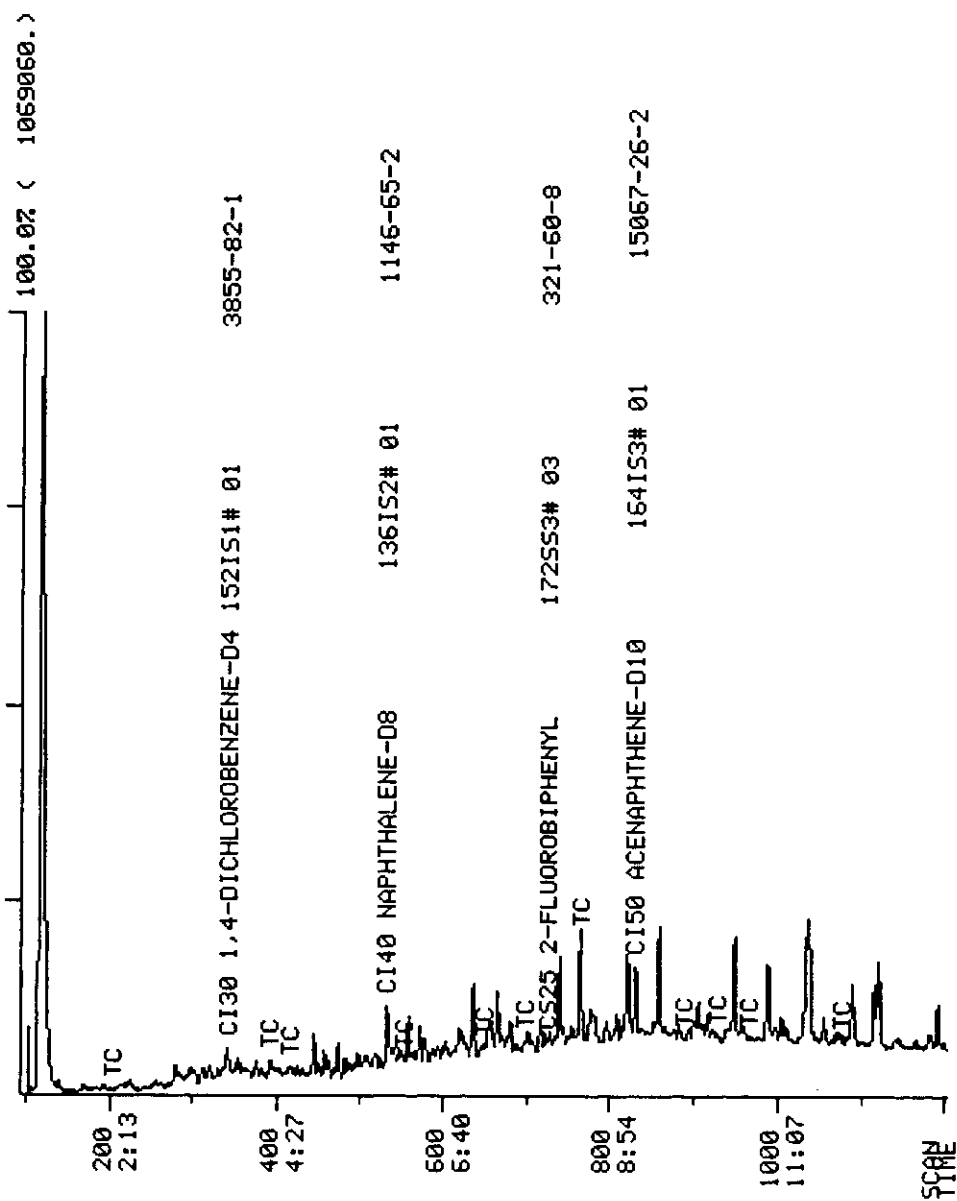
Data File: 30068101
Std Id: ST16980831

Sample: S-MM5-2 1/3SA/100M INST. ID: F16
Client: PACIF1 Date Analyzed: 08/31/98 19:00
Analyst: DAT Instrument ID: F16 Run Factor: 300.
Quan List Threshold: 0.95 Surrogate Vol.: 100.

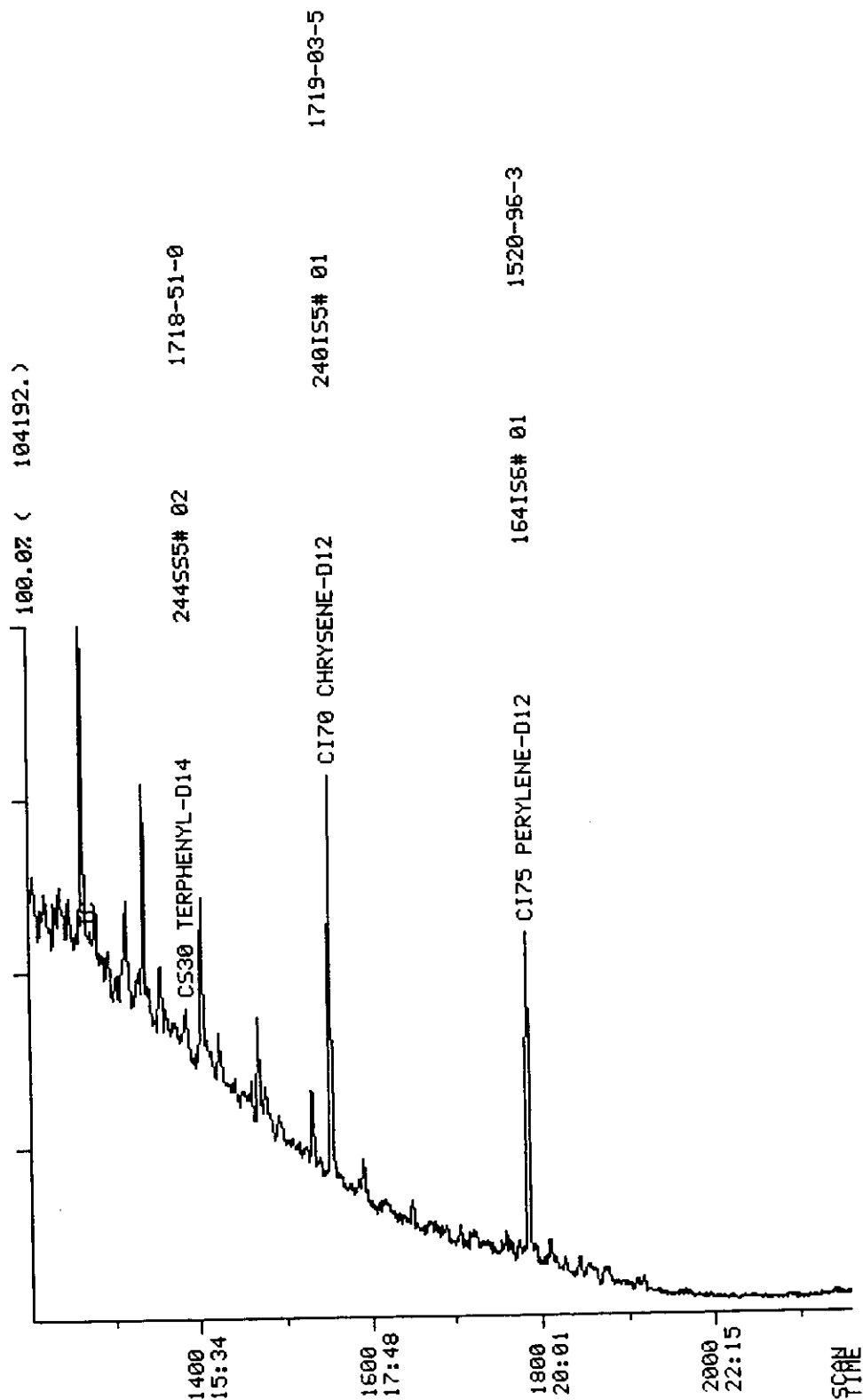
Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C775 BENZO(A)PYRENE		ND	UG/A	3000.0
C780 INDENO(1,2,3-CD)PYRENE		ND	UG/A	3000.0
C785 DIBENZ(A,H)ANTHRACENE		ND	UG/A	3000.0
C790 BENZO(G,H,I)PERYLENE		ND	UG/A	3000.0

DATA FROM FILE: 30068101 SCANS 93 TO 1203 ACQUIRED: 08/31/98 19:00:00
 SAMPLE: S-MM5-2 1/3SA/100M CALI: 30068101 #3
 COND.: UC/ML *100ML *1002/1002 *(NA/NA >)/1/3SA NA M INST. ID: F16



DATA FROM FILE: 30068101 SCANS 1203 TO 2158 ACQUIRED: 08/31/98 19:00:00
 CALI: 30068101 #3
 SAMPLE: S-MM5-2 1/35A/100M INST. ID: F16
 CONDS.: UG/ML *100ML *100%/100% *(NA/NA) >1/35A NA M



ADDITIONAL STANDARD USED FOR THIS REPORT IS: ST16980831A

QUANTERRA QUANTITATION SUMMARY

File: 30068101

Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Analyst: DAT Instrument Id: F16 Analyzed: 08/31/98 19:00

Compounds with amounts less than 0.20 reported as NOT FOUND

Library							Units: UG/ML
No Entry	Name	Mass Meth	Scan Ref	Area RRF(L)	Amount		
1 S1#	1 C130 1,4-DICHLOROBENZENE	152 A BV	341 1	25865. 1.000	40.000		
2 S2#	1 C140 NAPHTHALENE-D8	136 A BB	530 2	96183. 1.000	40.000		
3 S3#	1 C150 ACENAPHTHENE-D10	164 A BB	829 3	53596. 1.000	40.000		
4 S4#	1 C160 PHENANTHRENE-D10	188 A BB	1087 4	95242. 1.000	40.000		
5 S5#	1 C170 CHRYSENE-D12	240 A BB	1552 5	57183. 1.000	40.000		
6 S6#	1 C175 PERYLENE-D12	264 A BB	1783 6	44948. 1.000	40.000		
7 S2#	2 CS20 NITROBENZENE-D5	82 A BB	424 2	355. 0.470	0.314		
8 S3#	3 CS25 2-FLUOROBIPHENYL	172 A BB	723 3	1150. 1.234	0.695		
9 S5#	2 CS30 TERPHENYL-D14	244 A BB	1384 5	1924. 0.995	1.352		
10 S1#	3 CS45 PHENOL-D5	99	1 NOT FOUND				
11 S1#	2 CS50 2-FLUOROPHENOL	112	1 NOT FOUND				
12 S3#	2 CS55 2,4,6-TRIBROMOPHENO	330 A BB	971 3	380. 0.236	1.203		
13 S4#	4 HEXACHLOROBENZENE-C13	294	4 NOT FOUND				
14 S1#	15 C310 N-NITROSODIMETHYLAM	74	1 NOT FOUND				
15 S1#	20 PYRIDINE	79	1 NOT FOUND				
16 S1#	30 2-PICOLINE	93	1 NOT FOUND				
17 S1#	40 N-NITROSOMETHYLETHYLAMIN	42 A VB	180 1	405. 0.811	0.772		
18 S1#	60 METHYLMETHANESULFONATE	80 A BB	201 1	448. 0.636	1.089	73.5 3267	
19 S1#	70 N-NITROSODIETHYLAMINE	102	1 NOT FOUND				
20 S1#	85 ETHYLMETHANE SULFONATE	79	1 NOT FOUND				
21 S1#	95 PENTACHLOROETHANE	117	1 NOT FOUND				
22 S1#	100 C320 ANILINE	93	1 NOT FOUND				
23 S1#	105 C315 PHENOL	94	1 NOT FOUND				
24 S1#	110 C325 BIS(2-CHLOROETHYL)E	93 A BB	319 1	216. 1.368	0.244		
25 S1#	115 C330 2-CHLOROPHENOL	128	1 NOT FOUND				
26 S1#	125 C335 1,3-DICHLOROBENZENE	146	1 NOT FOUND				
27 S1#	130 C340 1,4-DICHLOROBENZENE	146	1 NOT FOUND				
28 S1#	145 C345 BENZYL ALCOHOL	108 A VB	374 1	268. 0.854	0.485		
29 S1#	150 C350 1,2-DICHLOROBENZENE	146	1 NOT FOUND				
30 S1#	160 C355 2-METHYLPHENOL	108 A BV	392 1	1488. 1.175	1.959		
31 S1#	165 C360 2,2'-OXYBIS(1-CLPRO	45 A BB	394 1	1758. 2.539	1.071		
32 S1#	170 C361 ACETOPHENONE	105	1 NOT FOUND				
33 S1#	175 N-NITROSOPYRROLIDINE	100	1 NOT FOUND				
34 S1#	180 N-NITROSOMORPHOLINE	56	1 NOT FOUND				
35 S1#	182 3-METHYL PHENOL	108 A VB	423 1	363. 2.252	0.249		
36 S1#	185 C365 4-METHYLPHENOL	108 A VV	419 1	452. 1.276	0.548		
37 S1#	190 C370 N-NITROSO-DI-N-PROP	70 A BB	414 1	1123. 0.906	1.917		
38 S1#	195 O-TOLUIDINE	106 A BB	413 1	332. 2.160	0.238		
39 S1#	200 C375 HEXACHLOROETHANE	117	1 NOT FOUND				
40 S2#	10 C410 NITROBENZENE	77	2 NOT FOUND				
41 S2#	15 N-NITROSOPIPERIDINE	42	2 NOT FOUND				
42 S2#	20 C415 ISOPHORONE	82	2 NOT FOUND				
43 S2#	25 C420 2-NITROPHENOL	139	2 NOT FOUND				
44 S2#	30 C425 2,4-DIMETHYLPHENOL	107 A BB	500 2	714. 0.330	0.900		
45 S2#	35 C435 BIS(2-CHLOROETHOXY)	93	2 NOT FOUND				

46	S2# 40	C440	2,4-DICHLOROPHENOL	162	A BB	519	2	200.	0.308	0.270
47	S2# 45	C430	BENZOIC ACID	122		2	NOT FOUND			
48	S2# 50	AA-DIMETHYLPHENETHYLAMIN	58			2	NOT FOUND			
49	S2# 55	C445	1,2,4-TRICHLOROBENZ	180		2	NOT FOUND			

50 S2# 60	C450 NAPHTHALENE	128 A BV	533 2	13016. 1.022	5.295
51 S2# 80	C455 4-CHLOROANALINE	127 A BB	556 2	389. 0.454	0.354
52 S2# 85	2,6-DICHLOROPHENOL	162 A BV	552 2	1164. 0.257	1.880
53 S2# 90	HEXACHLOROPROPENE	213	2 NOT FOUND		
54 S2# 95	C460 HEXACHLOROBUTADIENE	225	2 NOT FOUND		
55 S2#115	P-PHENYLENE DIAMINE	108	2 NOT FOUND		
56 S2#120	N-NITROSODI-N-BUTYLAMINE	84	2 NOT FOUND		
57 S2#130	C465 4-CHLORO-3-METHYLPH	107	2 NOT FOUND		
58 S2#140	SAFROLE	162 A BB	648 2	852. 0.248	1.431
59 S2#145	C470 2-METHYLNAPHTHALENE	142 A BB	652 2	21609. 0.751	11.962
60 S3# 10	1,2,4,5-TETRACHLOROBENZE	214	3 NOT FOUND		
61 S3# 15	ISOSAFROLE (#1)	162 A BB	697 3	418. 0.044	7.156
62 S3# 20	C510 HEXACHLOROCYCLOPENT	237	3 NOT FOUND		
63 S3# 25	C515 2,4,6-TRICHLOROPHEN	196	3 NOT FOUND		
64 S3# 30	C520 2,4,5-TRICHLOROPHEN	196	3 NOT FOUND		
65 S3# 35	ISOSAFROLE (#2)	104 A VB	738 3	394. 0.195	1.508
66 S3# 40	C525 2-CHLORONAPHTHALENE	162	3 NOT FOUND		
67 S3# 42	1-CHLORONAPHTHALENE	162	3 NOT FOUND		
68 S3# 45	C530 2-NITROANALINE	65 A BB	764 3	755. 0.502	1.122
69 S3# 50	1,4-NAPHTHOQUINONE	158 A VB	772 3	450. 0.415	0.810
70 S3# 55	C535 DIMETHYLPHTHALATE	163	3 NOT FOUND		
71 S3# 60	1,3-DINITROBENZENE	168	3 NOT FOUND		
72 S3# 65	C540 ACENAPHTHYLENE	152 A BB	808 3	533. 1.838	0.216
73 S3# 70	C543 2,6-DINITROTOLUENE	165 A BB	816 3	267. 0.324	0.615
74 S3# 75	C545 3-NITROANILINE	138	3 NOT FOUND		
75 S3# 80	C550 ACENAPHTHENE	153 A BV	834 3	943. 1.159	0.608
76 S3# 85	C555 2,4-DINITROPHENOL	184 A BB	857 3	70. 0.189	0.276
77 S3# 90	C565 DIBENZOFURAN	168 A BB	867 3	838. 1.669	0.375
78 S3# 95	C560 4-NITROPHENOL	109 A BV	887 3	343. 0.173	1.483
79 S3#100	PENTACHLOROBENZENE	250	3 NOT FOUND		
80 S3#105	C570 2,4-DINITROTOLUENE	165 A BB	881 3	483. 0.397	0.908
81 S3#110	1-NAPHTHYLAMINE	143	3 NOT FOUND		
82 S3#115	2-NAPHTHYLAMINE	143	3 NOT FOUND		
83 S3#120	2,3,4,6-TETRACHLOROPHENO	232 A*BB	896 3	118. 0.297	0.297
84 S3#130	C580 DIETHYLPHTHALATE	149	3 NOT FOUND		
85 S3#135	C590 FLUORENE	166 A BB	926 3	2574. 1.298	1.480
86 S3#140	C585 4-CHLOROPHENYL-PHEN	204	3 NOT FOUND		
87 S3#145	5-NITRO-O-TOLUIDINE	152 A BB	941 3	2148. 0.356	4.509
88 S3#150	C595 4-NITROANALINE	138 A BB	948 3	373. 0.307	0.907
89 S4# 10	C610 4,6-DINITRO-2-METHY	198	4 NOT FOUND		
90 S4# 15	C615 N-NITROSODIPHENYLAM	169 A BB	963 4	2520. 0.551	1.922
91 S4# 20	C620 AZOBENZENE	77	4 NOT FOUND		
92 S4# 25	SYM-TRINITROBENZENE	75	4 NOT FOUND		
93 S4# 30	C625 4-BROMOPHENYL-PHENY	248 A BB	1022 4	320. 0.222	0.604
94 S4# 35	PHENACETIN	108	4 NOT FOUND		
95 S4# 37	DIALATE	234	4 NOT FOUND		
96 S4# 40	C630 HEXACHLOROBENZENE	284	4 NOT FOUND		
97 S4# 45	4-AMINOBIIPHENYL	169 A BB	1076 4	1557. 0.609	1.074
98 S4# 50	C635 PENTACHLOROPHENOL	266 A VB	1063 4	156. 0.155	0.423
99 S4# 55	PRONAMIDE	173	4 NOT FOUND		
100 S4# 60	PENTACHLORONITROBENZENE	237 A BB	1064 4	54. 0.080	0.282
101 S4# 65	C640 PHENANTHRENE	178 A BB	1091 4	7792. 1.033	3.169
102 S4# 70	C645 ANTHRACENE	178	4 NOT FOUND		
103 S4# 75	2SECBUTYL-4,6-DINITROPHE	211	4 NOT FOUND		
104 S4# 80	C647 CARBAZOLE	167 A BB	1148 4	364. 0.764	0.200
105 S4# 85	C650 DI-N-BUTYLPHTHALATE	149	4 NOT FOUND		
106 S4#100	4-NITROQUINOLINE-1-OXIDE	190	4 NOT FOUND		
107 S4#105	METHAPYRILENE	58	4 NOT FOUND		
108 S4#106	ISODRIN	193 A BB	1270 4	504. 0.127	1.671
109 S4#110	C655 FLUORANTHENE	202	4 NOT FOUND		
110 S4#120	CHLOROBENZILATE	139	4 NOT FOUND		

1588

11.962

1.480

950.7

111 S5# 10 C710 BENZIDINE	184	5	NOT FOUND		
112 S5# 15 C715 PYRENE	202 A BB 1335	5	1175.	1.323	0.622
113 S5# 20 ARAMITE (#1)	185	5	NOT FOUND		
114 S5# 25 ARAMITE (#2)	185	5	NOT FOUND		
115 S5# 30 P-DIMETHYLAMINOAZOBENZEN	120	5	NOT FOUND		
116 S5# 35 3,3'-DIMETHYLBENZIDINE	212 A BV 1472	5	120.	0.418	0.201
117 S5# 37 KEPONE	272	5	NOT FOUND		
118 S5# 40 C720 BUTYLBENZYLPHTHALAT	149	5	NOT FOUND		
119 S5# 45 2-ACETYLAMINOFLUORENE	181	5	NOT FOUND		
120 S5# 50 C730 BENZO(A)ANTHRACENE	228	5	NOT FOUND		
121 S5# 55 C725 3,3'-DICHLOROBENZID	252	5	NOT FOUND		
122 S5# 60 C740 CHRYSENE	228 A VB 1554	5	825.	1.014	0.569
123 S5# 65 C745 BIS(2-ETHYLHEXYL)PH	149	5	NOT FOUND		
124 S5# 85 3-METHYLCHOLANTHRENE	268	5	NOT FOUND		
125 S6# 10 C760 DI-N-OCTYL PHTHALAT	149	6	NOT FOUND		
126 S6# 15 C765 BENZO(B)FLUORANTHEN	252	6	NOT FOUND		
127 S6# 20 7,12-DIMETHYLBENZANTHRAC	256	6	NOT FOUND		
128 S6# 25 C770 BENZO(K)FLUORANTHEN	252	6	NOT FOUND		
129 S6# 30 HEXACHLOROPHENE	196	6	NOT FOUND		
130 S6# 35 C775 BENZO(A)PYRENE	252	6	NOT FOUND		
131 S6# 55 C780 INDENO(1,2,3-CD)PYR	276	6	NOT FOUND		
132 S6# 60 C785 DIBENZ(A,H)ANTHRACE	278	6	NOT FOUND		
133 S6# 65 C790 BENZO(G,H,I)PERYLEN	276	6	NOT FOUND		

QUANTERRA Internal Standard Check

Standard Filename: ST16980831

Analyzed: 08/31/98 12:31

Standard	IS# 1			IS# 2			IS# 3		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	28649.		3.78	100432.		5.90	54347.		9.21
Upper Limit	57297.=+200		4.28	200864.=+200		6.40	108695.=+200		9.71
Lower Limit	14324.=+ 50		3.28	50216.=+ 50		5.40	27174.=+ 50		8.71
Filename									
1 30068101	25865.= 90		3.79	96183.= 95		5.90	53596.= 98		9.22

IS# 1 = C130 1,4-DICHLOROBENZENE-D4

IS# 2 = C140 NAPHTHALENE-D8

IS# 3 = C150 ACENAPHTHENE-D10

* - indicates an I.S. is outside QC limit(s)

Standard	IS# 4			IS# 5			IS# 6		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	98520.		12.08	62289.		17.26	50750.		19.83
Upper Limit	197039.=+200		12.58	124578.=+200		17.76	101500.=+200		20.33
Lower Limit	49260.=+ 50		11.58	31145.=+ 50		16.76	25375.=+ 50		19.33
Filename									
1 30068101	95242.= 96		12.09	57183.= 91		17.26	44948.= 88		19.83

IS# 4 = C160 PHENANTHRENE-D10

IS# 5 = C170 CHRYSENE-D12

IS# 6 = C175 PERYLENE-D12

* - indicates an I.S. is outside QC limit(s)

TARGET COMPOUND COMPARISON

COMPOUND: C450 NAPHTHALENE

128 S2# 50

91-20-3

RAW DATA: 30068101 #533

08/31/98 19:00

SAMPLE: S-MM5-2 1/35A/100M

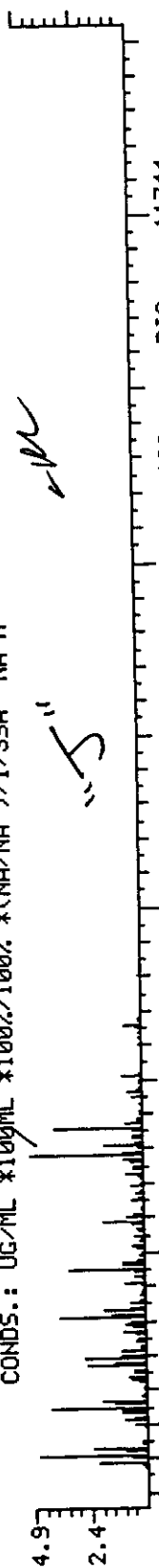
CONDS.: UG/ML *100%100% *(NA/NA) /1/35A NA M

INST. ID: F16

BASE M/Z: 128

RIC: 68096.

3854.

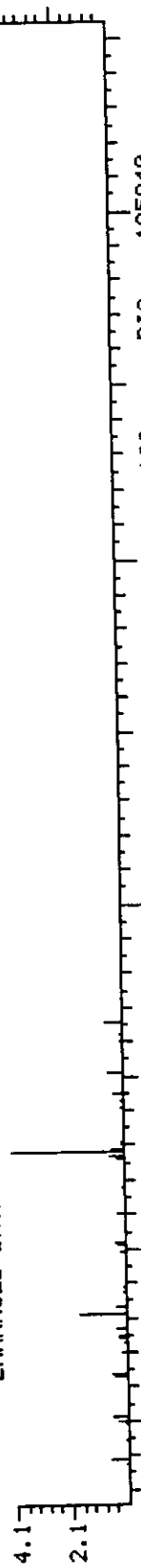


ENHANCED DATA: 30068101 #533

BASE M/Z: 128

RIC: 11744.

3254.



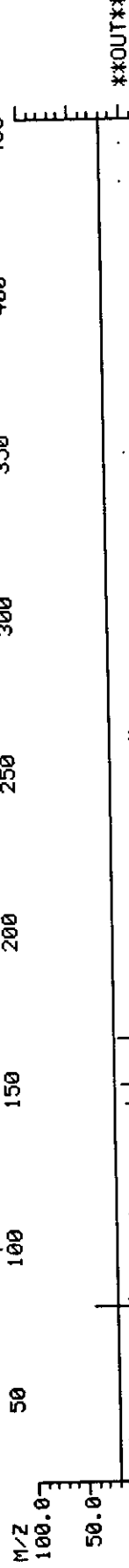
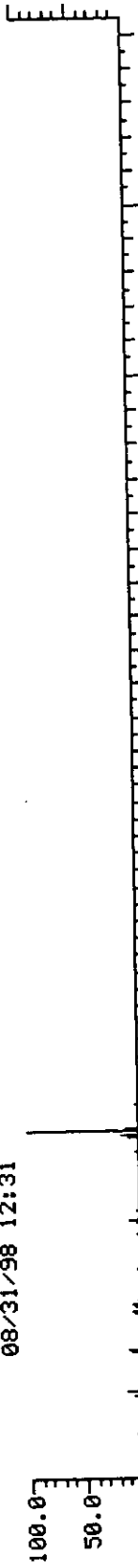
STANDARD FILE: ST16980831 #533

08/31/98 12:31

BASE M/Z: 128

RIC: 195840.

79232.



OUT

TARGET COMPOUND COMPARISON

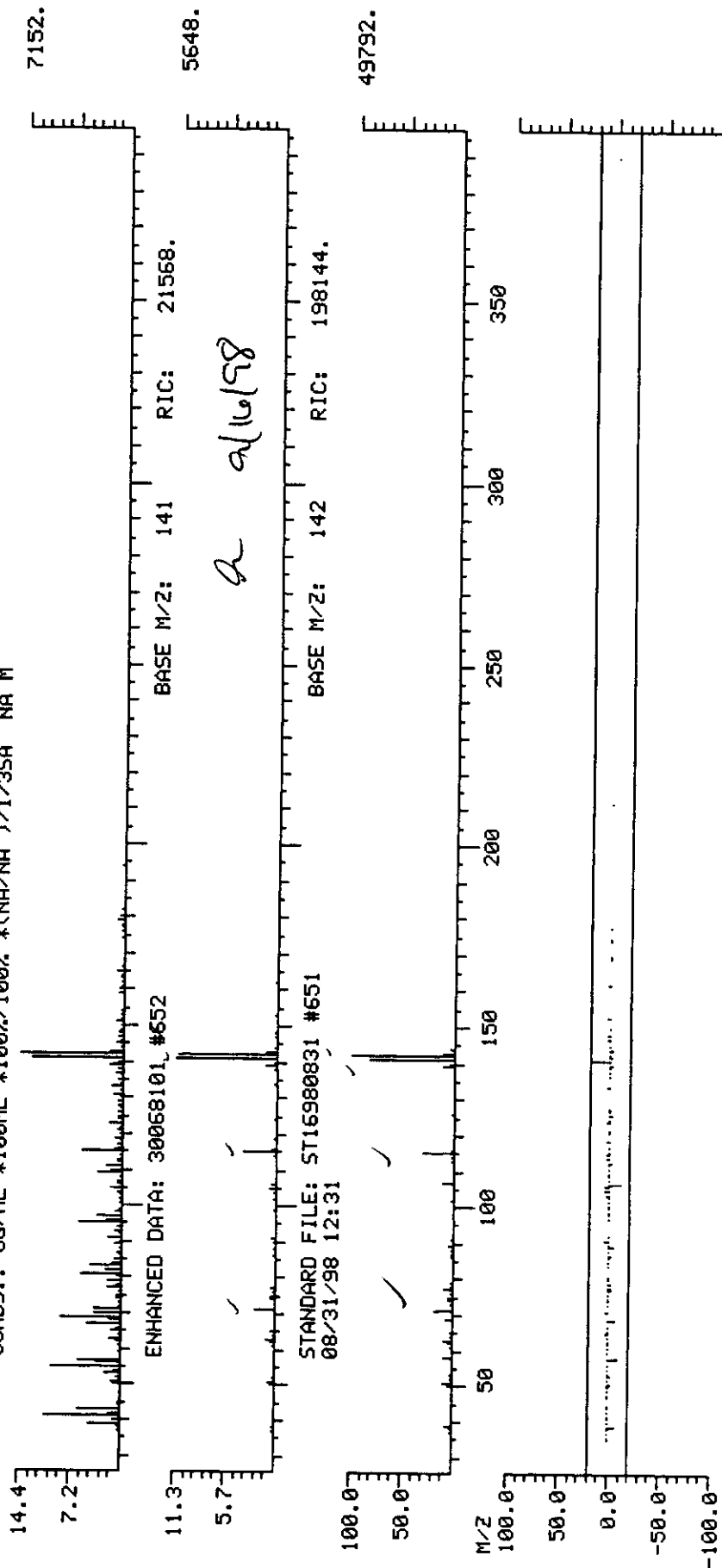
COMPOUND: C470 2-METHYLNAPHTHALENE 142 52#145 91-57-6

RAW DATA: 30068101 #552 BASE M/Z: 142 RIC: 97280.

08/31/98 19:00

SAMPLE: 5-MMS-2 1/35A/100M INST. ID: F16

CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M



TARGET COMPOUND COMPARISON

COMPOUND: C640 PHENANTHRENE

178 S4# 65

85-01-8

RAW DATA: 30068101 #1091

BASE M/Z: 41 RIC: 87680.

08/31/98 19:00

SAMPLE: S-MMS-2 1/35A/100M INST. ID: F16

COND.: UG/ML *100ML *100Z/100Z *(NA/NA) 1/35A NA M

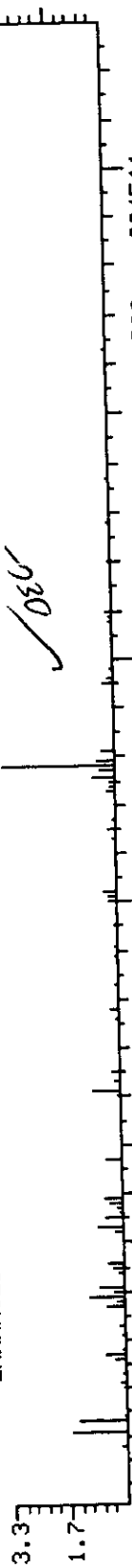
4200.



ENHANCED DATA: 30068101 #1091

BASE M/Z: 178 RIC: 15456.

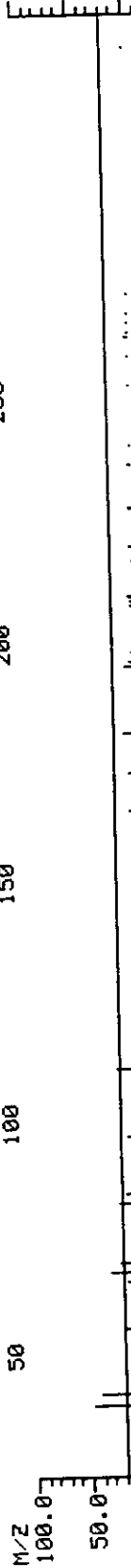
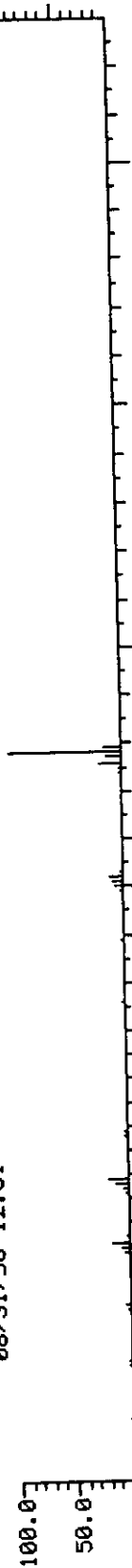
2448.



STANDARD FILE: ST16980831 #1091

BASE M/Z: 178 RIC: 204544.

73344.



OUT

Data Reduced by: *FW* Date: *9/16/98*
Data Reviewed by: *FW* Date: *9/18/98*

Data File: 30068101

QUANTERRA GC/MS TIC REPORT (Part 1)

Sample: S-MM5-2 1/3SA/100M
Analyst: DAT

INST. ID: F16
Date Analyzed: 08/31/98 19:00
Run Factor: 300.

# SCAN	Concentration in Sample (UG/A)	CAS #
X 444	7100.	74630-59-0
1-UNDECENE, 4-METHYL		

unkn. alkane 5746400

2 473	7900.	74962-98-4
2-TRIDECEN-1-OL, (E)		



X 556	6400.	1120-21-4
UNDECANE <i>or isomer</i>		

4 617	13000.	00-00-0
UNKNOWN		

5700

5 633	15000.	21078-85-9
4-DECANOL, 2-ETHYL		

576100

PAH

6 654	8900.	41446-67-7
3-TETRADECENE, (Z)		

or isomer

alkene 5720000
2/9/18/98

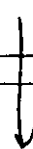
7 662	8600.	1120-21-4
UNDECANE <i>or isomer</i>		

X 678	6700.	00-00-0
UNKNOWN		

5700

9 700	8400.	00-00-0
UNKNOWN		

5700



X 714	6100.	00-00-0
UNKNOWN		

- 11 738 17000. 3891-98-3
DODECANE, 2,6,10-TRIMETHYL- or isomer
-
- 12 763 22000. 629-59-4
TETRADECANE or isomer
-
- 13 776 12000. 00-00-0 57620000
UNKNOWN
unkn. alkene
-
- 14 820 18000. 18344-37-1
HEPTADECANE, 2,6,10,14-TETRAMETHYL- or isomer
-
- 15 858 19000. 7045-71-8
UNDECANE, 2-METHYL- or isomer
-
- 16 904 13000. 52783-45-4
~~NONADECANOL~~ or isomer alkane 5746400
-
- 17 948 24000. 629-92-5
NONADECANE or isomer
-
- 18 987 15000. 54105-67-8
HEPTADECANE, 2,6-DIMETHYL- or isomer
-
- 19 1035 32000. 54105-67-8
HEPTADECANE, 2,6-DIMETHYL- or isomer
-
- 20 1114 10000. 54105-67-8
HEPTADECANE, 2,6-DIMETHYL- or isomer
-
- 21 1119 14000. 54105-67-8
HEPTADECANE, 2,6-DIMETHYL- or isomer
-
- 22 1191 7700. 629-92-5
NONADECANE or isomer
-

~~23~~ 1264 6400. 629-92-5
NONADECANE or *Isomer*

24 1335 9700. 18344-37-1
HEPTADECANE, 2,6,10,14-ETRAMETHYL-

25 1402 7500. 629-92-5
NONADECANE

DATA FILE: 30068101

QUANTERRA GC/MS TIC REPORT (Part 2)

CONCENTRATION = AREA(TIC)*CONC(IS)/AREA(IS)

#	FIT	PURITY	INT.		RRT	AREA	HEIGHT	AMOUNT		LIB	LIB #
			STD.	RT				AS ANALYZED	(UG/ML)		
1	947	652	1	4:56	0.838	178482.	55205.	23.760	NB	14793.	
2	862	667	1	5:15	0.892	198048.	49344.	26.364	NB	22482.	
3	964	839	1	6:11	1.049	160442.	59857.	21.358	NB	11607.	
4	883	598	1	6:51	1.164	331968.	36544.	44.192	UK	1.	
5	954	748	1	7:02	1.194	368896.	98560.	49.108	NB	19523.	
6	963	692	1	7:16	1.234	224000.	40832.	29.819	NB	21962.	
7	962	868	1	7:21	1.249	215296.	76160.	28.660	NB	11607.	
8	900	393	1	7:32	1.279	167552.	38144.	22.305	UK	1.	
9	887	401	2	7:47	0.844	197952.	27102.	27.953	UK	1.	
10	830	396	2	7:56	0.861	144707.	34489.	20.434	UK	1.	
11	981	834	2	8:12	0.890	411520.	117376.	58.111	NB	25991.	
12	959	739	2	8:29	0.920	522729.	149934.	73.815	NB	22530.	
13	946	550	2	8:37	0.936	294272.	40960.	41.554	UK	1.	
14	975	831	2	9:07	0.989	421760.	107008.	59.557	NB	42196.	
15	945	752	2	9:32	1.035	441984.	145920.	62.413	NB	15352.	
16	938	674	2	10:03	1.090	301824.	48512.	42.621	NB	40233.	
17	956	833	2	10:32	1.144	567424.	133632.	80.127	NB	37465.	
18	949	817	3	10:58	0.908	404224.	103040.	48.343	NB	37462.	
19	975	872	3	11:30	0.952	897408.	162304.	107.325	NB	37462.	
20	910	791	3	12:23	1.025	279936.	81024.	33.479	NB	37462.	
21	950	823	3	12:26	1.029	383872.	114688.	45.909	NB	37462.	
22	969	756	3	13:14	1.096	213440.	62304.	25.526	NB	37465.	
23	969	773	3	14:03	1.163	179442.	48179.	21.460	NB	37465.	
24	910	691	4	14:50	0.860	160448.	34827.	32.250	NB	42196.	
25	961	716	4	15:35	0.903	123776.	25984.	24.879	NB	37465.	

Library Search Data: 30068101 # 444 Base m/z: 43
 08/31/98 19:00:00 + 4:56 Cali: 30068101 # 3 RIC: 45952.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 740 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 14793 1-UNDECENE, 4-METHYL-
 2 19015 DECANE, 2,5,6-TRIMETHYL-
 3 11607 UNDECANE
 4 19523 1-DECANOL, 2-ETHYL-
 5 12074 1-HEPTANOL, 2-PROPYL-
 6 11602 OCTANE, 2,4,6-TRIMETHYL-
 7 8104 OCTANE, 3,5-DIMETHYL-
 8 15969 HYDROXYLAMINE, O-DECYL-
 9 8089 NONANE, 2-METHYL-

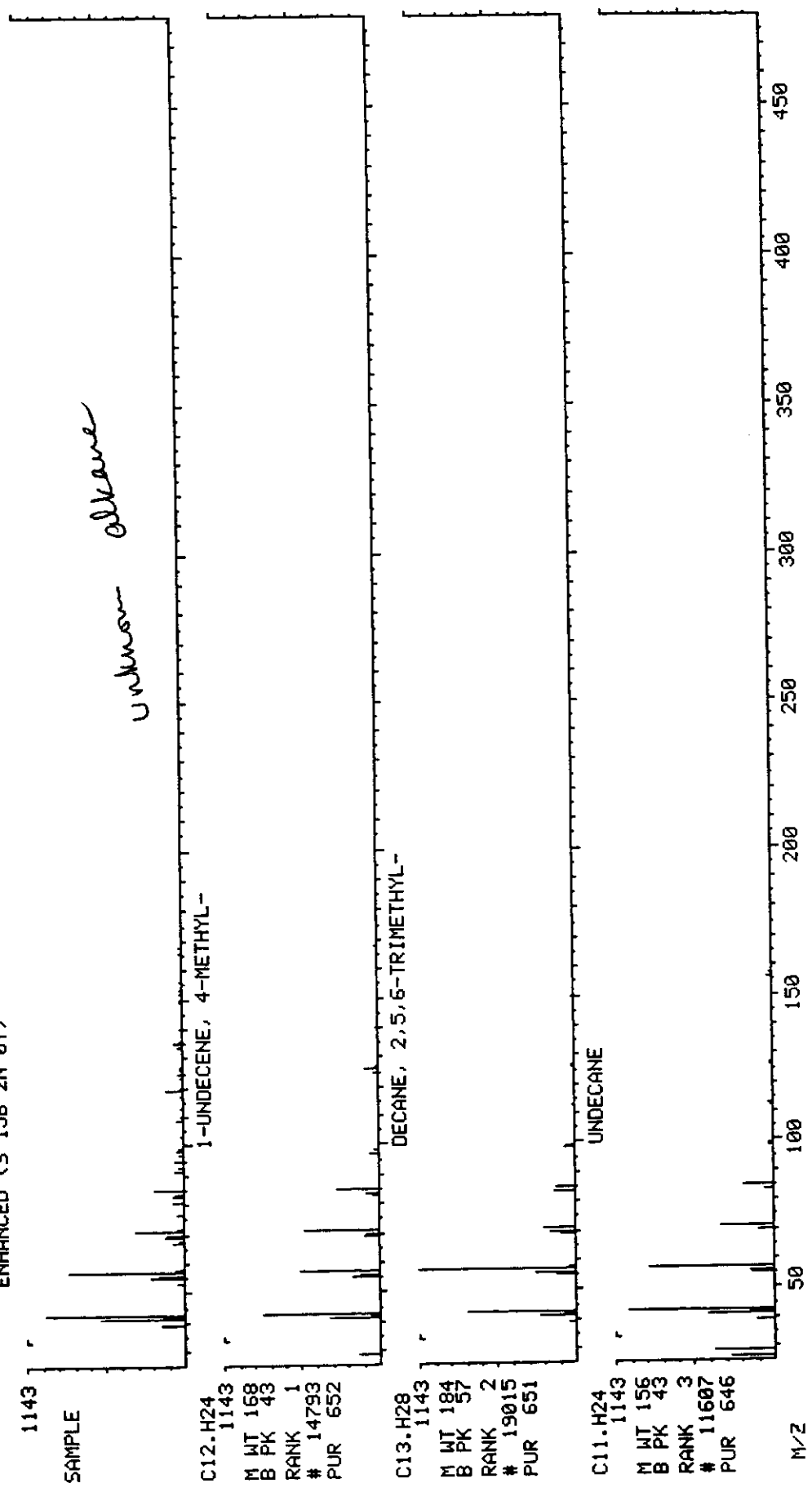
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H24	168	43	652	947	652
2	C13.H28	184	57	651	939	651
3	C11.H24	156	43	646	979	653
4	C12.H26.O	186	57	636	922	661
5	C10.H22.O	158	43	631	936	640
6	C11.H24	156	57	625	948	627
7	C10.H22	142	57	623	930	635
8	C10.H23.O.N	173	43	620	928	658
9	C10.H22	142	43	620	941	628

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	74630-39-0
2	—	—	—	—	62108-23-0
3	—	—	—	—	1120-21-4
4	—	—	—	—	21078-65-9
5	—	—	—	—	10042-59-8
6	—	—	—	—	62016-37-9
7	—	—	—	—	15869-93-9
8	—	—	—	—	29812-79-1
9	—	—	—	—	871-83-0

MID LIBRARY SEARCH (LIBRARY.NB)
 08/31/98 19:00:00 + 4:56
 SAMPLE: S-MMS-2 1/35A/100M
 COND.: UG/ML *100ML *100% *100% *(NA/NA) /1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068101 # 444
 CALI: 30068101 # 3

BASE M/Z: 43
 RIC: 45952.



Library Search Data: 30068101 # 473 Base m/z: 43
 08/31/98 19:00:00 + 5:16 Cali: 30068101 # 3 RIC: 40896.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 459 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 22482 2-TRIDECEN-1-OL, (E)-
 2 39230 9-EICOSYNE
 3 25971 OXIRANE, DODECYL-
 4 34800 (R)-(-)-(Z)-14-METHYL-8-HEXADECEN-1-OL
 5 37449 OCTADECANAL
 6 29222 (Z)6-PENTADECEN-1-OL
 7 39224 1-EICOSYNE
 8 37444 OXIRANE, HEXADECYL-
 9 28267 3-HEXADECYNE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H26.O	198	57	667	862	673
2	C20.H38	278	81	661	902	679
3	C14.H28.O	212	41	658	917	658
4	C17.H34.O	254	55	654	848	664
5	C18.H36.O	268	43	639	830	682
6	C15.H30.O	226	67	637	869	655
7	C20.H38	278	82	630	837	657
8	C18.H36.O	268	57	626	808	670
9	C16.H30	222	67	622	930	626

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	74962-98-4
2	—	—	—	—	71899-38-2
3	—	—	—	—	3234-28-4
4	—	—	—	—	30689-78-2
5	—	—	—	—	638-66-4
6	—	—	—	—	68797-95-5
7	—	—	—	—	765-27-5
8	—	—	—	—	7390-81-0
9	—	—	—	—	61886-62-2

DATA: 30068101 # 473
CALI: 30068101 # 3
BASE M/Z: 43
RIC: 40896.

MID LIBRARY SEARCH (LIBRARYNB)

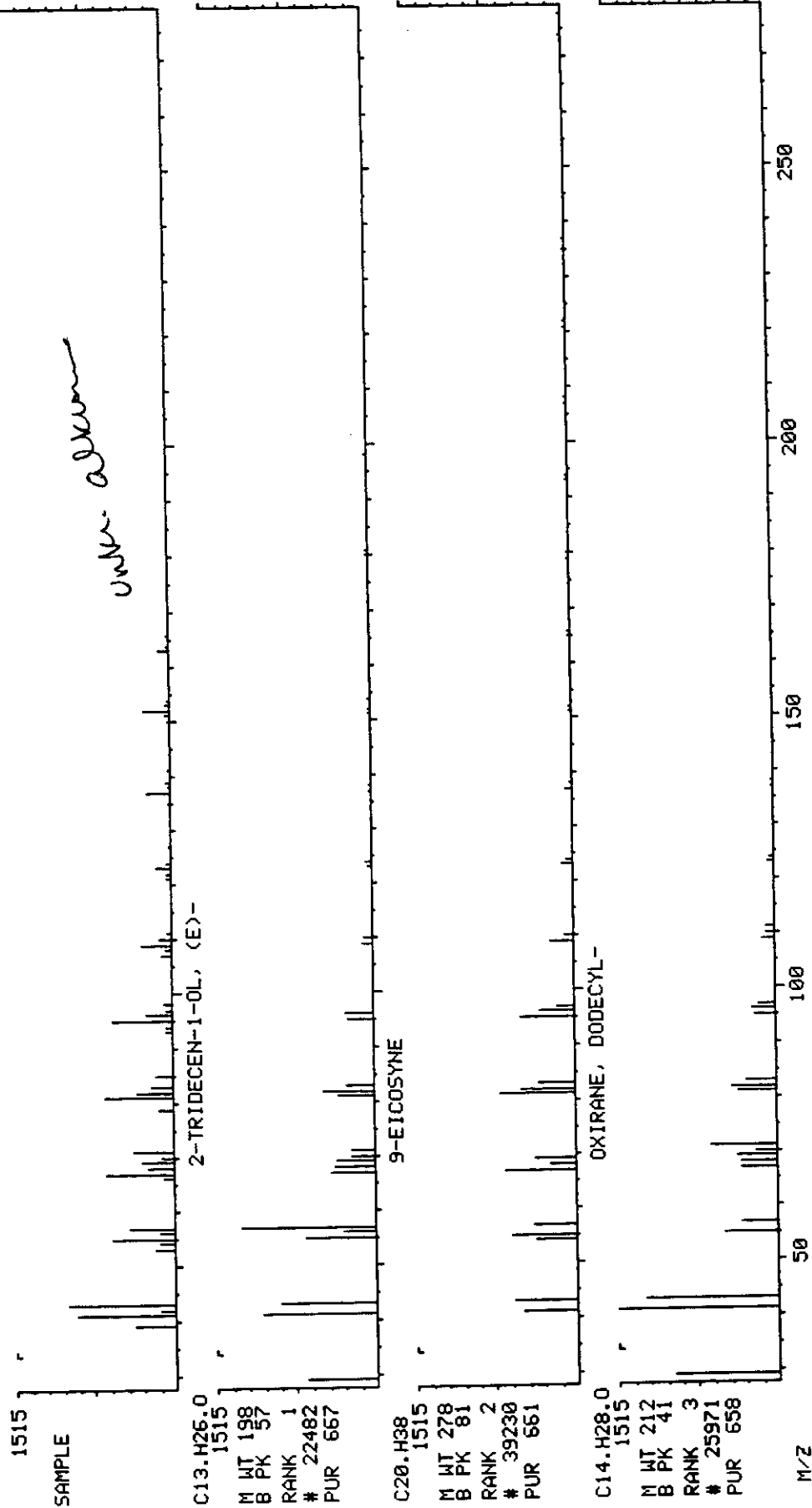
08/31/98 19:00:00 + 5:16

SAMPLE: S-MM5-2 1/35A/100M

INST. ID: F16

CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M

ENHANCED (S 158 2N 0T)



Library Search Data: 30068101 # 556 Base m/z: 43
 08/31/98 19:00:00 + 6:11 Cali: 30068101 # 3 RIC: 54144.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 890 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 11607 UNDECANE
 2 15353 2,6-DIMETHYLDECANE
 3 19016 UNDECANE, 4,7-DIMETHYL-
 4 15357 DECANE, 2,9-DIMETHYL-
 5 15352 UNDECANE, 2-METHYL-
 6 19015 DECANE, 2,5,6-TRIMETHYL-
 7 11602 OCTANE, 2,4,6-TRIMETHYL-
 8 19028 DECANE, 2,4,6-TRIMETHYL-
 9 22530 TETRADECANE

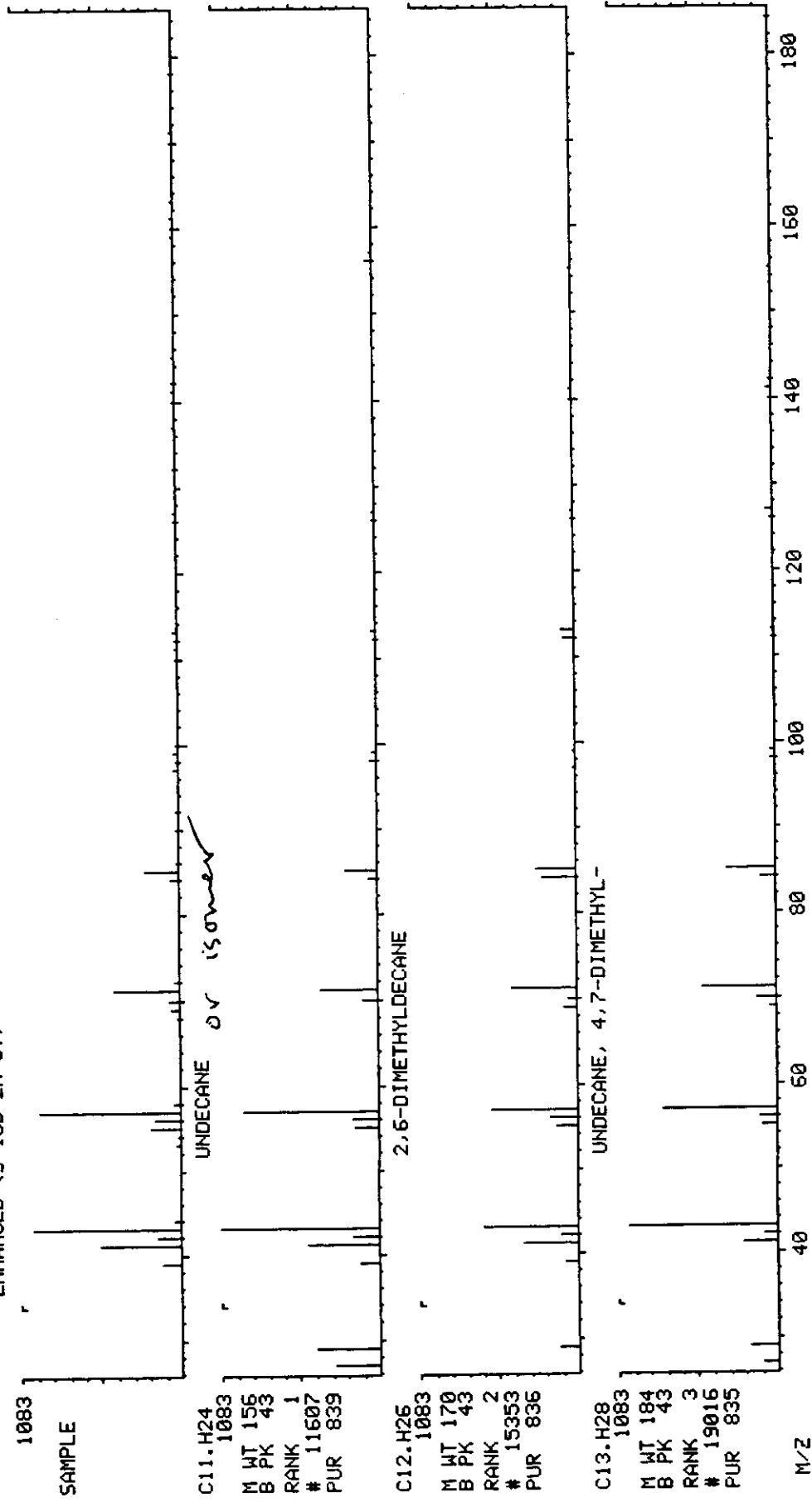
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C11.H24	156	43	839	964	864
2	C12.H26	170	43	836	924	842
3	C13.H28	184	43	835	959	836
4	C12.H26	170	43	832	967	842
5	C12.H26	170	43	829	949	845
6	C13.H28	184	57	825	929	826
7	C11.H24	156	57	822	967	822
8	C13.H28	184	43	822	941	828
9	C14.H30	198	43	821	923	876

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	1120-21-4
2	—	—	—	—	13150-81-7
3	—	—	—	—	17301-32-5
4	—	—	—	—	1002-17-1
5	—	—	—	—	7045-71-8
6	—	—	—	—	62108-23-0
7	—	—	—	—	62016-37-9
8	—	—	—	—	62108-27-4
9	—	—	—	—	629-59-4

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:00:00 + 6:11
 SAMPLE: S-MMS-2 1/35A/100M
 COND.: UG/ML *100ML *100Z/100Z *(NA/NA) /1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068101 # 556
 CALI: 30068101 # 3

BASE M/Z: 43
 RIC: 54144.



Library Search Data: 30068101 # 617 Base m/z: 41
 08/31/98 19:00:00 + 6:52 Cali: 30068101 # 3 RIC: 33216.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 207 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 24994 1,1'-BICYCLOHEXYL, 2-(1-METHYLETHYL)-, TRANS-
 2 33979 BORINIC ACID, DIETHYL-, 1-CYCLODODECEN-1-YL ESTER
 3 24993 1,1'-BICYCLOHEXYL, 2-(1-METHYLETHYL)-, CIS-
 4 39227 NAPHTHALENE, DECAHYDRO-2,6-DIMETHYL-3-OCTYL-
 5 39519 CYCLOHEXANE, 1-(1,5-DIMETHYLHEXYL)-4-(4-METHYLPENTYL)-
 6 24990 MUUROLANE-B
 7 31486 ACETAMIDE, N-METHYL-N-[4-[4-METHOXY-1-HEXAHYDROPIRIDYL]-2-BUTYNYL]-
 8 24987 AMORPHANE-A
 9 24991 AMORPHANE-B

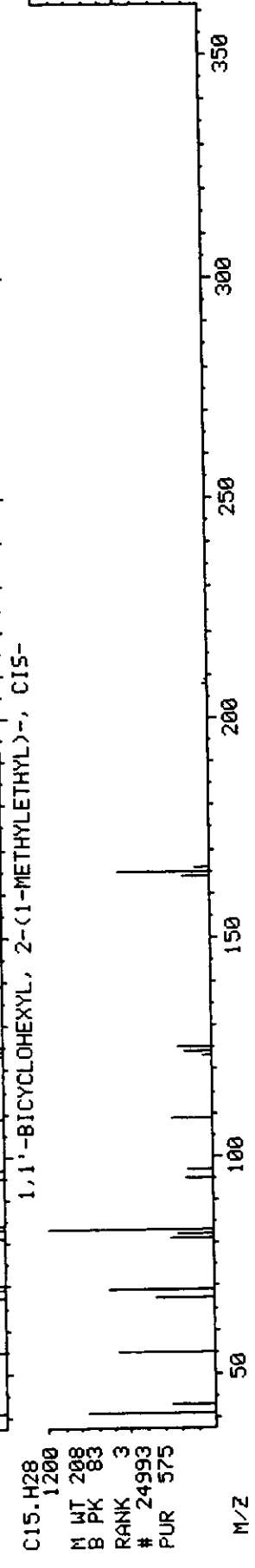
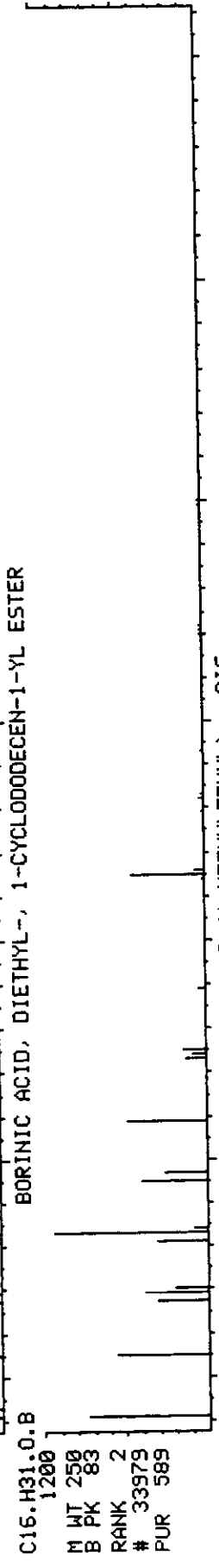
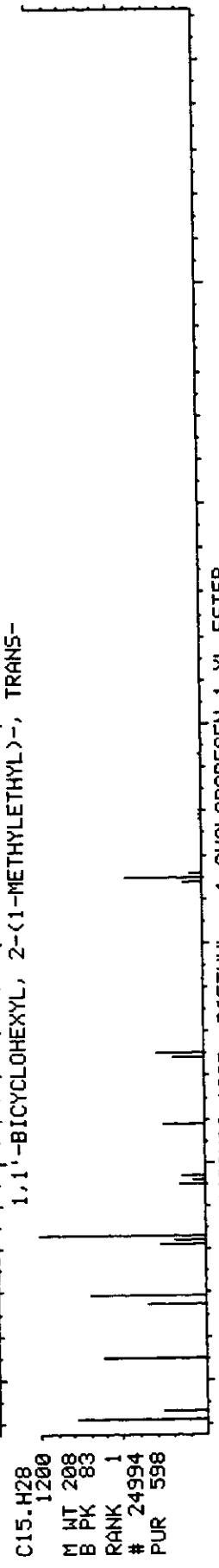
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C15.H28	208	83	598	883	622
2	C16.H31.O.B	250	83	589	911	617
3	C15.H28	208	83	575	868	615
4	C20.H38	278	109	539	820	606
5	C20.H40	280	43	537	741	636
6	C15.H28	208	109	534	830	586
7	C13.H22.O2.N2	238	43	527	732	606
8	C15.H28	208	109	519	817	552
9	C15.H28	208	95	518	804	581

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	50991-16-7
2	---	---	---	---	61142-73-2
3	---	---	---	---	50991-15-6
4	---	---	---	---	54964-85-1
5	---	---	---	---	56009-20-2
6	---	---	---	---	- -
7	---	---	---	---	- -
8	---	---	---	---	- -
9	---	---	---	---	- -

DATA: 30068101 # 617
 CALI: 30068101 # 3
 BASE M/Z: 41
 RIC: 33216.

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:00:00 + 6:52
 SAMPLE: 5-MM5-2 1/35A/100M
 COND.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (5 158 2N 0T)

INST. ID: F16



Library Search Data: 30068101 # 633 Base m/z: 57
 08/31/98 19:00:00 + 7:02 Cal: 30068101 # 3 RIC: 84096.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 919 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 19523 1-DECANOL, 2-ETHYL-
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 14799 1-DECENE, 3,4-DIMETHYL-
 4 26409 2-HEXYL-1-OCTANOL
 5 26408 2-ETHYL-1-DODECANOL
 6 32418 2-HEXYL-1-DECANOL
 7 22534 TRIDECANE, 7-METHYL-
 8 14751 2-UNDECENE, 5-METHYL-
 9 8539 1-PENTANOL, 4-METHYL-2-PROPYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H26.O	186	57	748	954	748
2	C19.H40	268	57	734	932	748
3	C12.H24	168	57	731	944	731
4	C14.H30.O	214	57	727	927	741
5	C14.H30.O	214	57	725	923	730
6	C16.H34.O	242	57	724	924	737
7	C14.H30	198	57	711	911	719
8	C12.H24	168	57	709	958	709
9	C9.H20.O	144	57	705	979	705

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	21078-65-9
2	—	—	—	—	54105-67-8
3	—	—	—	—	50871-03-9
4	—	—	—	—	- -
5	—	—	—	—	- -
6	—	—	—	—	- -
7	—	—	—	—	26730-14-3
8	—	—	—	—	56851-34-4
9	—	—	—	—	54004-41-0

DATA: 30068101 # 633
CALI: 30068101 # 3

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 19:00:00 + 7:02

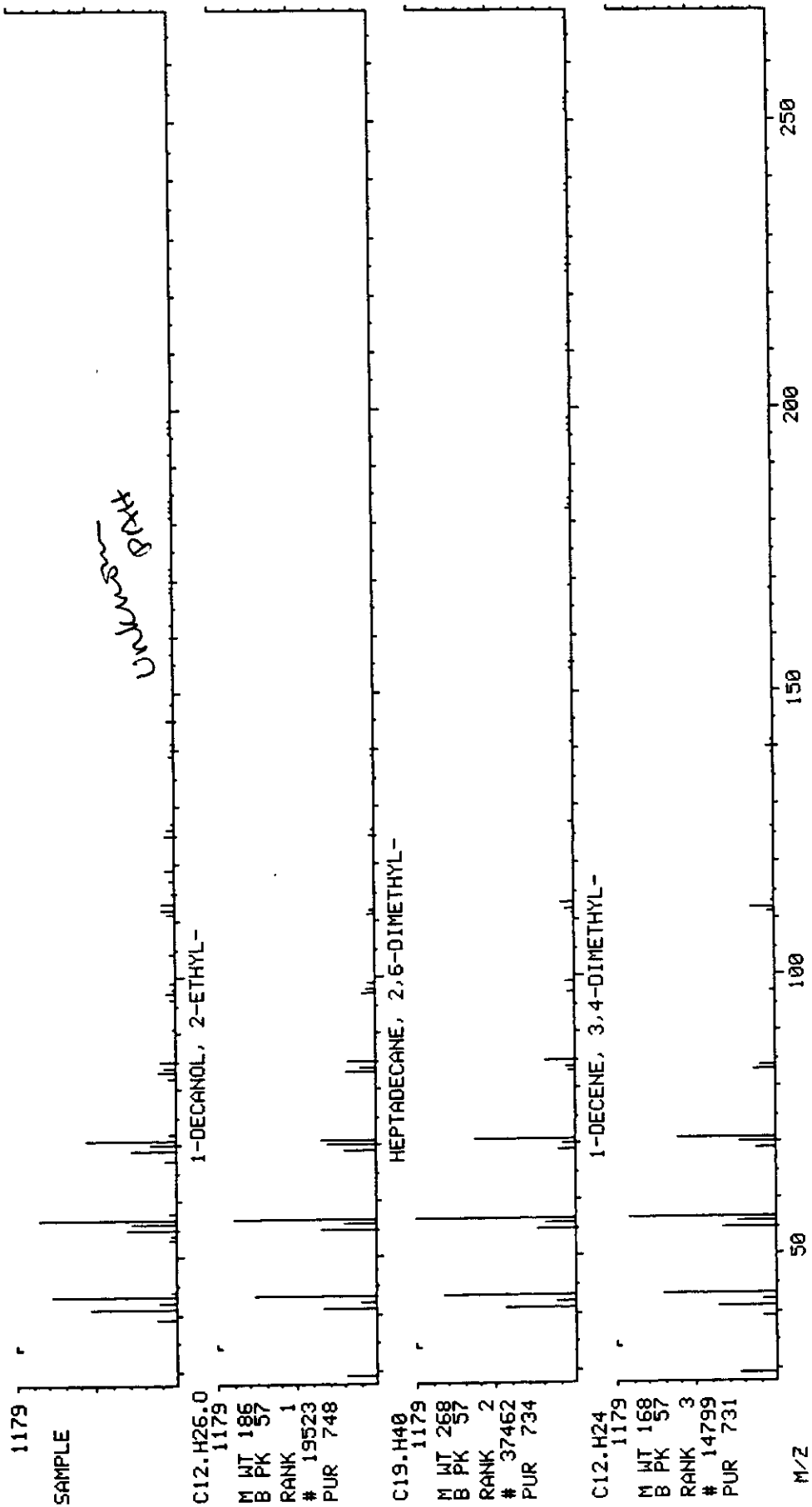
SAMPLE: S-MMS-2 1/35A/100M

COND.: UG/ML *100ML *100Z/100Z *(NA/NA) >1/35A NA M

ENHANCED (S 15B 2N 0T)

INST. ID: F16

BASE M/Z: 57
RIC: 84096.



632

Library Search Data: 30068101 # 654 Base m/z: 41
 08/31/98 19:00:00 + 7:16 Cali: 30068101 # 3 RIC: 39104.
 Sample: S-MMS-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

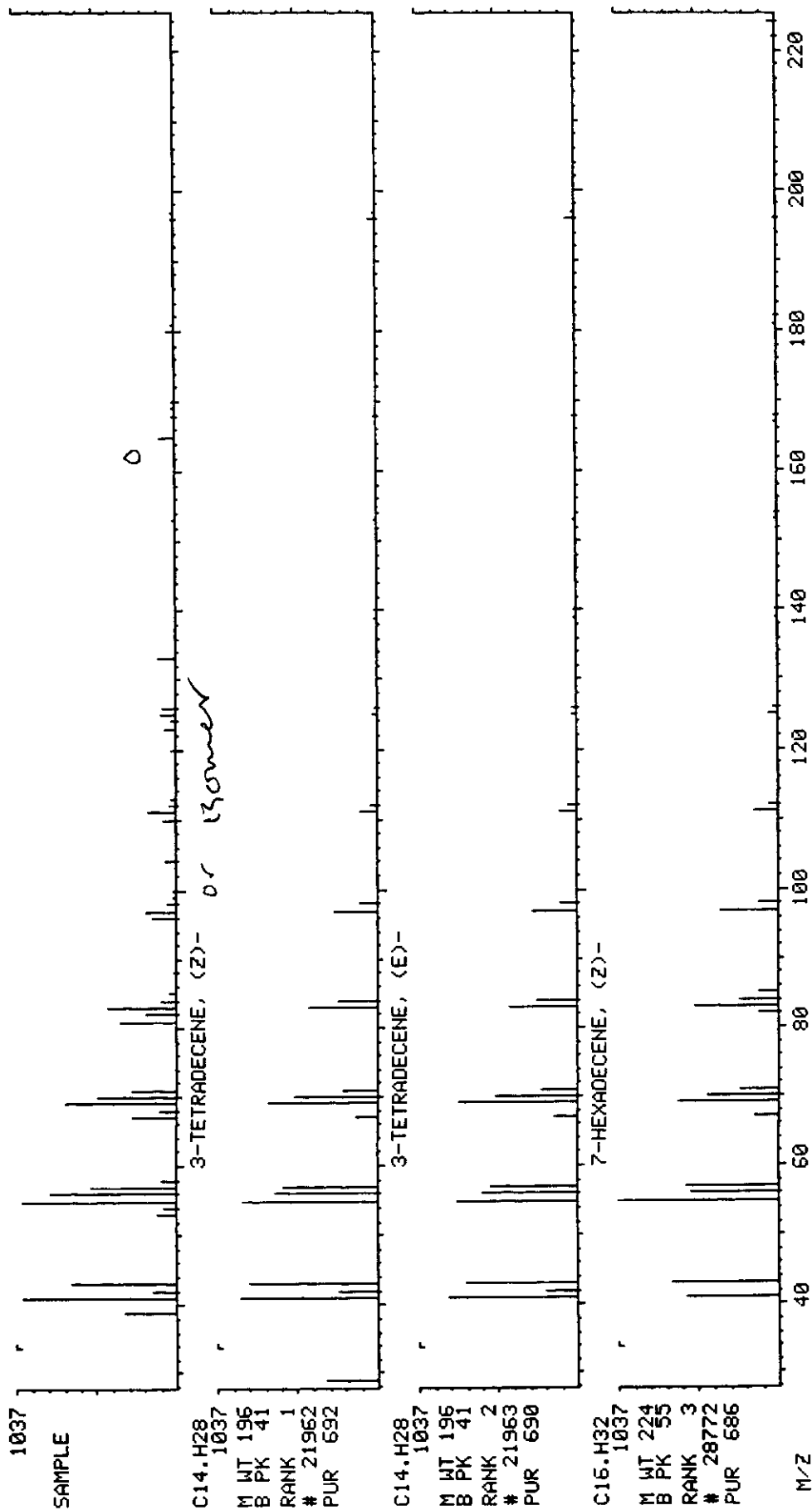
62231 spectra in LIBRARYNB searched for maximum PURITY
 687 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 21962 3-TETRADECENE, (Z)-
 2 21963 3-TETRADECENE, (E)-
 3 28772 7-HEXADECENE, (Z)-
 4 11084 CYCLOPROPANE, 1-HEPTYL-2-METHYL-
 5 28768 3-HEXADECENE, (Z)-
 6 21973 7-TETRADECENE, (Z)-
 7 34411 5-OCTADECENE, (E)-
 8 34410 3-OCTADECENE, (E)-
 9 21961 5-TETRADECENE, (E)-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C14.H28	196	41	692	963	695
2	C14.H28	196	41	690	959	693
3	C16.H32	224	55	686	923	710
4	C11.H22	154	56	685	943	689
5	C16.H32	224	55	685	925	711
6	C14.H28	196	55	683	952	687
7	C18.H36	252	55	682	917	708
8	C18.H36	252	69	681	904	708
9	C14.H28	196	55	680	949	683

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	41446-67-7
2	—	—	—	—	41446-68-8
3	—	—	—	—	35507-09-6
4	—	—	—	—	74663-91-5
5	—	—	—	—	34303-81-6
6	—	—	—	—	41446-60-0
7	—	—	—	—	7206-21-5
8	—	—	—	—	7206-19-1
9	—	—	—	—	41446-66-6

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:00:00 + 7:16
 SAMPLE: S-MM5-2 1/35A/100M
 COND.: UG/ML *100ML *100Z/100Z *KNA/NA >1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30068101 # 654
 CALI: 30068101 # 3
 BASE M/Z: 41
 RIC: 39104.



Library Search Data: 30068101 # 662 Base m/z: 57
 08/31/98 19:00:00 + 7:22 Cali: 30068101 # 3 RIC: 75264.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

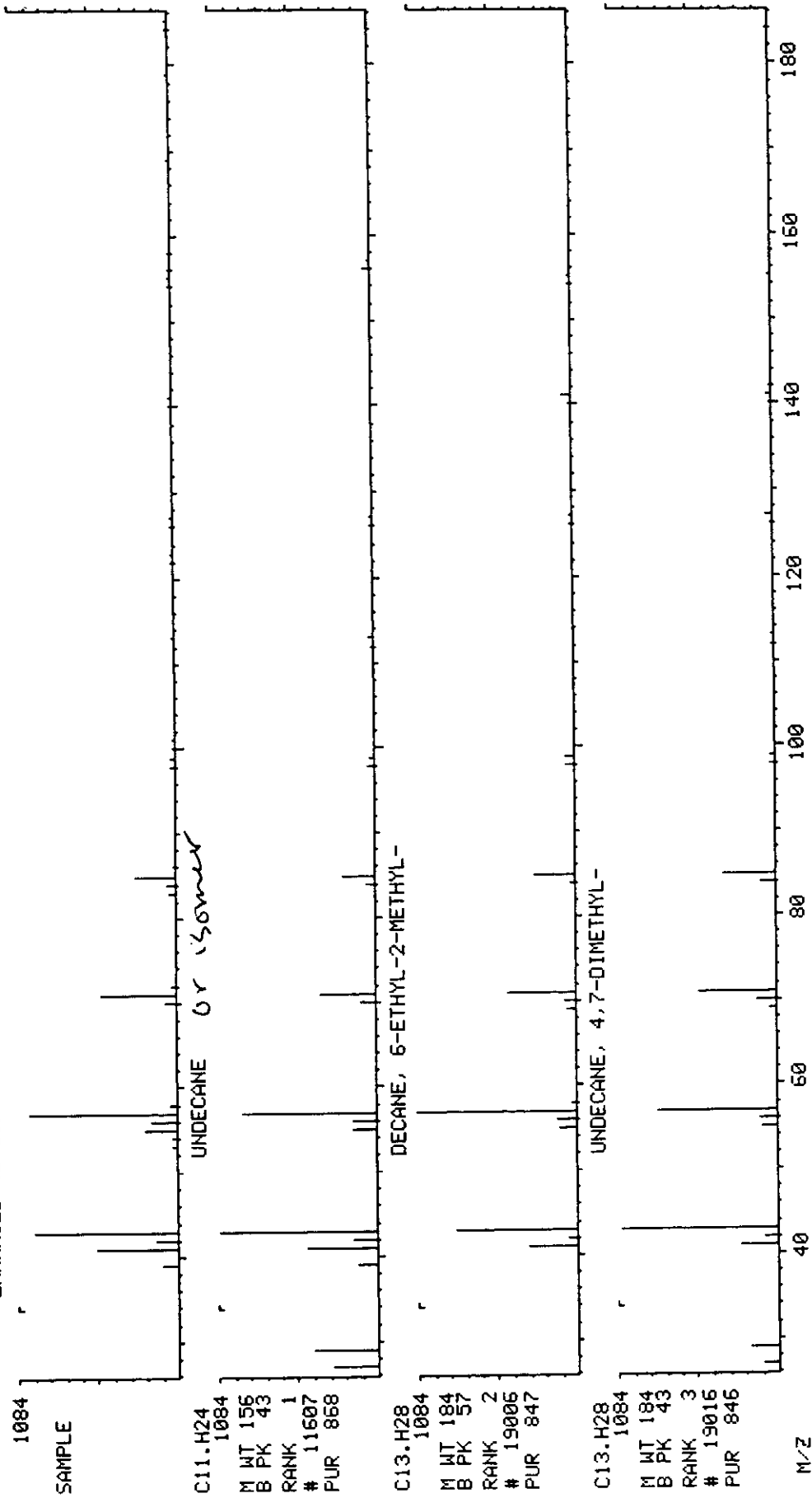
62231 spectra in LIBRARYNB searched for maximum PURITY
 632 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 11607 UNDECANE
 2 19006 DECANE, 6-ETHYL-2-METHYL-
 3 19016 UNDECANE, 4,7-DIMETHYL-
 4 22530 TETRADECANE
 5 11602 OCTANE, 2,4,6-TRIMETHYL-
 6 15357 DECANE, 2,9-DIMETHYL-
 7 18998 UNDECANE, 3,7-DIMETHYL-
 8 19028 DECANE, 2,4,6-TRIMETHYL-
 9 18985 TRIDECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C11.H24	156	43	868	962	878
2	C13.H28	184	57	847	943	862
3	C13.H28	184	43	846	934	863
4	C14.H30	198	43	831	908	897
5	C11.H24	156	57	830	945	850
6	C12.H26	170	43	826	943	848
7	C13.H28	184	43	826	925	850
8	C13.H28	184	43	825	915	858
9	C13.H28	184	57	819	923	851

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	1120-21-4
2	—	—	—	—	62108-21-8
3	—	—	—	—	17301-32-5
4	—	—	—	—	629-59-4
5	—	—	—	—	62016-37-9
6	—	—	—	—	1002-17-1
7	—	—	—	—	17301-29-0
8	—	—	—	—	62108-27-4
9	—	—	—	—	629-50-5

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:00:00 + 7:22
 SAMPLE: 5-MMS-2 1/35A/100M
 CONDS.: UC/ML #100ML #100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30068101 # 662
 CALI: 30068101 # 3
 BASE M/Z: 57
 RIC: 75264.



Library Search Data: 30068101 # 678 Base m/z: 57
 08/31/98 19:00:00 + 7:32 Cali: 30068101 # 3 RIC: 36736.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 413 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 11597 NONANE, 3,7-DIMETHYL-
 2 5159 NONANE
 3 37251 1-iodo-2-methylnonane
 4 5041 3-HEXANONE, 2,4-DIMETHYL-
 5 12534 NAPHTHALENE, 1,2,3,4-TETRAHYDRO-2,6-DIMETHYL-
 6 15343 DODECANE
 7 11607 UNDECANE
 8 19006 DECANE, 6-ETHYL-2-METHYL-
 9 8093 OCTANE, 2,2-DIMETHYL-

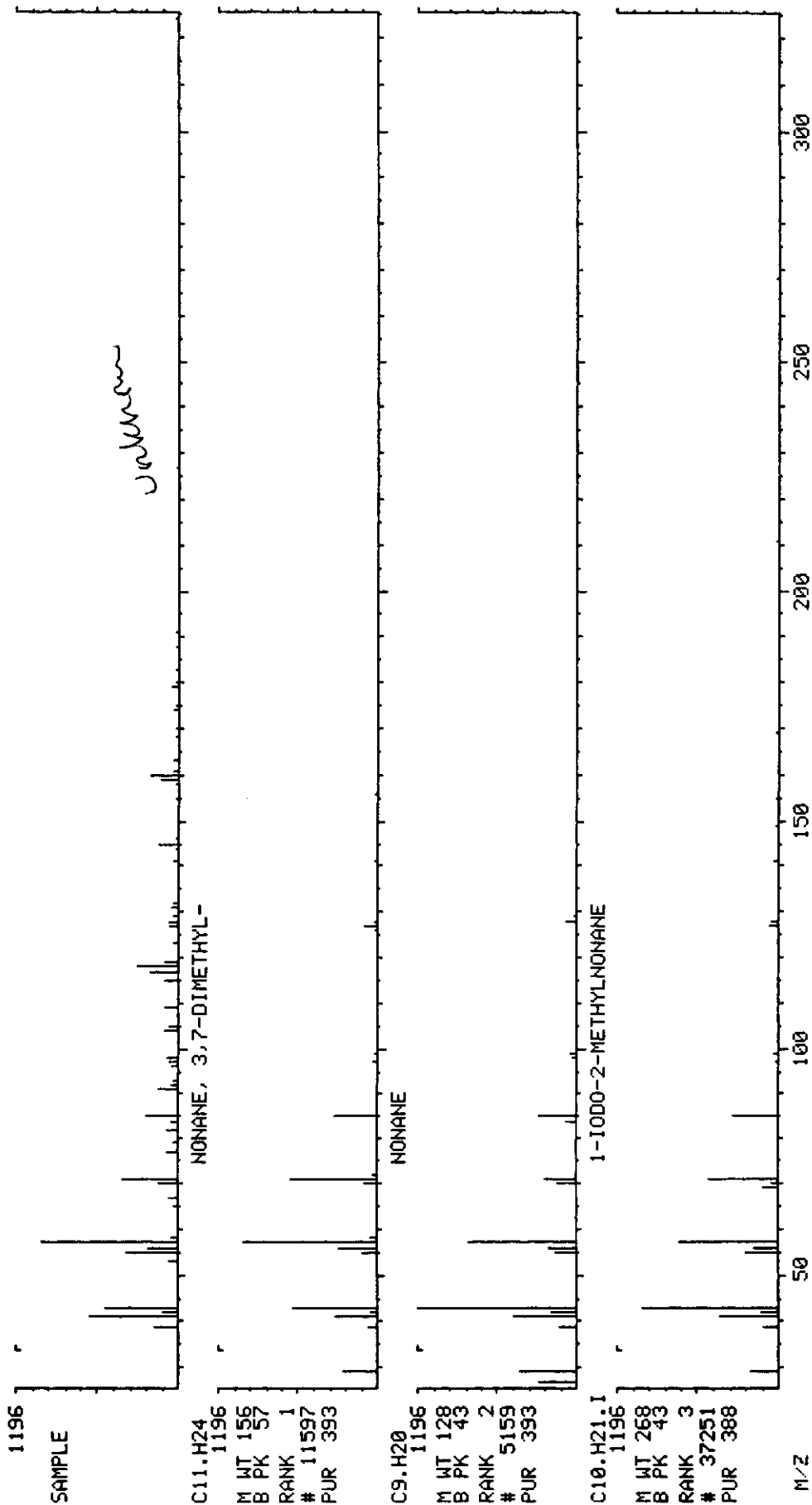
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C11.H24	156	57	393	900	404
2	C9.H20	128	43	393	912	402
3	C10.H21.I	268	43	388	897	412
4	C8.H16.O	128	57	385	918	389
5	C12.H16	160	118	375	821	387
6	C12.H26	170	57	373	870	399
7	C11.H24	156	43	373	885	401
8	C13.H28	184	57	372	881	395
9	C10.H22	142	57	371	894	376

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	17302-32-8
2	—	—	—	—	111-84-2
3	—	—	—	—	- -
4	—	—	—	—	18641-70-8
5	—	—	—	—	7524-63-2
6	—	—	—	—	112-40-3
7	—	—	—	—	1120-21-4
8	—	—	—	—	62108-21-8
9	—	—	—	—	15869-87-1

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:00:00 + 7:32
 SAMPLE: S-MMS-2 1/35A/100M
 COND.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068101 # 678
 CALI: 30068101 # 3

BASE M/Z: 57
 RIC: 36735.



Library Search Data: 30068101 # 700 Base m/z: 81
 08/31/98 19:00:00 + 7:47 Cali: 30068101 # 3 RIC: 25344.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 294 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 6996 1-ETHYNYL-1-CYCLOHEPTANOL
 2 28273 8-HEXADECYNE
 3 36677 7-OCTADECYNE, 2-METHYL-
 4 18371 5,7-DIMETHYLOCTAHYDROCOUMARIN
 5 10423 3-UNDECYNE
 6 28267 3-HEXADECYNE
 7 14158 3-DODECYNE
 8 25001 CYCLOHEXANE, 1-(CYCLOHEXYLMETHYL)-2-ETHYL-, TRANS-
 9 34010 3-OCTADECYNE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C9.H14.O	138	81	401	887	401
2	C16.H30	222	81	395	777	436
3	C19.H36	264	81	394	749	439
4	C11.H18.O2	182	95	392	863	412
5	C11.H20	152	67	389	893	389
6	C16.H30	222	67	388	732	405
7	C12.H22	166	67	378	832	380
8	C15.H28	208	55	375	707	426
9	C18.H34	250	67	374	810	395

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	2809-78-1
2	—	—	—	—	19781-86-3
3	—	—	—	—	35354-38-2
4	—	—	—	—	—
5	—	—	—	—	60212-30-8
6	—	—	—	—	61886-62-2
7	—	—	—	—	6790-27-8
8	—	—	—	—	54934-92-8
9	—	—	—	—	61886-64-4

BASE M/Z: 81
RIC: 25344.

DATA: 30068101 # 700
CALI: 30068101 # 3

MID LIBRARY SEARCH (LIBRARYNB)

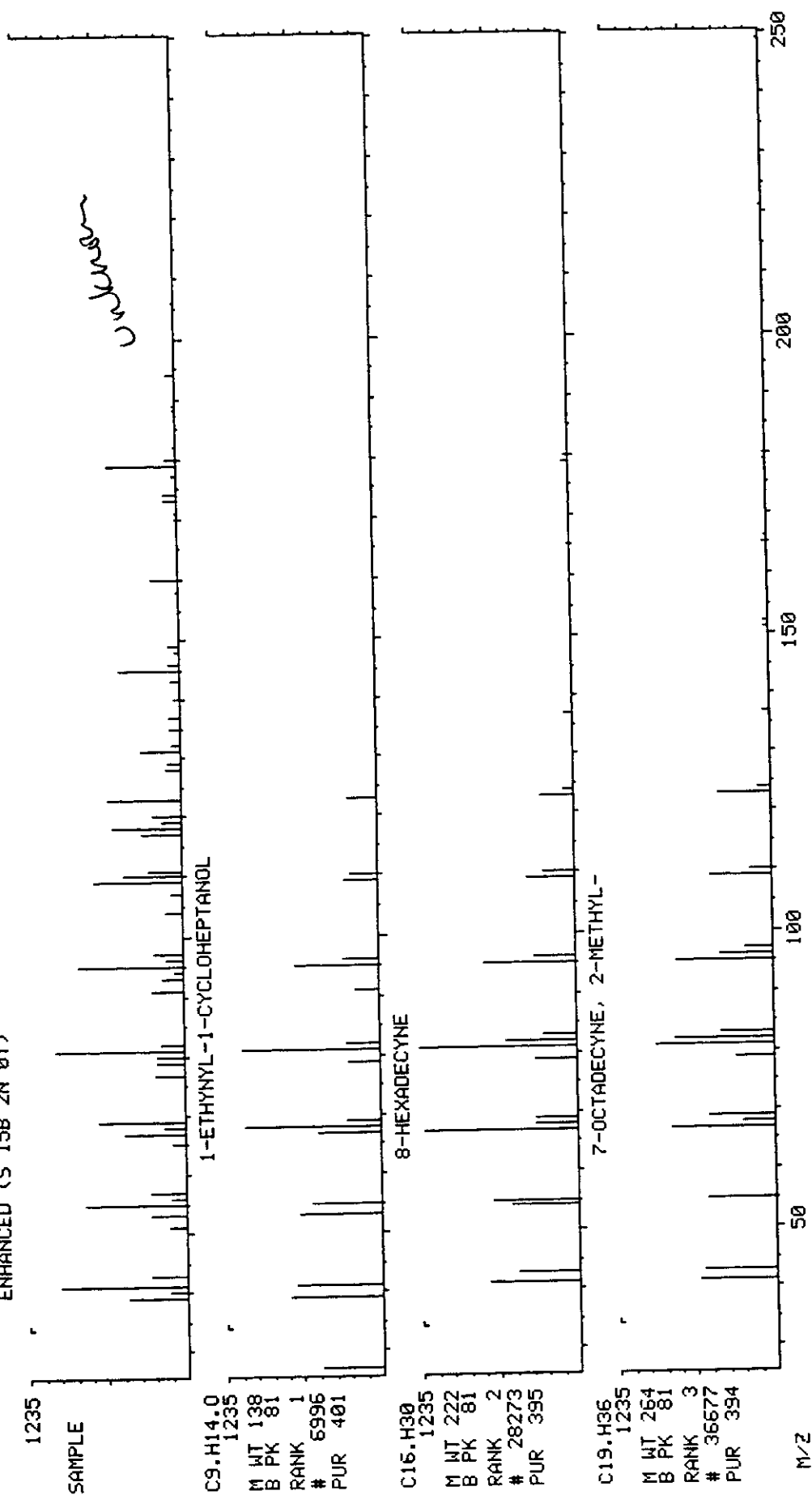
08/31/98 19:00:00 + 7:47

SAMPLE: S-MM5-2 1/35A/100M

COND.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M

ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068101 # 714 Base m/z: 159
 08/31/98 19:00:00 + 7:57 Cali: 30068101 # 3 RIC: 35264.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

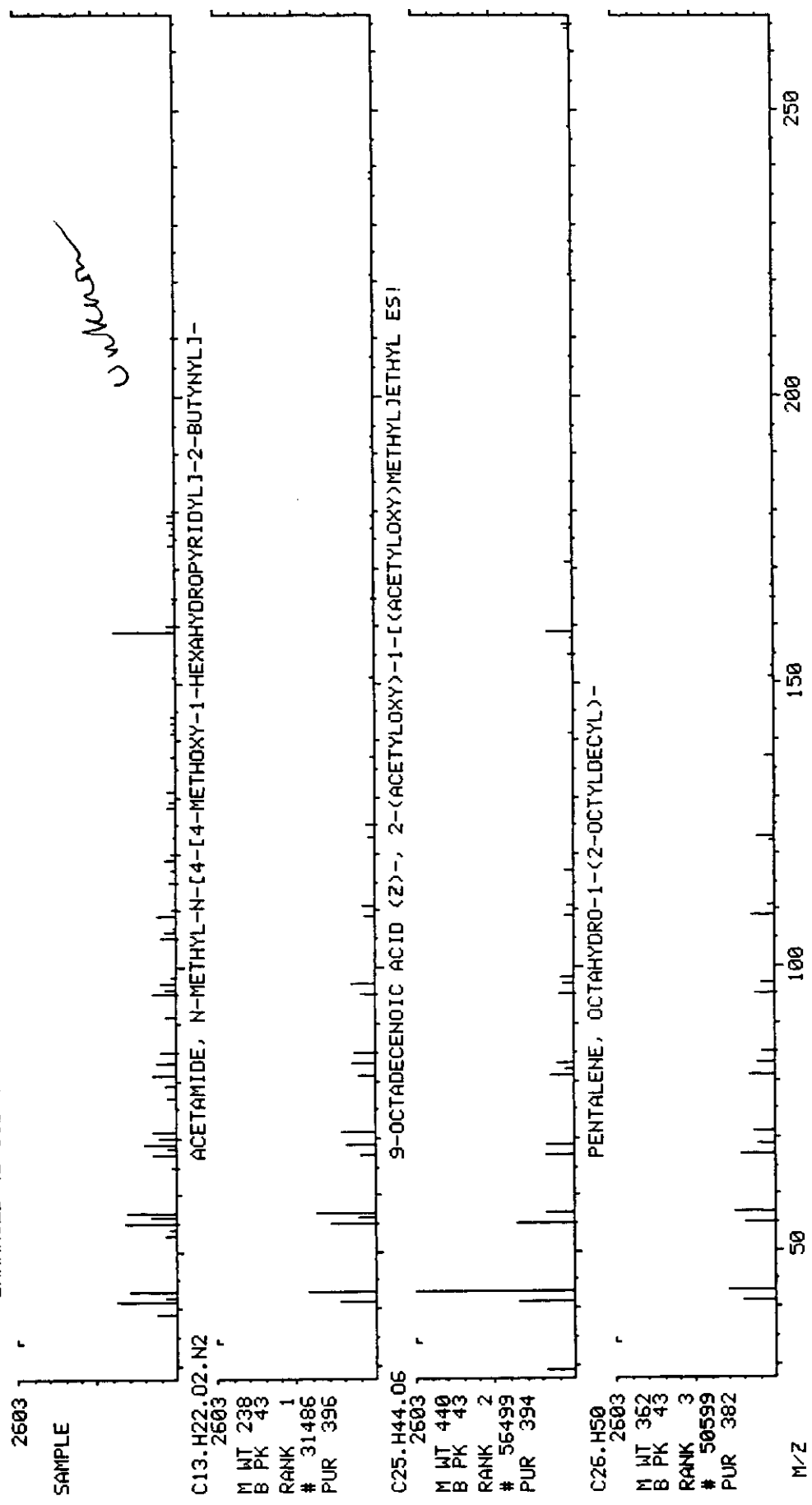
62231 spectra in LIBRARYNB searched for maximum PURITY
 310 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 31486 ACETAMIDE, N-METHYL-N-[4-[4-METHOXY-1-HEXAHYDROPYRIDYL]-2-BUTYNYL]-
 2 56499 9-OCTADECENOIC ACID (Z)-, 2-(ACETYLOXY)-1-[(ACETYLOXY)METHYL]ETHYL *
 3 50599 PENTALENE, OCTAHYDRO-1-(2-OCTYLDECYL)-
 4 56413 9,12-OCTADECADIENOIC ACID (Z,Z)-, 2-(ACETYLOXY)-1-[(ACETYLOXY)METHY*
 5 56303 9,12,15-OCTADECATRIENOIC ACID, 2-(ACETYLOXY)-1-[(ACETYLOXY)METHYL]E*
 6 40193 CIS-9,10-EPOXYOCTADECAN-1-OL
 7 32052 OXIRANE, TETRADECYL-
 8 46251 DODECANE, 1,2-DIBROMO-
 9 5032 3,4-DIMETHYLCYCLOHEXANOL

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H22.O2.N2	238	43	396	830	457
2	C25.H44.O6	440	43	394	705	496
3	C26.H50	362	43	382	847	415
4	C25.H42.O6	438	43	380	708	454
5	C25.H40.O6	436	43	369	718	449
6	C18.H36.O2	284	55	367	848	415
7	C16.H32.O	240	41	366	847	411
8	C12.H24.BR2	326	41	351	778	432
9	C8.H16.O	128	71	339	834	357

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	—
2	—	—	—	—	55401-63-3
3	—	—	—	—	55401-65-5
4	—	—	—	—	55320-03-1
5	—	—	—	—	55320-01-9
6	—	—	—	—	13980-12-6
7	—	—	—	—	7320-37-8
8	—	—	—	—	55334-42-4
9	—	—	—	—	5715-23-1

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:00:00 + 7:57
 SAMPLE: 5-MM5-2 1/35A/100M
 COND5.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068101 # 714
 CALI: 30068101 # 3
 BASE M/Z: 159
 RIC: 35264.



Library Search Data: 30068101 # 738 Base m/z: 57
 08/31/98 19:00:00 + 8:13 Cal: 30068101 # 3 RIC: 103296.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

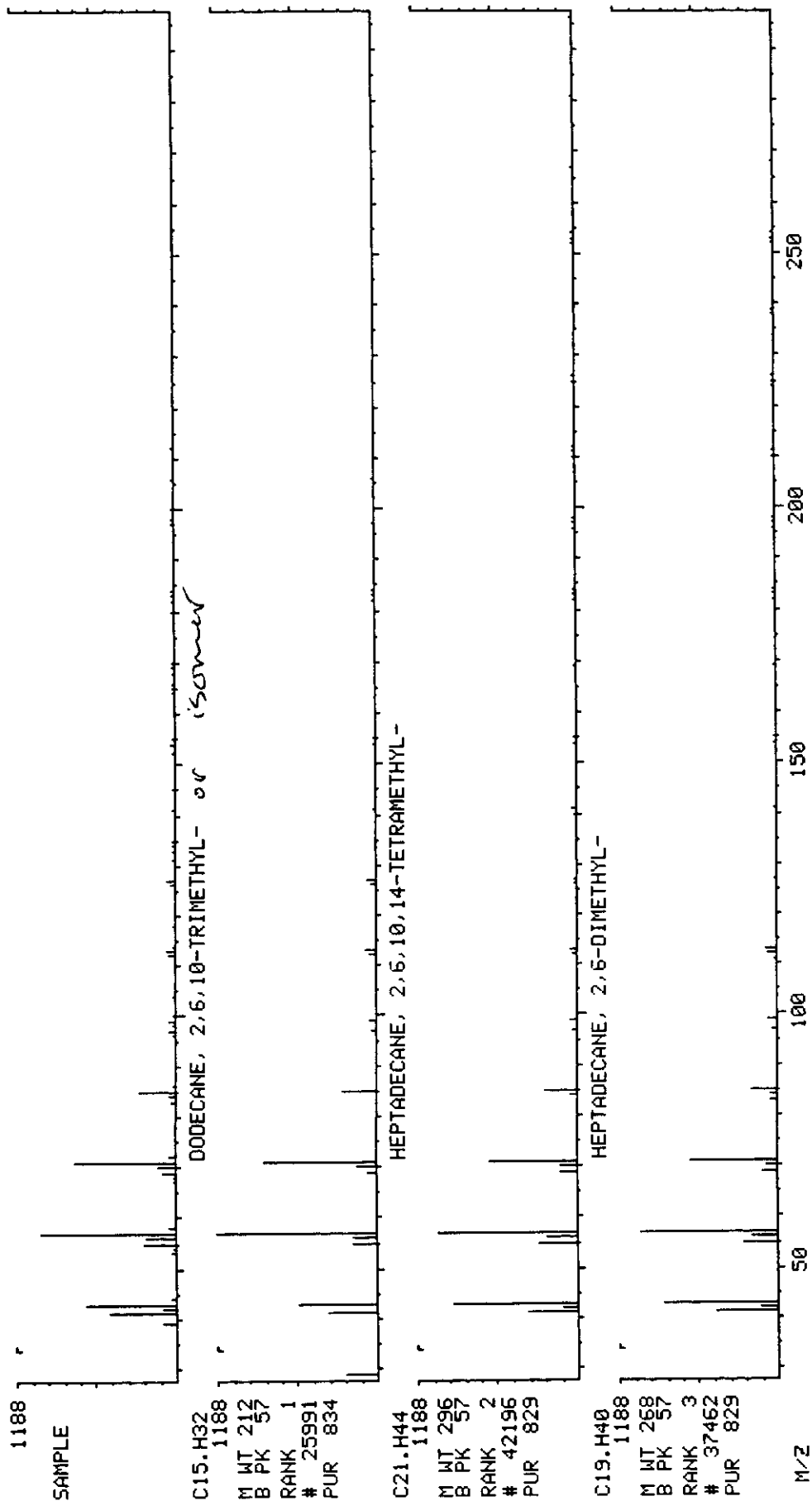
62231 spectra in LIBRARYNB searched for maximum PURITY
 712 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 25991 DODECANE, 2,6,10-TRIMETHYL-
 2 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 3 37462 HEPTADECANE, 2,6-DIMETHYL-
 4 25997 PENTADECANE
 5 26001 DODECANE, 2,7,10-TRIMETHYL-
 6 22535 DODECANE, 4,6-DIMETHYL-
 7 37465 NONADECANE
 8 25994 DODECANE, 2,6,11-TRIMETHYL-
 9 22530 TETRADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C15.H32	212	57	834	981	834
2	C21.H44	296	57	829	944	850
3	C19.H40	268	57	829	950	846
4	C15.H32	212	57	817	948	829
5	C15.H32	212	57	815	982	821
6	C14.H30	198	57	811	953	817
7	C19.H40	268	57	811	929	841
8	C15.H32	212	57	805	968	814
9	C14.H30	198	43	802	928	813

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	3891-98-3
2	—	—	—	—	18344-37-1
3	—	—	—	—	54105-67-8
4	—	—	—	—	629-62-9
5	—	—	—	—	74645-98-0
6	—	—	—	—	61141-72-8
7	—	—	—	—	629-92-5
8	—	—	—	—	31295-56-4
9	—	—	—	—	629-59-4

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:00:00 + 8:13
 SAMPLE: S-MM5-2 1/35A/100M
 COND.: UG/ML *100ML *100% /100% * (NA/NA) /1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30058101 # 738
 CALI: 30058101 # 3
 BASE M/Z: 57
 RIC: 103296.



Library Search Data: 30068101 # 763 Base m/z: 43
 08/31/98 19:00:00 + 8:29 Cali: 30068101 # 3 RIC: 136960.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

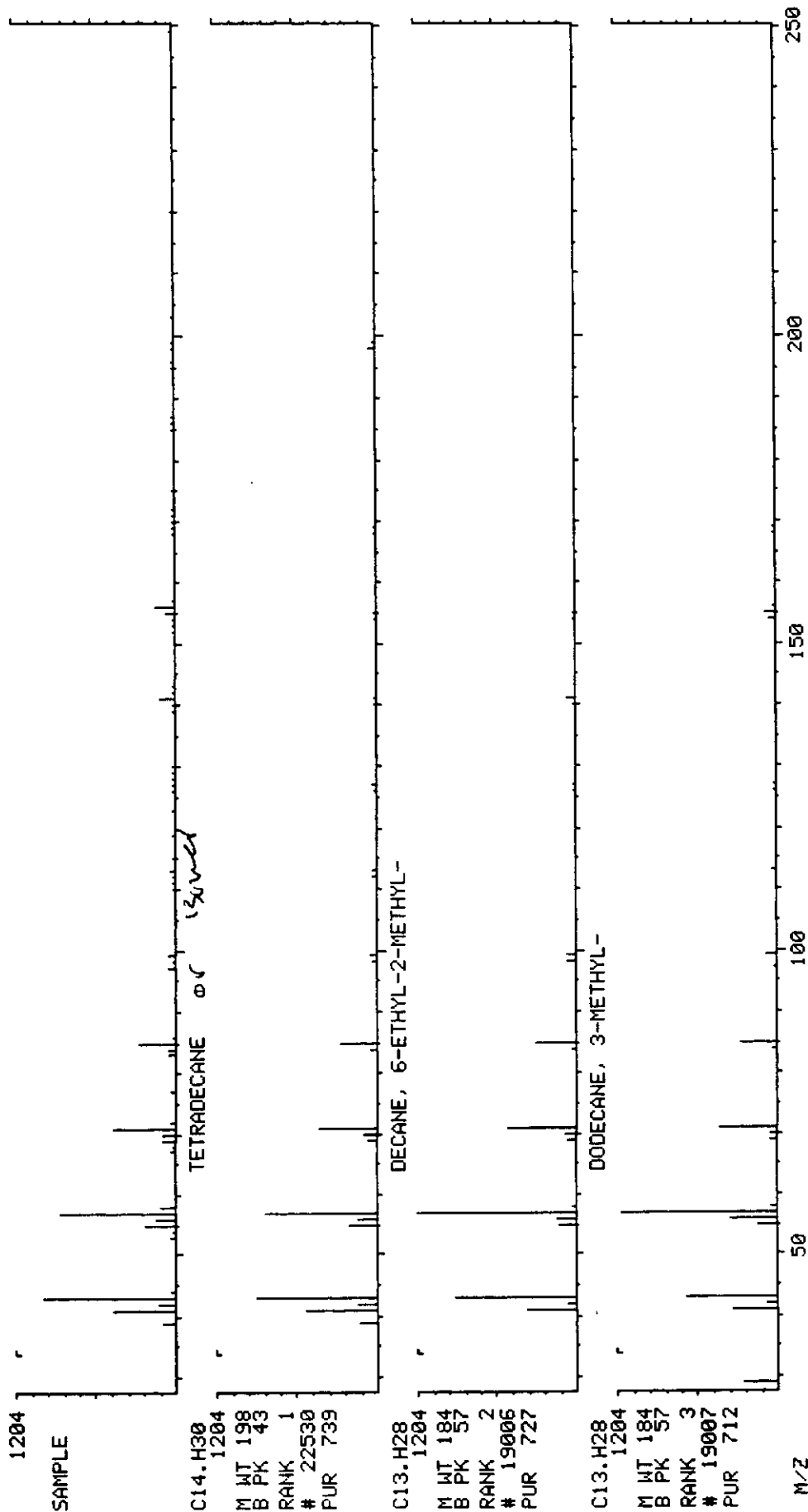
62231 spectra in LIBRARYNB searched for maximum PURITY
 611 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 22530 TETRADECANE
 2 19006 DECANE, 6-ETHYL-2-METHYL-
 3 19007 DODECANE, 3-METHYL-
 4 11607 UNDECANE
 5 11612 NONANE, 2,5-DIMETHYL-
 6 15358 UNDECANE, 3-METHYL-
 7 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 8 18990 UNDECANE, 2,9-DIMETHYL-
 9 39681 1-iodoundecane

Rank	Formula	M.Wt	B.Pk	Purity	Fit	Rfit
1	C14.H30	198	43	739	959	739
2	C13.H28	184	57	727	964	727
3	C13.H28	184	57	712	904	712
4	C11.H24	156	43	711	968	711
5	C11.H24	156	57	708	871	708
6	C12.H26	170	57	703	924	703
7	C21.H44	296	57	700	918	729
8	C13.H28	184	57	699	876	699
9	C11.H23.I	282	57	697	875	704

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-59-4
2	—	—	—	—	62108-21-8
3	—	—	—	—	17312-57-1
4	—	—	—	—	1120-21-4
5	—	—	—	—	17302-27-1
6	—	—	—	—	1002-43-3
7	—	—	—	—	18344-37-1
8	—	—	—	—	17301-26-7
9	—	—	—	—	4282-44-4

MID LIBRARY SEARCH (LIBRARYB)
 08/31/98 19:00:00 + 8:29
 SAMPLE: S-MM5-2 1/35A/100M
 COND5.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30068101 # 763
 CALI: 30068101 # 3
 BASE M/Z: 43
 RIC: 136960.



Library Search Data: 30068101 # 776 Base m/z: 156
 08/31/98 19:00:00 + 8:38 Cali: 30068101 # 3 RIC: 36864.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

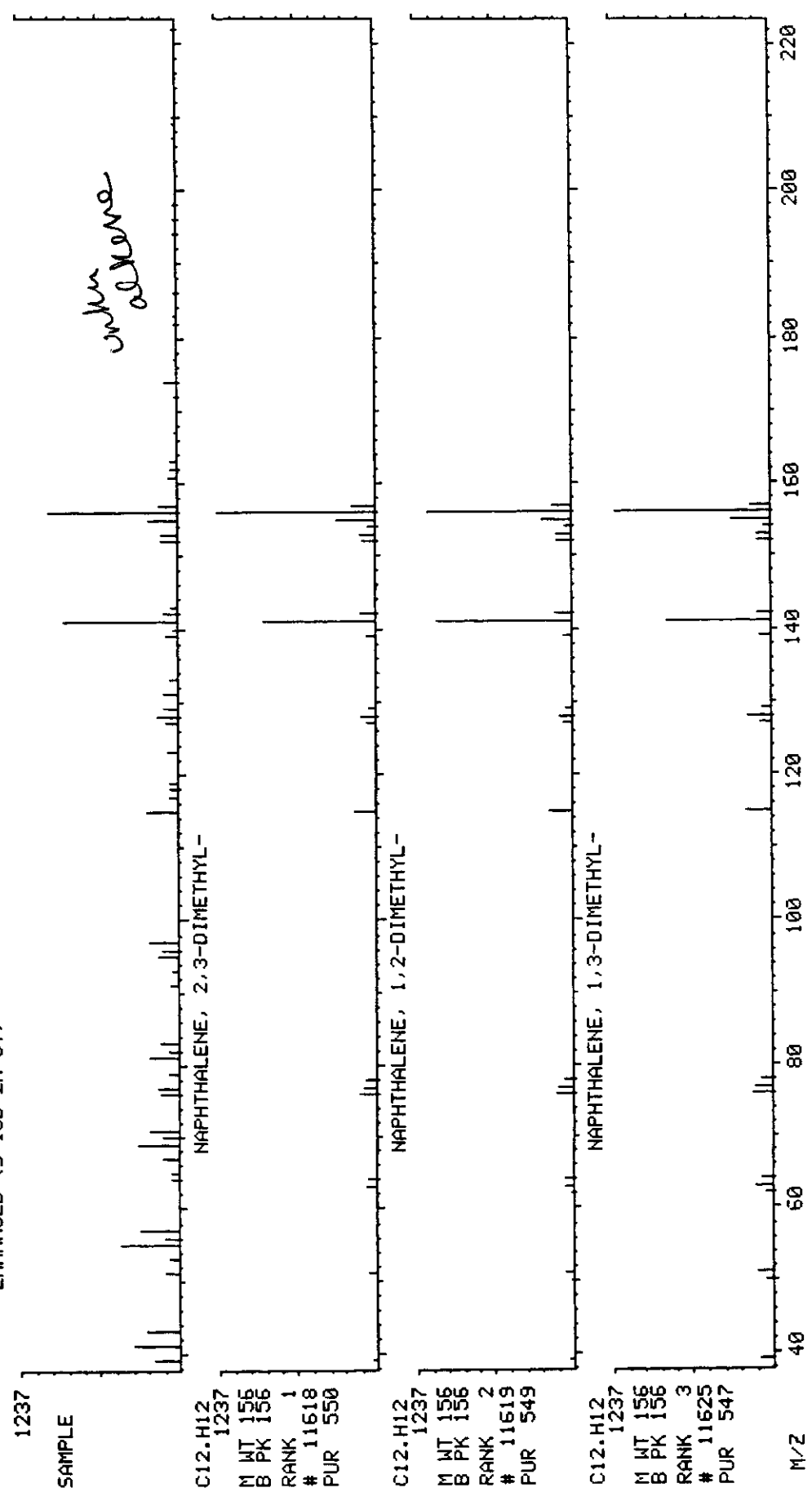
62231 spectra in LIBRARYNB searched for maximum PURITY
 205 matched at least 5 of the 16 largest peaks in the unknown

Rank In.	Name
1	11618 NAPHTHALENE, 2,3-DIMETHYL-
2	11619 NAPHTHALENE, 1,2-DIMETHYL-
3	11625 NAPHTHALENE, 1,3-DIMETHYL-
4	11629 NAPHTHALENE, 1,5-DIMETHYL-
5	11624 NAPHTHALENE, 1,7-DIMETHYL-
6	11620 NAPHTHALENE, 2,6-DIMETHYL-
7	11628 NAPHTHALENE, 1,6-DIMETHYL-
8	11615 NAPHTHALENE, 2,7-DIMETHYL-
9	11626 NAPHTHALENE, 1,4-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H12	156	156	550	946	569
2	C12.H12	156	156	549	944	570
3	C12.H12	156	156	547	941	576
4	C12.H12	156	156	540	928	571
5	C12.H12	156	156	539	927	574
6	C12.H12	156	156	537	938	568
7	C12.H12	156	156	537	924	564
8	C12.H12	156	156	537	923	564
9	C12.H12	156	141	536	936	567

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	581-40-8
2	—	—	—	—	573-98-8
3	—	—	—	—	575-41-7
4	—	—	—	—	571-61-9
5	—	—	—	—	575-37-1
6	—	—	—	—	581-42-0
7	—	—	—	—	575-43-9
8	—	—	—	—	582-16-1
9	—	—	—	—	571-58-4

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:00:00 + 8:38
 SAMPLE: S-MM5-2 1/35A/100M
 COND.S.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30068101 # 776
 CALI: 30068101 # 3
 BASE M/Z: 156
 RIC: 36864.



Library Search Data: 30068101 # 820 Base m/z: 57
 08/31/98 19:00:00 + 9:07 Cali: 30068101 # 3 RIC: 98176.
 Sample: S-MMS-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

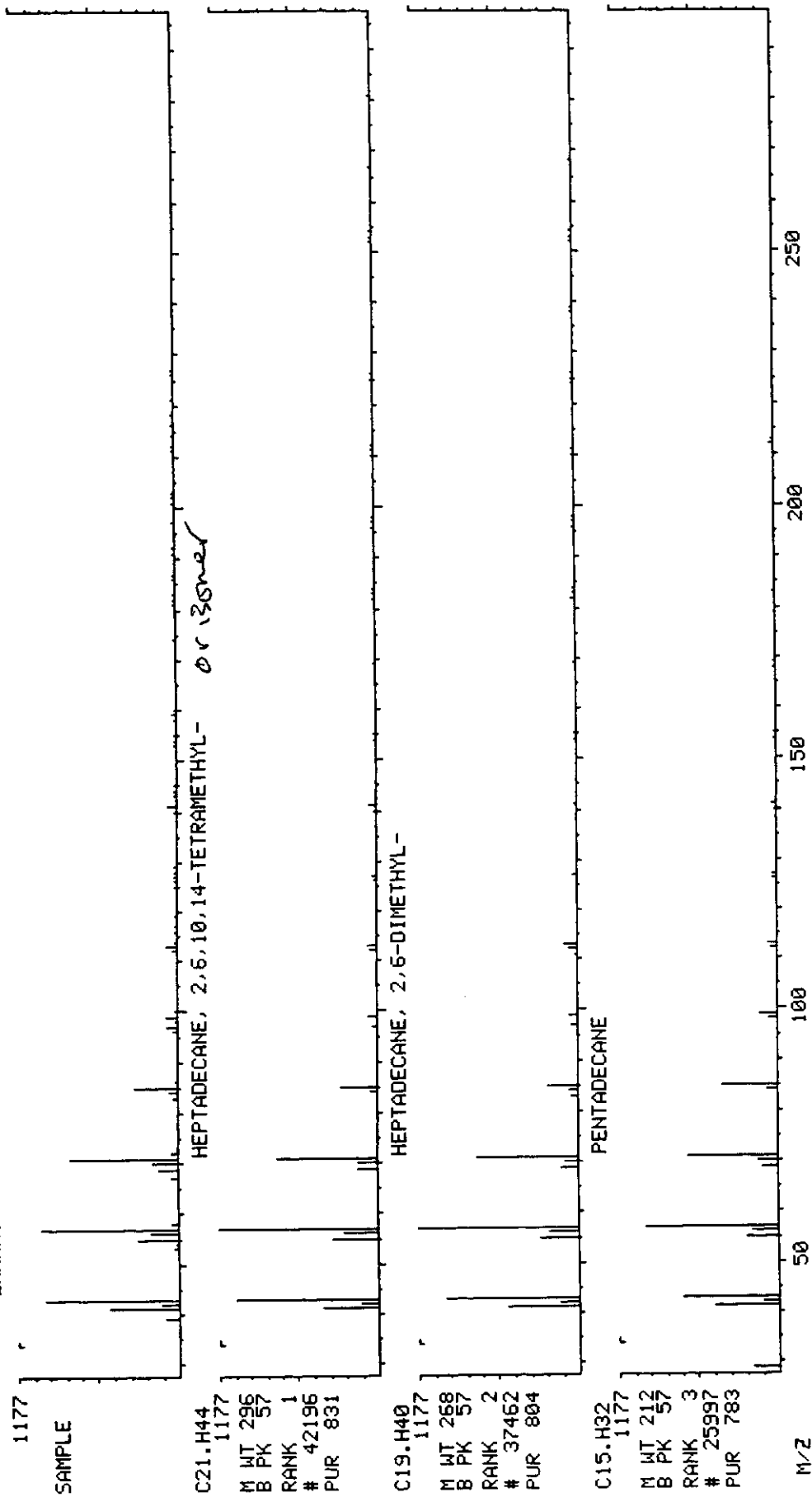
62231 spectra in LIBRARYNB searched for maximum PURITY
 760 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 25997 PENTADECANE
 4 37465 NONADECANE
 5 19016 UNDECANE, 4,7-DIMETHYL-
 6 25991 DODECANE, 2,6,10-TRIMETHYL-
 7 18987 UNDECANE, 2,8-DIMETHYL-
 8 15353 2,6-DIMETHYLDECANE
 9 29263 HEXADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C21.H44	296	57	831	975	842
2	C19.H40	268	57	804	949	823
3	C15.H32	212	57	783	945	795
4	C19.H40	268	57	783	930	815
5	C13.H28	184	43	779	954	780
6	C15.H32	212	57	778	958	781
7	C13.H28	184	43	778	965	782
8	C12.H26	170	43	774	920	783
9	C16.H34	226	57	768	915	784

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	18344-37-1
2	—	—	—	—	54105-67-8
3	—	—	—	—	629-62-9
4	—	—	—	—	629-92-5
5	—	—	—	—	17301-32-5
6	—	—	—	—	3891-98-3
7	—	—	—	—	17301-25-6
8	—	—	—	—	13150-81-7
9	—	—	—	—	544-76-3

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:00:00 + 9:07
 SAMPLE: S-MMS-2 1/35A/100M
 CONDS.: UC/ML *100ML *100Z/100Z *(NA/NA) /1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068101 # 820
 CALI: 30068101 # 3
 BASE M/Z: 57
 RIC: 98176.



Library Search Data: 30068101 # 858 Base m/z: 43
 08/31/98 19:00:00 + 9:33 Cali: 30068101 # 3 RIC: 130176.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 835 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 15352 UNDECANE, 2-METHYL-
 2 19007 DODECANE, 3-METHYL-
 3 25997 PENTADECANE
 4 19006 DECANE, 6-ETHYL-2-METHYL-
 5 18998 UNDECANE, 3,7-DIMETHYL-
 6 22530 TETRADECANE
 7 15353 2,6-DIMETHYLDECANE
 8 15343 DODECANE
 9 37465 NONADECANE

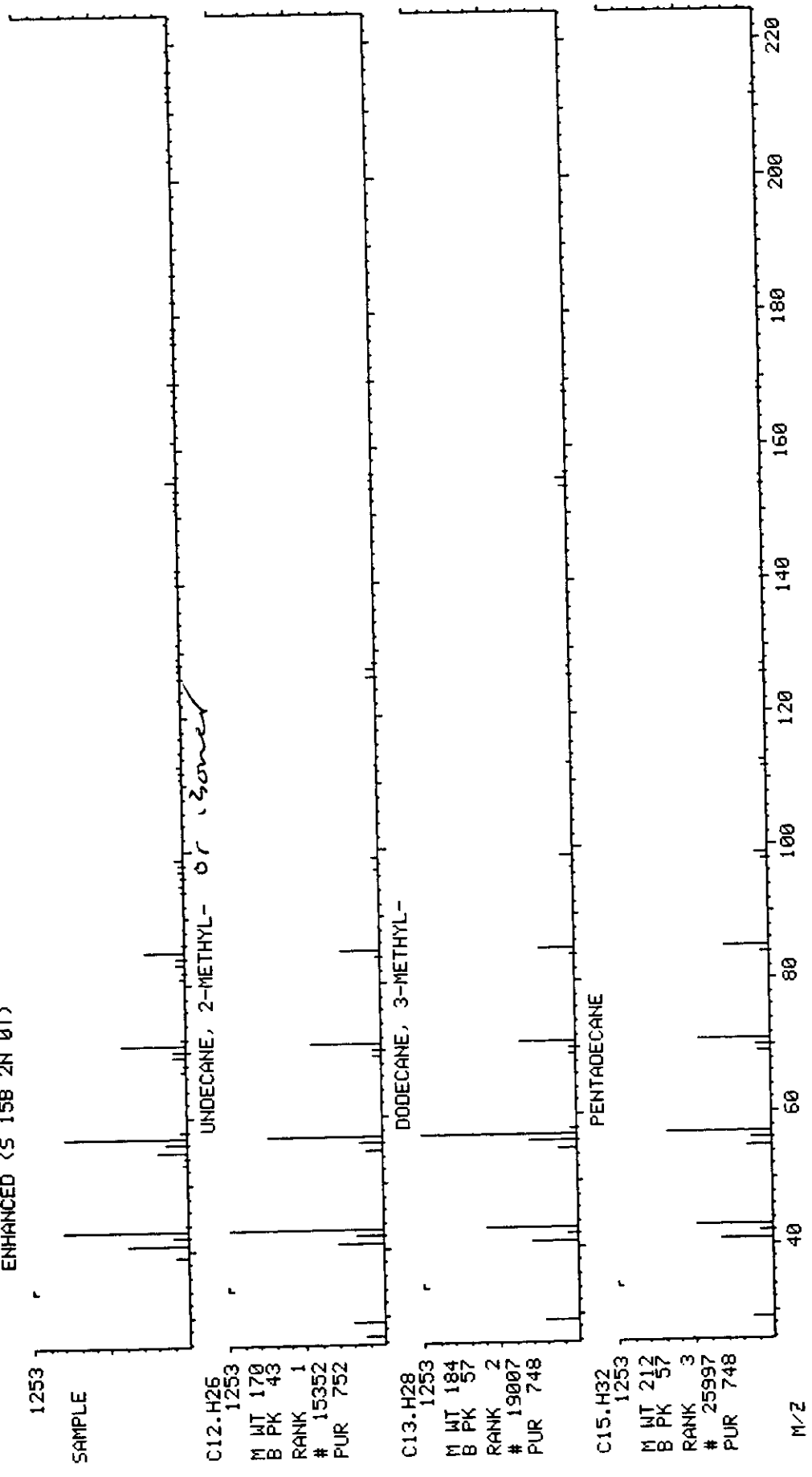
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H26	170	43	752	945	752
2	C13.H28	184	57	748	962	749
3	C15.H32	212	57	748	954	753
4	C13.H28	184	57	742	965	742
5	C13.H28	184	43	738	950	738
6	C14.H30	198	43	738	936	772
7	C12.H26	170	43	737	914	737
8	C12.H26	170	57	734	984	739
9	C19.H40	268	57	732	925	759

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	7045-71-8
2	—	—	—	—	17312-57-1
3	—	—	—	—	629-62-9
4	—	—	—	—	62108-21-8
5	—	—	—	—	17301-29-0
6	—	—	—	—	629-59-4
7	—	—	—	—	13150-81-7
8	—	—	—	—	112-40-3
9	—	—	—	—	629-92-5

DATA: 30068101 # 858
CALI: 30068101 # 3
BASE M/Z: 43
RIC: 130176.

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 19:00:00 + 9:33
SAMPLE: 5-MM5-2 1/35A/100M
CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 158 2N 0T)

INST. ID: F16



Library Search Data: 30068101 # 904 Base m/z: 43
 08/31/98 19:00:00 + 10:03 Cali: 30068101 # 3 RIC: 47040.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 751 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 40233 NONADECANOL
 2 42521 1-EICOSANOL
 3 42180 OCTADECANE, 1-(ETHENYLOXY)-
 4 15969 HYDROXYLAMINE, O-DECYL-
 5 25971 OXIRANE, DODECYL-
 6 58380 DECANEDIOIC ACID, DIDECYL ESTER
 7 37444 OXIRANE, HEXADECYL-
 8 37449 OCTADECANAL
 9 19520 1-OCTANOL, 2-BUTYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40.O	284	43	674	938	703
2	C20.H42.O	298	43	670	931	705
3	C20.H40.O	296	43	647	941	677
4	C10.H23.O.N	173	43	638	906	649
5	C14.H28.O	212	41	636	902	648
6	C30.H58.O4	482	57	634	919	649
7	C18.H36.O	268	57	634	841	687
8	C18.H36.O	268	43	633	852	687
9	C12.H26.O	186	57	633	921	642

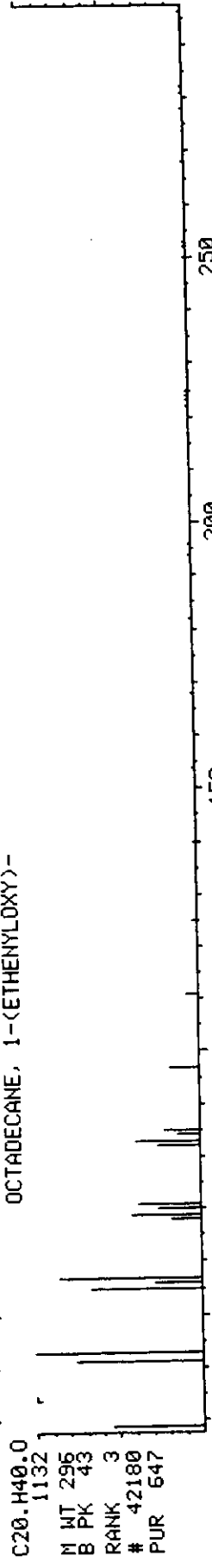
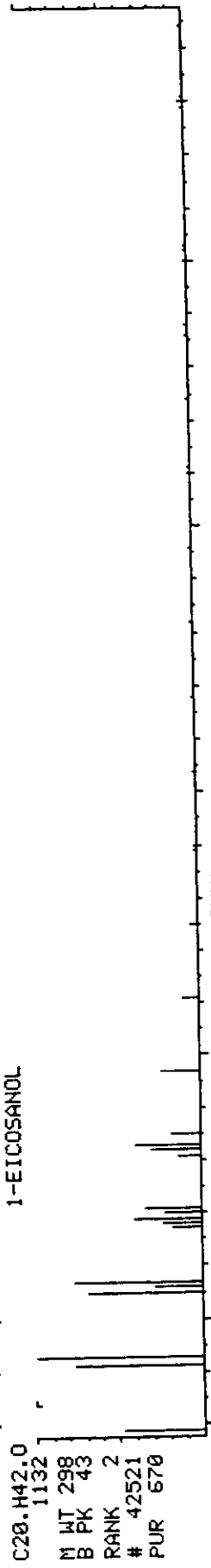
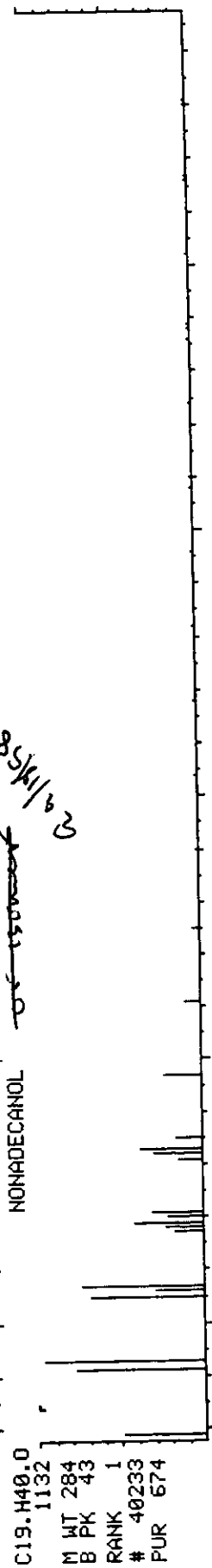
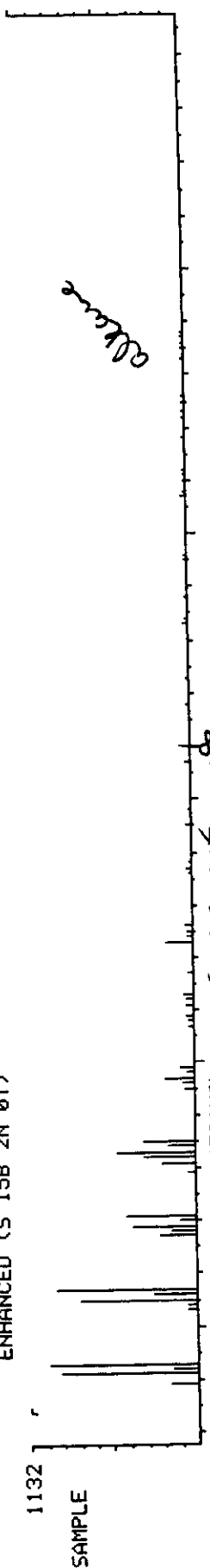
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	52783-43-4
2	—	—	—	—	629-96-9
3	—	—	—	—	930-02-9
4	—	—	—	—	29812-79-1
5	—	—	—	—	3234-28-4
6	—	—	—	—	2432-89-5
7	—	—	—	—	7390-81-0
8	—	—	—	—	638-66-4
9	—	—	—	—	3913-02-8

DATA: 30068101 # 904
CALI: 30068101 # 3

BASE M/Z: 43
RIC: 47040.

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 19:00:00 + 10:03
SAMPLE: S-MM5-2 1/35A/100M
CONDOS.: UG/ML *100ML *100Z/100Z *(NA/NA >)/1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



M/Z

Library Search Data: 30068101 # 948 Base m/z: 57
 08/31/98 19:00:00 + 10:33 Cali: 30068101 # 3 RIC: 114816.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 343 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 15969 HYDROXYLAMINE, O-DECYL-
 4 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 5 25997 PENTADECANE
 6 19016 UNDECANE, 4,7-DIMETHYL-
 7 25994 DODECANE, 2,6,11-TRIMETHYL-
 8 26001 DODECANE, 2,7,10-TRIMETHYL-
 9 19523 1-DECANOL, 2-ETHYL-

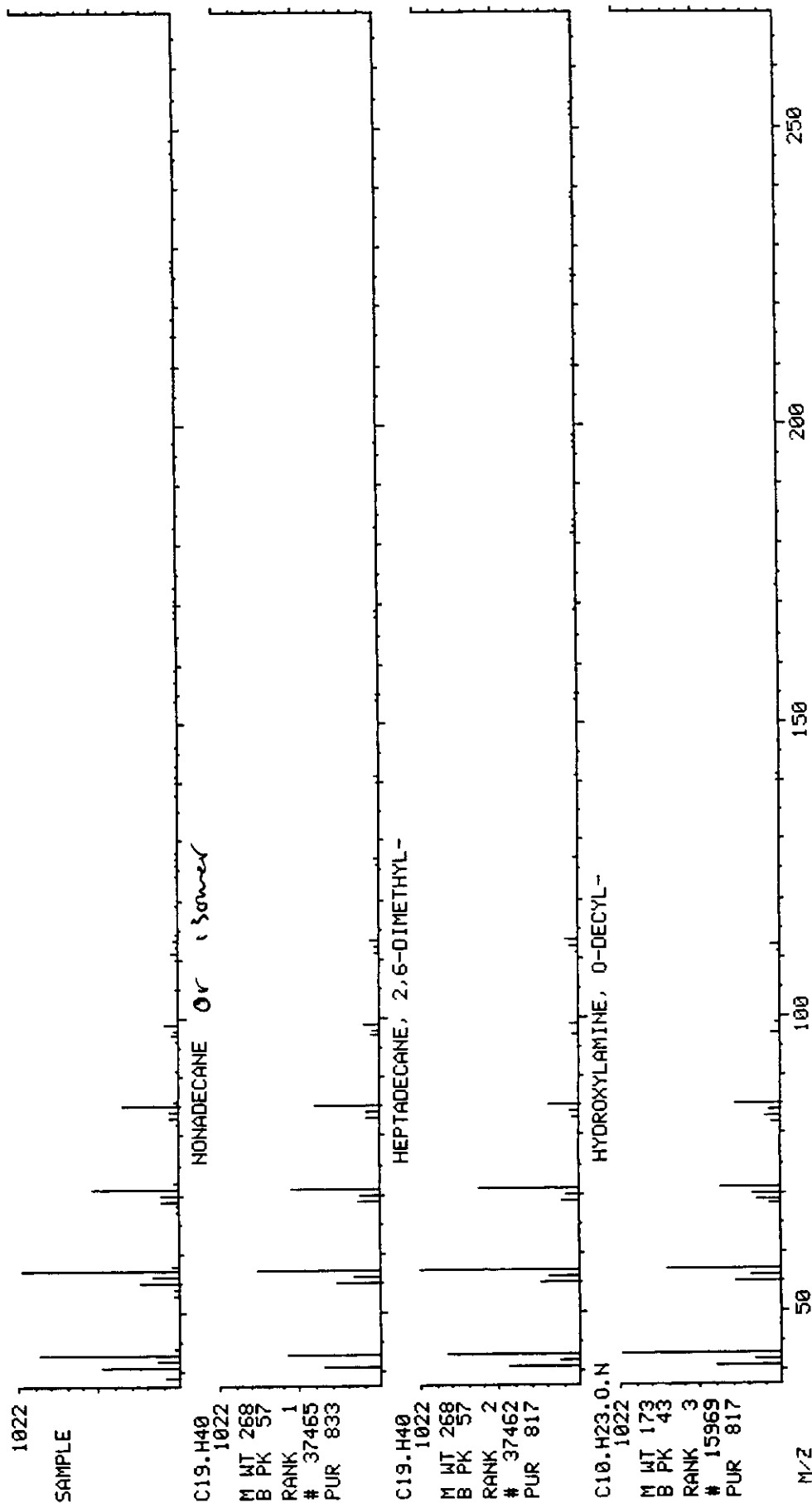
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	833	956	854
2	C19.H40	268	57	817	936	844
3	C10.H23.O.N	173	43	817	966	817
4	C21.H44	296	57	806	944	830
5	C15.H32	212	57	803	949	833
6	C13.H28	184	43	799	966	800
7	C15.H32	212	57	792	961	795
8	C15.H32	212	57	783	956	788
9	C12.H26.O	186	57	782	925	795

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	29812-79-1
4	—	—	—	—	18344-37-1
5	—	—	—	—	629-62-9
6	—	—	—	—	17301-32-5
7	—	—	—	—	31295-56-4
8	—	—	—	—	74645-98-0
9	—	—	—	—	21078-65-9

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:00:00 + 10:33
 SAMPLE: S-MM5-2 1/35A/100M
 COND.: UG/ML *100ML *100Z/100Z *(NA/NA) 1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30058101 # 948
 CALI: 30058101 # 3

BASE M/Z: 57
 RIC: 114816.



Library Search Data: 30068101 # 987 Base m/z: 57
 08/31/98 19:00:00 + 10:59 Cali: 30068101 # 3 RIC: 96384.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 644 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 37465 NONADECANE
 3 25997 PENTADECANE
 4 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 5 25996 TRIDECANE, 4,8-DIMETHYL-
 6 22530 TETRADECANE
 7 32059 HEPTADECANE
 8 29263 HEXADECANE
 9 22527 DODECANE, 2,5-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	817	949	825
2	C19.H40	268	57	812	938	831
3	C15.H32	212	57	806	952	810
4	C21.H44	296	57	805	951	812
5	C15.H32	212	57	803	936	814
6	C14.H30	198	43	801	942	806
7	C17.H36	240	57	801	935	811
8	C16.H34	226	57	799	934	808
9	C14.H30	198	57	796	942	797

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	629-92-5
3	—	—	—	—	629-62-9
4	—	—	—	—	18344-37-1
5	—	—	—	—	55030-62-1
6	—	—	—	—	629-59-4
7	—	—	—	—	629-78-7
8	—	—	—	—	544-76-3
9	—	—	—	—	56292-65-0

BASE M/Z: 57
RIC: 96384.

DATA: 30068101 # 987
CALI: 30068101 # 3

MID LIBRARY SEARCH <LIBRARYNB>

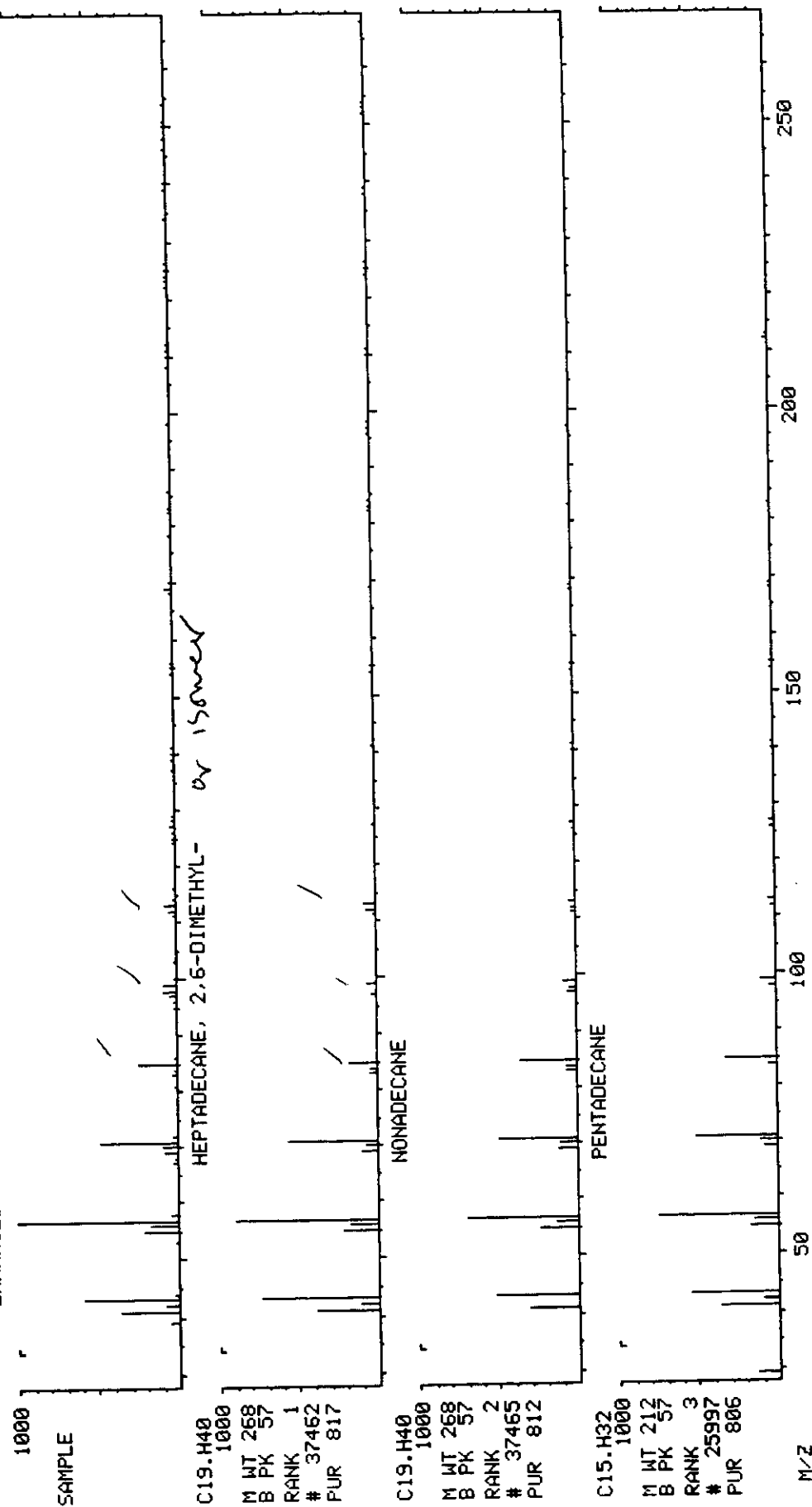
08/31/98 19:00:00 + 10:59

SAMPLE: S-MM5-2 1/35A/100M

COND.: UG/ML *100ML *100Z/100Z *(NA/NA) /1/35A NA M

ENHANCED <S 158 2N 0T>

INST. ID: F16



Library Search Data: 30068101 #1035 Base m/z: 57
 08/31/98 19:00:00 + 11:31 Cali: 30068101 # 3 RIC: 149504.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 713 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 3 37465 NONADECANE
 4 25991 DODECANE, 2,6,10-TRIMETHYL-
 5 22535 DODECANE, 4,6-DIMETHYL-
 6 26001 DODECANE, 2,7,10-TRIMETHYL-
 7 37466 PENTADECANE, 2,6,10,14-TETRAMETHYL-
 8 25994 DODECANE, 2,6,11-TRIMETHYL-
 9 46161 TRICOSANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	872	975	879
2	C21.H44	296	57	836	950	860
3	C19.H40	268	57	827	944	844
4	C15.H32	212	57	822	972	826
5	C14.H30	198	57	821	967	822
6	C15.H32	212	57	821	971	825
7	C19.H40	268	71	815	941	833
8	C15.H32	212	57	802	948	810
9	C23.H48	324	43	794	892	858

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	18344-37-1
3	—	—	—	—	629-92-5
4	—	—	—	—	3891-98-3
5	—	—	—	—	61141-72-8
6	—	—	—	—	74645-98-0
7	—	—	—	—	1921-70-6
8	—	—	—	—	31295-56-4
9	—	—	—	—	638-67-5

BASE M/Z: 57
RIC: 149504.

DATA: 30068101 #1035
CALI: 30068101 # 3

MID LIBRARY SEARCH (LIBRARYNB)

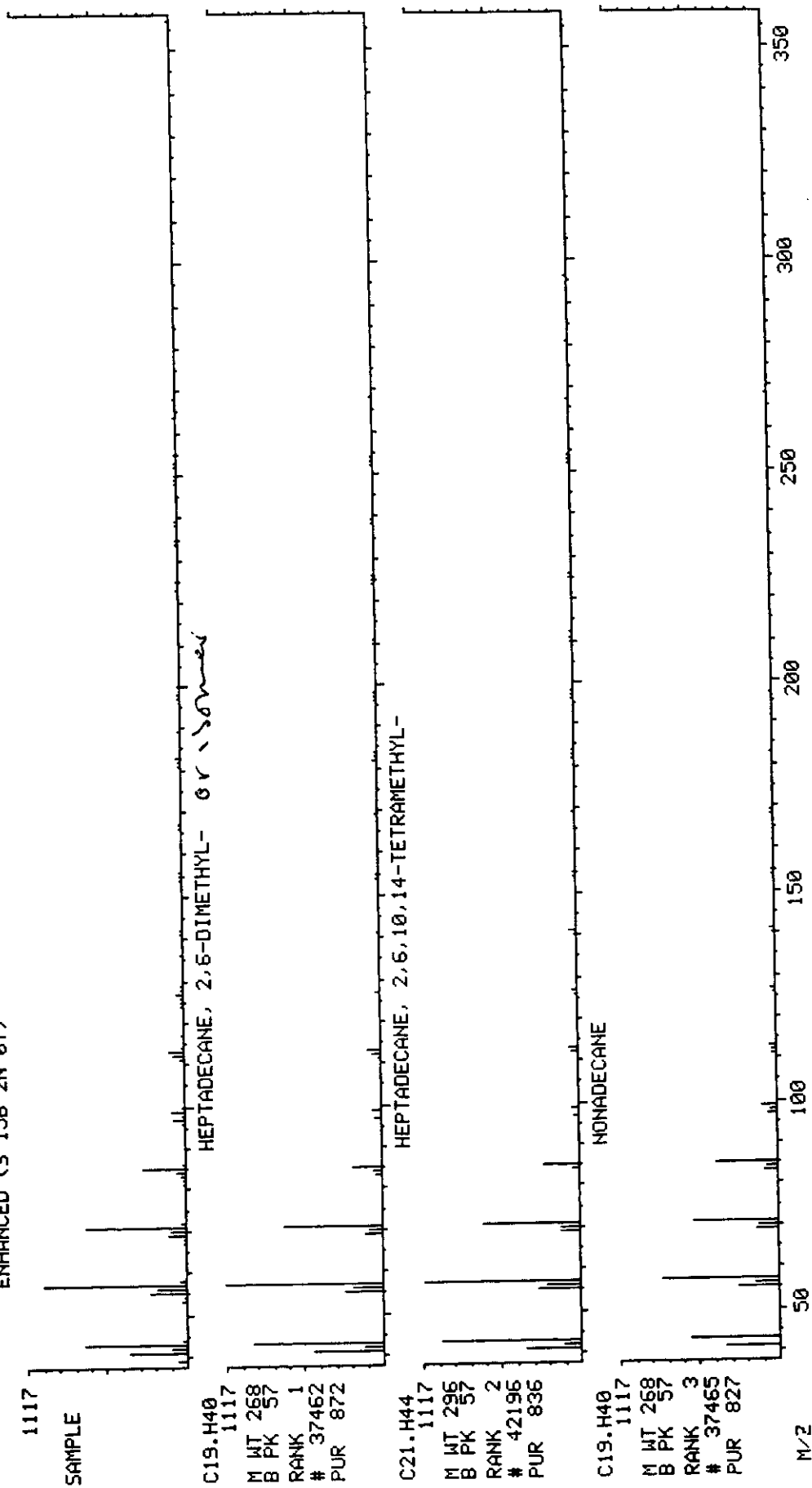
08/31/98 19:00:00 + 11:31

SAMPLE: S-MMS-2 1/35A/100M

COND.: UG/ML *100ML *100%/100% *(NA/NA) /1/35A NA M

ENHANCED (5 158 2N 0T)

INST. ID: F16



Library Search Data: 30068101 #1114 Base m/z: 57
 08/31/98 19:00:00 + 12:23 Cali: 30068101 # 3 RIC: 60416.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 651 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 37465 NONADECANE
 3 19016 UNDECANE, 4,7-DIMETHYL-
 4 19006 DECANE, 6-ETHYL-2-METHYL-
 5 19056 UNDECANE, 2,7-DIMETHYL-
 6 8102 HEXANE, 2,2,3,3-TETRAMETHYL-
 7 19028 DECANE, 2,4,6-TRIMETHYL-
 8 15356 DECANE, 2,4-DIMETHYL-
 9 22530 TETRADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	791	910	847
2	C19.H40	268	57	775	896	852
3	C13.H28	184	43	769	922	805
4	C13.H28	184	57	768	934	799
5	C13.H28	184	43	767	942	779
6	C10.H22	142	57	767	920	774
7	C13.H28	184	43	767	938	789
8	C12.H26	170	43	765	944	776
9	C14.H30	198	43	763	899	841

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	629-92-5
3	—	—	—	—	17301-32-5
4	—	—	—	—	62108-21-8
5	—	—	—	—	17301-24-5
6	—	—	—	—	13475-81-5
7	—	—	—	—	62108-27-4
8	—	—	—	—	2801-84-5
9	—	—	—	—	629-59-4

BASE M/Z: 57
RIC: 60416.

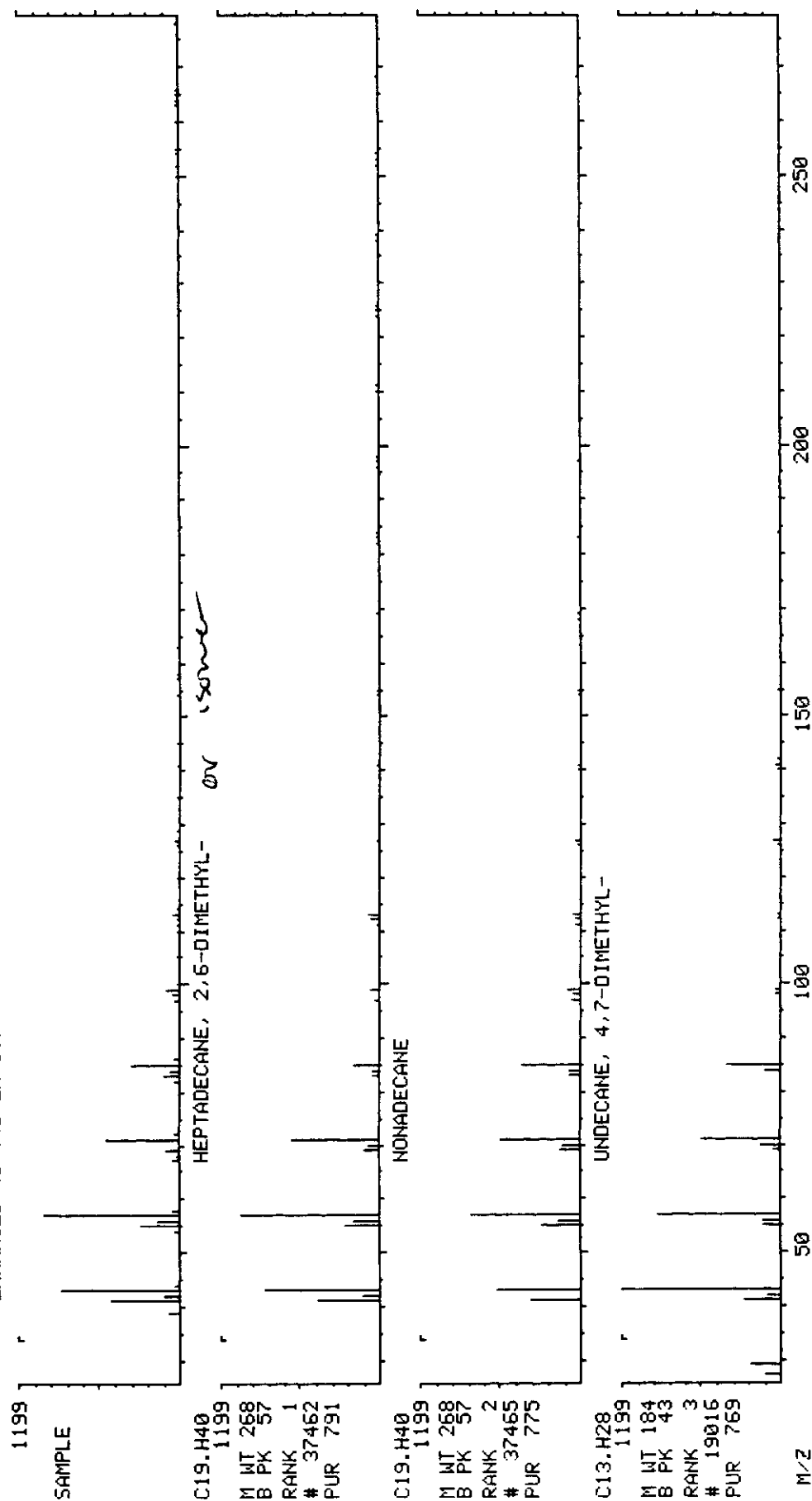
MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 19:00:00 + 12:23

SAMPLE: S-MM5-2 1/35A/100M

SAMPLE: S-MM5-2 1/35A/100M INST. ID: F16

SAMPLE: S-MM5-2 1/35A/100M INST. ID: F16
CONDOS.: UG/ML *100ML *100%/100% *(NA/NA)/1/35A NA M



Library Search Data: 30068101 #1119 Base m/z: 57
 08/31/98 19:00:00 + 12:27 Cali: 30068101 # 3 RIC: 88576.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

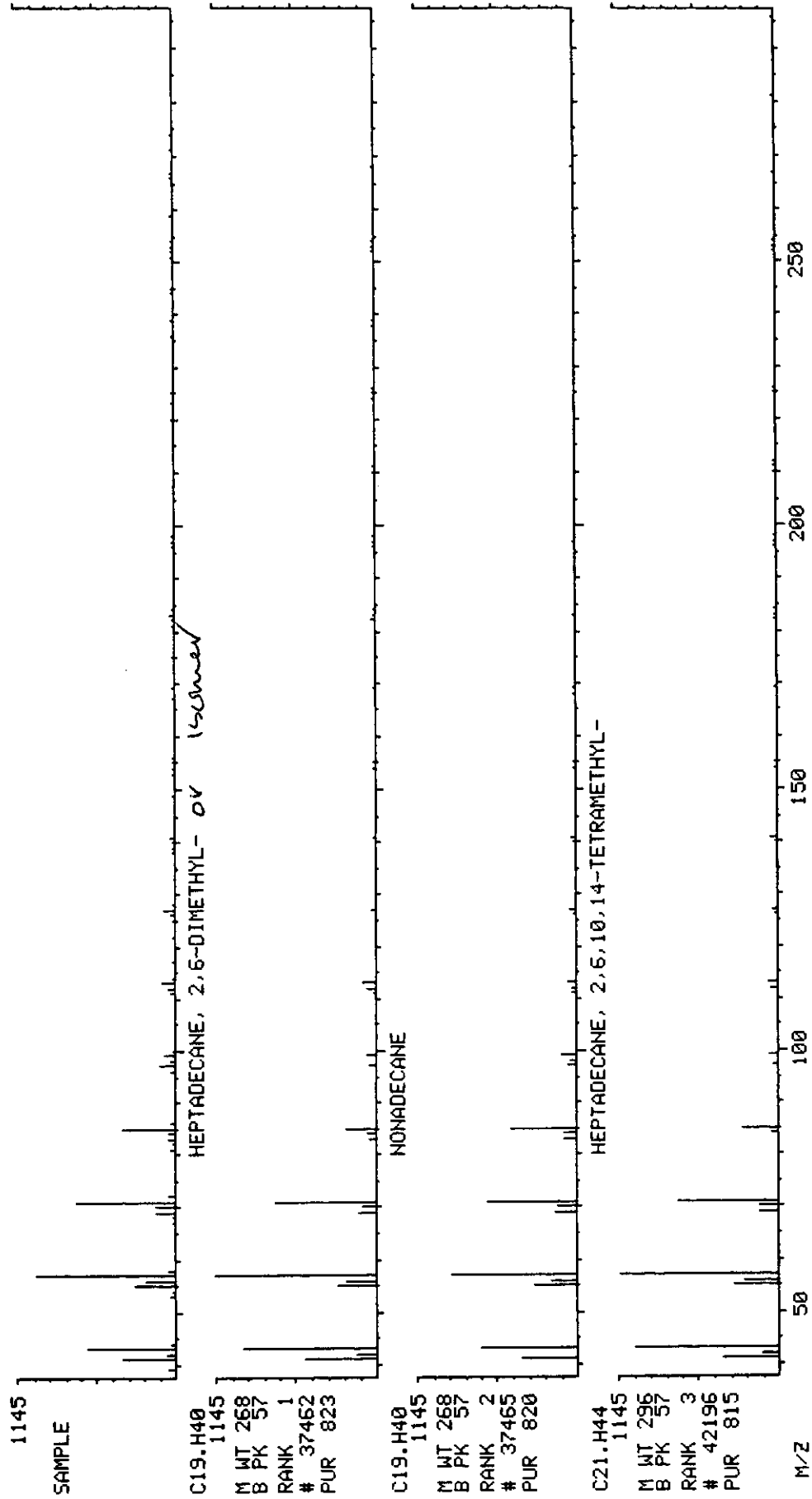
62231 spectra in LIBRARYNB searched for maximum PURITY
 590 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 37465 NONADECANE
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 46161 TRICOSANE
 5 25991 DODECANE, 2,6,10-TRIMETHYL-
 6 37466 PENTADECANE, 2,6,10,14-TETRAMETHYL-
 7 42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-
 8 26001 DODECANE, 2,7,10-TRIMETHYL-
 9 32059 HEPTADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	823	950	828
2	C19.H40	268	57	820	956	838
3	C21.H44	296	57	815	944	828
4	C23.H48	324	43	804	921	845
5	C15.H32	212	57	797	970	798
6	C19.H40	268	71	795	939	801
7	C21.H44	296	57	794	956	819
8	C15.H32	212	57	789	971	789
9	C17.H36	240	57	788	927	820

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	629-92-5
3	—	—	—	—	18344-37-1
4	—	—	—	—	638-67-5
5	—	—	—	—	3891-98-3
6	—	—	—	—	1921-70-6
7	—	—	—	—	54833-48-6
8	—	—	—	—	74645-98-0
9	—	—	—	—	629-78-7

MID LIBRARY SEARCH <LIBRARYNB>
 08/31/98 19:00:00 + 12:27
 SAMPLE: S-MM5-2 1/35A/100M
 CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068101 #1119
 CALI: 30068101 # 3
 BASE M/Z: 57
 RIC: 88576.



Library Search Data: 30068101 #1191 Base m/z: 57
 08/31/98 19:00:00 + 13:15 Cal: 30068101 # 3 RIC: 54912.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 294 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 15969 HYDROXYLAMINE, O-DECYL-
 5 25997 PENTADECANE
 6 46161 TRICOSANE
 7 19016 UNDECANE, 4,7-DIMETHYL-
 8 25994 DODECANE, 2,6,11-TRIMETHYL-
 9 19523 1-DECANOL, 2-ETHYL-

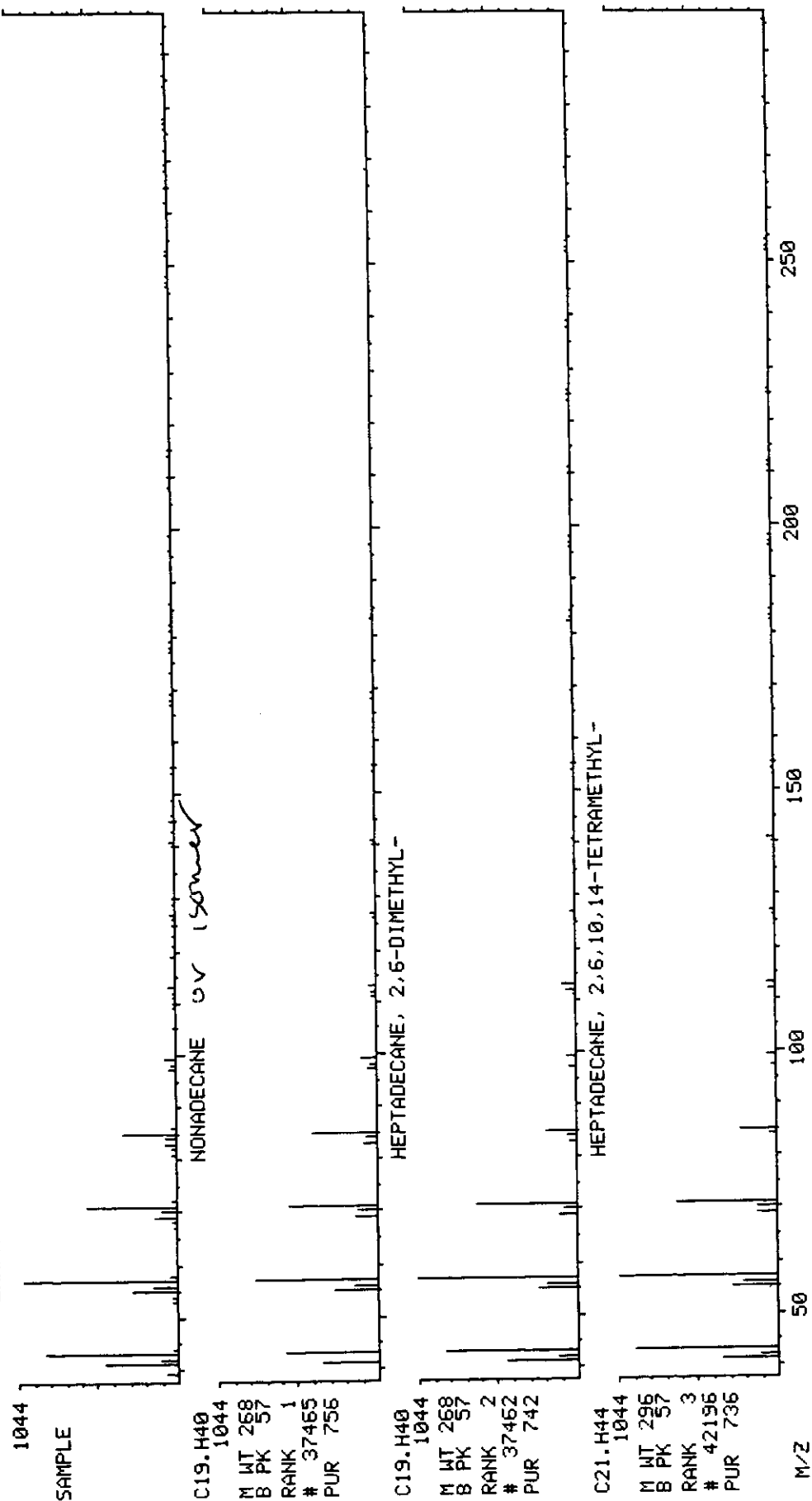
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	756	969	766
2	C19.H40	268	57	742	941	760
3	C21.H44	296	57	736	952	758
4	C10.H23.O.N	173	43	729	968	729
5	C15.H32	212	57	725	958	747
6	C23.H48	324	43	724	910	786
7	C13.H28	184	43	719	962	720
8	C15.H32	212	57	706	958	709
9	C12.H26.O	186	57	703	937	707

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	18344-37-1
4	—	—	—	—	29812-79-1
5	—	—	—	—	629-62-9
6	—	—	—	—	638-67-5
7	—	—	—	—	17301-32-5
8	—	—	—	—	31295-56-4
9	—	—	—	—	21078-65-9

DATA: 30068101 #1191
CALI: 30068101 # 3
BASE M/Z: 57
RIC: 54912.

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 19:00:00 + 13:15
SAMPLE: S-MMS-2 1/35A/100M
COND.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068101 #1264 Base m/z: 57
 08/31/98 19:00:00 + 14:04 Cali: 30068101 # 3 RIC: 45120.
 Sample: S-MMS-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

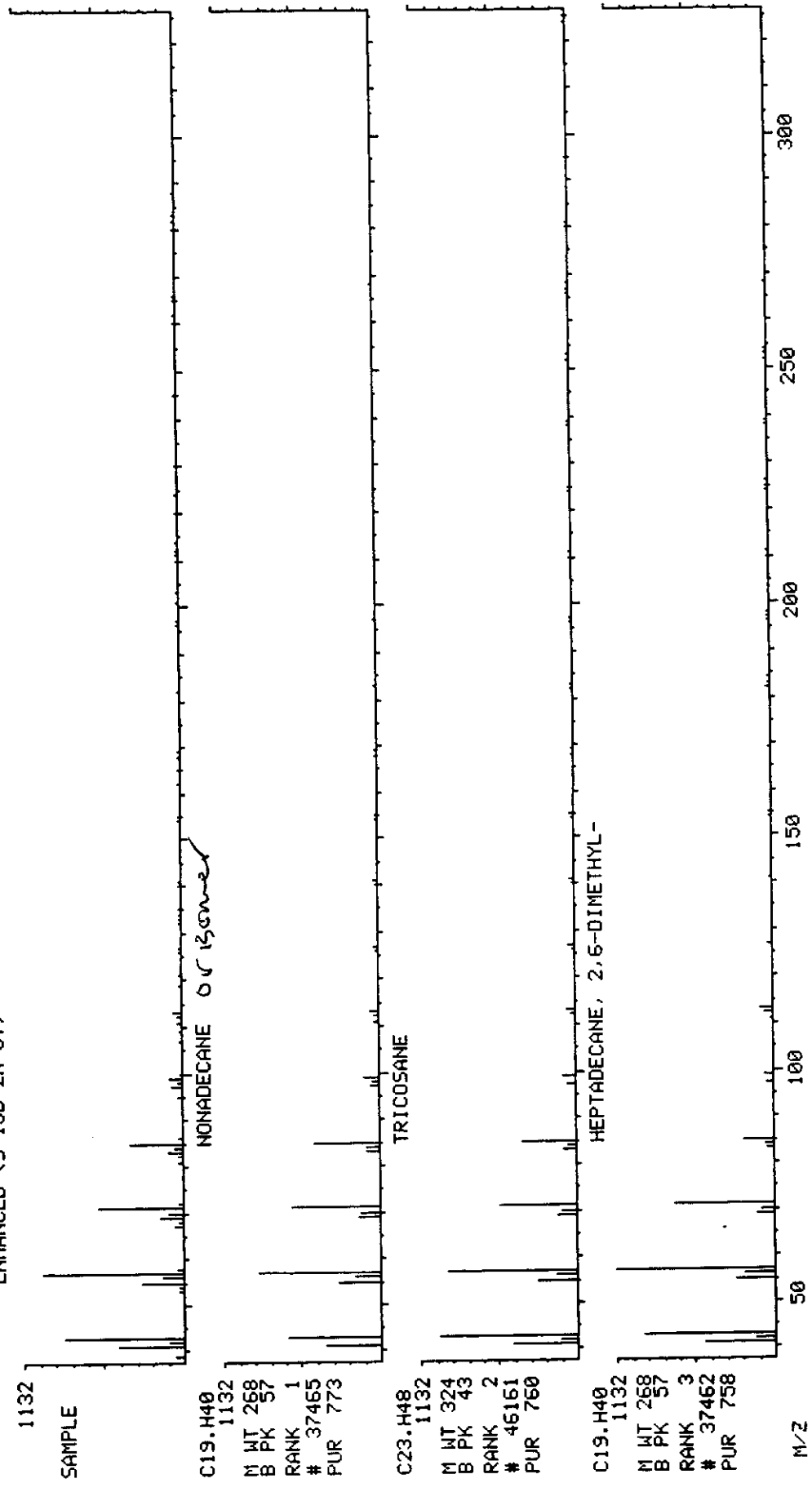
62231 spectra in LIBRARYNB searched for maximum PURITY
 298 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 46161 TRICOSANE
 3 37462 HEPTADECANE, 2,6-DIMETHYL-
 4 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 5 25997 PENTADECANE
 6 49555 PENTACOSANE
 7 42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-
 8 32059 HEPTADECANE
 9 15969 HYDROXYLAMINE, O-DECYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	773	969	791
2	C23.H48	324	43	760	923	816
3	C19.H40	268	57	758	931	778
4	C21.H44	296	57	755	947	771
5	C15.H32	212	57	748	982	753
6	C25.H52	352	43	745	901	819
7	C21.H44	296	57	731	932	759
8	C17.H36	240	57	731	928	759
9	C10.H23.O.N	173	43	730	967	732

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	638-67-5
3	—	—	—	—	54105-67-8
4	—	—	—	—	18344-37-1
5	—	—	—	—	629-62-9
6	—	—	—	—	629-99-2
7	—	—	—	—	54833-48-6
8	—	—	—	—	629-78-7
9	—	—	—	—	29812-79-1

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:00:00 + 14:04
 SAMPLE: 5-MMS-2 1/35A/100M
 CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 01)
 DATA: 30068101 #1264
 CALI: 30068101 # 3
 BASE M/Z: 57
 RIC: 45120.



Library Search Data: 30068101 #1335 Base m/z: 57
 08/31/98 19:00:00 + 14:51 Cali: 30068101 # 3 RIC: 29728.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 474 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 2 37456 2-METHYLOCTADECANE
 3 37462 HEPTADECANE, 2,6-DIMETHYL-
 4 29263 HEXADECANE
 5 37465 NONADECANE
 6 25997 PENTADECANE
 7 32058 HEXADECANE, 3-METHYL-
 8 46161 TRICOSANE
 9 42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C21.H44	296	57	691	910	713
2	C19.H40	268	43	689	925	709
3	C19.H40	268	57	683	889	712
4	C16.H34	226	57	683	939	700
5	C19.H40	268	57	682	951	707
6	C15.H32	212	57	672	968	680
7	C17.H36	240	57	664	919	691
8	C23.H48	324	43	660	896	725
9	C21.H44	296	57	657	922	687

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	18344-37-1
2	—	—	—	—	- -
3	—	—	—	—	54105-67-8
4	—	—	—	—	544-76-3
5	—	—	—	—	629-92-5
6	—	—	—	—	629-62-9
7	—	—	—	—	6418-43-5
8	—	—	—	—	638-67-5
9	—	—	—	—	54833-48-6

BASE M/Z: 57
RIC: 29728.

DATA: 30068101 #1335
CALI: 30068101 # 3

MID LIBRARY SEARCH (LIBRARYNB)

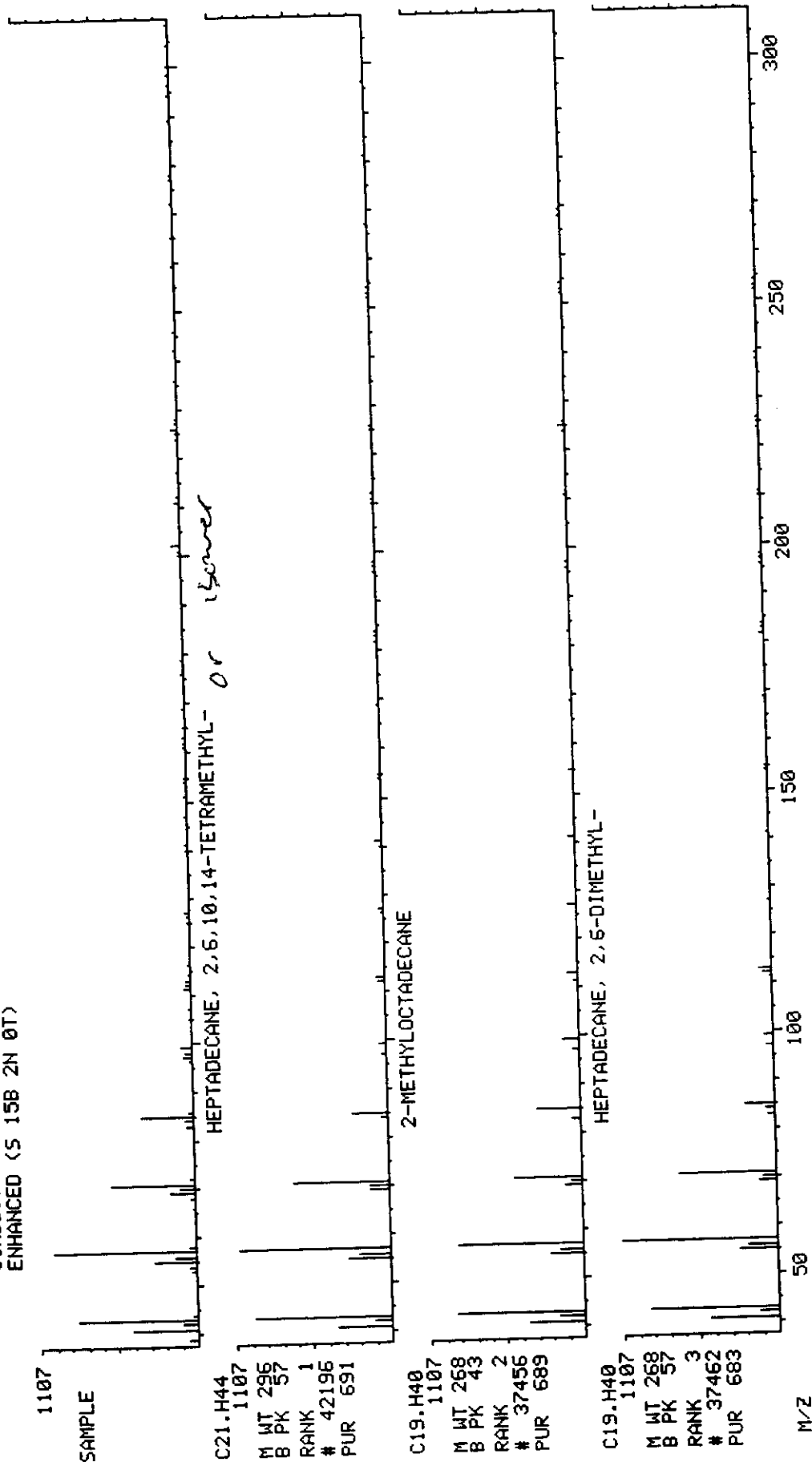
08/31/98 19:00:00 + 14:51

SAMPLE: S-MM5-2 1/35A/100M

COND.S.: UG/ML *100ML *100Z/100Z *(NA/NA) >1/35A NA M

ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068101 #1402 Base m/z: 57
 08/31/98 19:00:00 + 15:36 Cali: 30068101 # 3 RIC: 23168.
 Sample: S-MM5-2 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 850 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 46161 TRICOSANE
 5 37456 2-METHYLOCTADECANE
 6 25997 PENTADECANE
 7 42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-
 8 15969 HYDROXYLAMINE, O-DECYL-
 9 49555 PENTACOSANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	716	961	738
2	C19.H40	268	57	702	934	715
3	C21.H44	296	57	696	945	709
4	C23.H48	324	43	694	920	749
5	C19.H40	268	43	685	942	711
6	C15.H32	212	57	684	967	691
7	C21.H44	296	57	680	940	703
8	C10.H23.O.N	173	43	676	966	678
9	C25.H52	352	43	673	885	756

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	629-92-5
2	---	---	---	---	54105-67-8
3	---	---	---	---	18344-37-1
4	---	---	---	---	638-67-5
5	---	---	---	---	- -
6	---	---	---	---	629-62-9
7	---	---	---	---	54833-48-6
8	---	---	---	---	29812-79-1
9	---	---	---	---	629-99-2

BASE M/Z: 57
RIC: 23158.

DATA: 30068101 #1402
CALI: 30068101 # 3

MID LIBRARY SEARCH (LIBRARYNB)

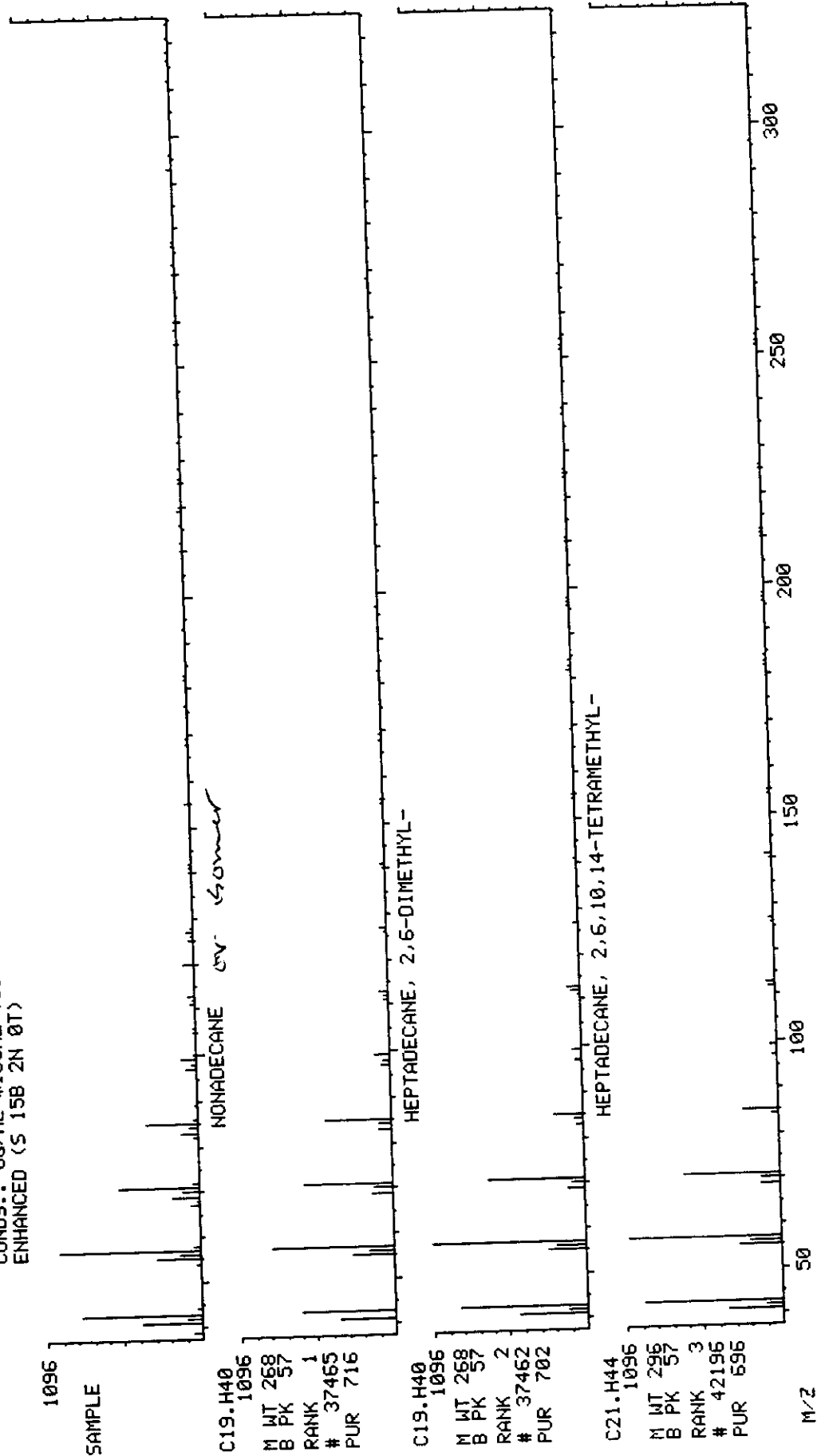
08/31/98 19:00:00 + 15:36

SAMPLE: S-MM5-2 1/35A/100M

COND.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M

ENHANCED (S 158 2N 0T)

INST. ID: F16



TIC SELECTION REPORT

DATA FILE: 30068101

THE FOLLOWING PEAKS WERE REJECTED BECAUSE
AT LEAST 40 % OF THEIR SIZE WAS ACCOUNTED FOR BY
TARGET COMPOUNDS ELUTING WITHIN 4 SCANS OF THE
PEAK TOP.

SCAN	SIZE	AMOUNT
341	170240.	22.662
530	300480.	40.000
829	283264.	40.000
1087	334464.	40.000
1552	199007.	40.000
1783	185564.	40.000

TOTAL NUMBER OF UNIDENTIFIED PEAKS WITH SIZE
GREATER THAN 10 % OF THE CLOSEST INTERNAL STANDARD
THAT DOES NOT HAVE INTERFERENCES = 94

INTERNAL STANDARDS THAT HAVE RIC SIZE LESS THAN
50 % OR GREATER THAN 200 % OF THE ESTIMATED RIC SIZE
ARE CONSIDERED TO HAVE INTERFERENCES AND WILL NOT BE USED
FOR QUANTITATION.

#	INTERNAL STANDARD	PERCENT OF ESTIMATED		SCAN
		RIC SIZE	RIC SIZE	
1	C140 NAPHTHALENE-D8	300480.	196	530
2	C150 ACENAPHTHENE-D10	283264.	147	829
3	C160 PHENANTHRENE-D10	334464.	154	1087
4	C170 CHRYSENE-D12	199007.	142	1552
5	C175 PERYLENE-D12	185564.	154	1783

* INDICATES INTERFERENCE

SIZE = AREA

Semivolatile Organics
Method 0010/8270

Client Name: Pacific Environmental Services
Client ID: S-MM5-1B-F, FH, XAD, COND, BH
LAB ID: 300681-0002-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 25 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 100

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
Phenol	ND	ug/Sample	3000	RG
bis(2-Chloroethyl)ether	ND	ug/Sample	3000	
2-Chlorophenol	ND	ug/Sample	3000	
1,3-Dichlorobenzene	ND	ug/Sample	3000	
1,4-Dichlorobenzene	ND	ug/Sample	3000	
Benzyl alcohol	ND	ug/Sample	3000	
1,2-Dichlorobenzene	ND	ug/Sample	3000	
2-Methylphenol	ND	ug/Sample	3000	
2,2'-Oxybis(1-chloropropane)	ND	ug/Sample	3000	
3/4-Methylphenol	ND	ug/Sample	3000	
N-Nitroso-di-n-propylamine	ND	ug/Sample	3000	
Hexachloroethane	ND	ug/Sample	3000	
Nitrobenzene	ND	ug/Sample	3000	
Isophorone	ND	ug/Sample	3000	
2-Nitrophenol	ND	ug/Sample	3000	
2,4-Dimethylphenol	ND	ug/Sample	3000	
Benzoic acid	ND	ug/Sample	15000	
bis(2-Chloroethoxy)-methane	ND	ug/Sample	3000	
2,4-Dichlorophenol	ND	ug/Sample	3000	
1,2,4-Trichlorobenzene	ND	ug/Sample	3000	
Naphthalene	ND	ug/Sample	3000	
4-Chloroaniline	ND	ug/Sample	3000	
Hexachlorobutadiene	ND	ug/Sample	3000	
4-Chloro-3-methylphenol	ND	ug/Sample	3000	
2-Methylnaphthalene	4700	ug/Sample	3000	
Hexachlorocyclopentadiene	ND	ug/Sample	3000	
2,4,6-Trichlorophenol	ND	ug/Sample	3000	
2,4,5-Trichlorophenol	ND	ug/Sample	15000	
2-Chloronaphthalene	ND	ug/Sample	3000	
2-Nitroaniline	ND	ug/Sample	3000	
Dimethyl phthalate	ND	ug/Sample	3000	
Acenaphthylene	ND	ug/Sample	15000	
3-Nitroaniline	ND	ug/Sample	3000	
Acenaphthene	ND	ug/Sample	15000	
2,4-Dinitrophenol	ND	ug/Sample	15000	
4-Nitrophenol	ND	ug/Sample	15000	

Note G = Reporting limit(s) raised due to matrix interference.
Note R = Reporting limit(s) raised due to sample volume limitations.
ND = Not Detected

724

Reported By: Emily Uebelhoer

Approved By: Mike Orbanosky

The cover letter is an integral part of this report.
Rev 230787

Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: S-MM5-1B-F, FH, XAD, COND, BH
LAB ID: 300681-0002-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 25 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 100

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
Dibenzofuran	ND	ug/Sample	3000	
2,4-Dinitrotoluene	ND	ug/Sample	3000	
2,6-Dinitrotoluene	ND	ug/Sample	3000	
Diethyl phthalate	ND	ug/Sample	3000	
4-Chlorophenyl phenyl ether	ND	ug/Sample	3000	
Fluorene	ND	ug/Sample	3000	
4-Nitroaniline	ND	ug/Sample	15000	
4,6-Dinitro-2-methylphenol	ND	ug/Sample	15000	
N-Nitrosodiphenylamine	ND	ug/Sample	3000	
4-Bromophenyl phenyl ether	ND	ug/Sample	3000	
Hexachlorobenzene	ND	ug/Sample	3000	
Pentachlorophenol	ND	ug/Sample	15000	
Phenanthrene	ND	ug/Sample	3000	
Anthracene	ND	ug/Sample	3000	
Di-n-butyl phthalate	ND	ug/Sample	3000	
Fluoranthene	ND	ug/Sample	3000	
Pyrene	ND	ug/Sample	3000	
Butyl benzyl phthalate	ND	ug/Sample	3000	
3,3'-Dichlorobenzidine	ND	ug/Sample	6000	
Benzo(a)anthracene	ND	ug/Sample	3000	
bis(2-Ethylhexyl)-phthalate	ND	ug/Sample	3000	
Chrysene	ND	ug/Sample	3000	
Di-n-octyl phthalate	ND	ug/Sample	3000	
Benzo(b)fluoranthene	ND	ug/Sample	3000	
Benzo(k)fluoranthene	ND	ug/Sample	3000	
Benzo(a)pyrene	ND	ug/Sample	3000	
Indeno(1,2,3-cd)pyrene	ND	ug/Sample	3000	
Dibenz(a,h)anthracene	ND	ug/Sample	3000	
Benzo(g,h,i)perylene	ND	ug/Sample	3000	
Acetophenone	ND	ug/Sample	3000	
4-Aminobiphenyl	ND	ug/Sample	15000	
Aniline	ND	ug/Sample	3000	
Benzidine	ND	ug/Sample	30000	
3,3'-Dimethylbenzidine	ND	ug/Sample	6000	
N-Nitrosodimethylamine	ND	ug/Sample	3000	
N-Nitrosomorpholine	ND	ug/Sample	3000	
Pentachloronitrobenzene (PCNB)	ND	ug/Sample	15000	
o-Toluidine	ND	ug/Sample	6000	

735

ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Mike Orbanosky

The cover letter is an integral part of this report.
Rev 230787

Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: S-MM5-1B-F, FH, XAD, COND, BH
LAB ID: 300681-0002-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 25 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 100

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
2-Methoxybenzenamine	ND	ug/Sample	--	
Biphenyl	ND	ug/Sample	--	
Chloroacetophenone	ND	ug/Sample	--	
Cumene	ND	ug/Sample	--	
DBCP (1,2-Dibromo-3-chloropropane)	ND	ug/Sample	--	
Benzo(e)pyrene	ND	ug/Sample	--	
N-N-Diethylaniline	ND	ug/Sample	--	
Dimethylaniline	ND	ug/Sample	--	
3,3'-Dimethoxybenzidine	ND	ug/Sample	--	
Hydroquinone	ND	ug/Sample	--	
4,4'-Methyl-bis(2-chloroaniline)	ND	ug/Sample	--	
4-Nitrodiphenyl	ND	ug/Sample	--	
Trifluralin	ND	ug/Sample	--	

Surrogate	Recovery	Acceptable Range	
Nitrobenzene-d5	ND %	45 - 107	H
2-Fluorobiphenyl	ND %	62 - 110	
Terphenyl-d14	ND %	58 - 135	
Phenol-d5	ND %	43 - 130	
2-Fluorophenol	ND %	36 - 111	
2,4,6-Tribromophenol	ND %	58 - 131	

Note H = Spiked analyte not detected because of required sample dilution.
ND = Not Detected

736

Reported By: Emily Uebelhoer

Approved By: Mike Orbanosky

The cover letter is an integral part of this report.
Rev 230787

Semivolatiles Library Search (20 Compound TID)
Method 8270

Client Name: Pacific Environmental Services

Client ID: S-MM5-1B-F, FH, XAD, COND, BH

LAB ID: 300681-0002-SA

Matrix: AIRTRAIN

Authorized: 30 JUL 98

Sampled: 25 JUL 98

Prepared: NA

Received: 30 JUL 98

Analyzed: 31 AUG 98

Dilution Factor: 100

Parameter	Result	Units	Reporting Limit	Qualifier
Unknown alkane	7800	ug/Sample	--	
Dodecane	7800	ug/Sample	--	0
Unknown	15000	ug/Sample	--	
Unknown hydrocarbon	17000	ug/Sample	--	
3-Hexadecene, (z)-	10000	ug/Sample	--	0
Undecane	11000	ug/Sample	--	0
Unknown alkane	20000	ug/Sample	--	
Unknown alkane	26000	ug/Sample	--	
Unknown	14000	ug/Sample	--	
Heptadecane, 2,6,10,14 -tetramethyl-	22000	ug/Sample	--	0
Dodecane, 3-methyl-	22000	ug/Sample	--	0
Unknown alkane	9900	ug/Sample	--	
Nonadecane	26000	ug/Sample	--	0
Nonadecane	18000	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	40000	ug/Sample	--	0
Unknown alkane	12000	ug/Sample	--	
Nonadecane	17000	ug/Sample	--	0
Nonadecane	8100	ug/Sample	--	0
Nonadecane	9800	ug/Sample	--	0
Nonadecane	7800	ug/Sample	--	0

Note 0 = Or structurally similar compound (isomer).
NA = Not Applicable

727

Reported By: Emily Uebelhoer

Approved By: Mike Orbanosky

The cover letter is an integral part of this report.

Rev 230787

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068102
Std Id: ST16980831

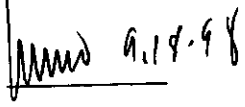
Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 19:31
Analyst: DAT Instrument ID: F16 Run Factor: 300.
Quan List Threshold: 0.95 Surrogate Vol.: 100.

Surrogate Spike Recoveries
8270-G, LIMS 10/10/96

Surrogate	Surrogate Amount (ug)		% Recovery		QC limits
	Spiked	Measured	Measured		
CS20 NITROBENZENE-D5	50.0	BDL	0*	45	107
CS25 2-FLUOROBIPHENY	50.0	BDL	0*	62	110
CS30 TERPHENYL-D14	50.0	BDL	0*	58	135
CS45 PHENOL-D5	100.	BDL	0*	43	130
CS50 2-FLUOROPHENOL	100.0	140.0	140.	*	36 111
CS55 2,4,6-TRIBROMOP	100.	BDL	0*	58	131

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting
				Limit
HEXACHLOROBENZENE-C13		ND	UG/A	3000.0
C310 N-NITROSODIMETHYLAMINE		ND	UG/A	3000.0
PYRIDINE		ND	UG/A	6000.0
2-PICOLINE		ND	UG/A	3000.0
N-NITROSOMETHYLETHYLAMINE		ND	UG/A	3000.0
METHYLMETHANESULFONATE		ND	UG/A	3000.0
N-NITROSODIETHYLAMINE		ND	UG/A	3000.0
ETHYLMETHANE SULFONATE		ND	UG/A	3000.0
PENTACHLOROETHANE		ND	UG/A	3000.0
C320 ANILINE		ND	UG/A	3000.0
C315 PHENOL		ND	UG/A	3000.0
C325 BIS(2-CHLOROETHYL)ETHER		ND	UG/A	3000.0
C330 2-CHLOROPHENOL		ND	UG/A	3000.0
C335 1,3-DICHLOROBENZENE		ND	UG/A	3000.0
C340 1,4-DICHLOROBENZENE		ND	UG/A	3000.0
C345 BENZYL ALCOHOL		ND	UG/A	3000.0
C350 1,2-DICHLOROBENZENE		ND	UG/A	3000.0
C355 2-METHYLPHENOL		ND	UG/A	3000.0
C360 2,2'-OXYBIS(1-CLPROPAN)		ND	UG/A	3000.0
C361 ACETOPHENONE		ND	UG/A	3000.0
N-NITROSOPYRROLIDINE		ND	UG/A	3000.0
N-NITROSOMORPHOLINE		ND	UG/A	3000.0
3-METHYL PHENOL		ND	UG/A	3000.0
C365 4-METHYLPHENOL		ND	UG/A	3000.0
C370 N-NITROSO-DI-N-PROPYLAM		ND	UG/A	3000.0
O-TOLUIDINE		ND	UG/A	3000.0
C375 HEXACHLOROETHANE		ND	UG/A	3000.0
C410 NITROBENZENE		ND	UG/A	3000.0
N-NITROSOPIPERIDINE		ND	UG/A	3000.0

Reviewed by:  9.14.98

733

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068102
Std Id: ST16980831

Sample: S-MM5-18 1/3SA/100M INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 19:31
Analyst: DAT Instrument ID: F16 Run Factor: 300.
Quan List Threshold: 0.95 Surrogate Vol.: 100.

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C415 ISOPHORONE		ND	UG/A	3000.0
C420 2-NITROPHENOL		ND	UG/A	3000.0
C425 2,4-DIMETHYLPHENOL		ND	UG/A	3000.0
C435 BIS(2-CHLOROETHOXY)METH		ND	UG/A	3000.0
C440 2,4-DICHLOROPHENOL		ND	UG/A	3000.0
C430 BENZOIC ACID		ND	UG/A	15000.0
AA-DIMETHYLPHENETHYLAMINE		ND	UG/A	3000.0
C445 1,2,4-TRICHLOROBENZENE		ND	UG/A	3000.0
C450 NAPHTHALENE		ND	UG/A	3000.0
C455 4-CHLOROANALINE		ND	UG/A	3000.0
2,6-DICHLOROPHENOL		ND	UG/A	3000.0
HEXACHLOROPROPENE		ND	UG/A	3000.0
C460 HEXACHLOROBUTADIENE		ND	UG/A	3000.0
P-PHENYLENE DIAMINE		ND	UG/A	3000.0
N-NITROSODI-N-BUTYLAMINE		ND	UG/A	3000.0
C465 4-CHLORO-3-METHYLPHENO		ND	UG/A	3000.0
SAFROLE		ND	UG/A	3000.0
C470 2-METHYLNAPHTHALENE	652	4700.	UG/A	3000.0
1,2,4,5-TETRACHLOROBENZENE		ND	UG/A	3000.0
ISOSAFROLE (#1)		ND	UG/A	6000.0
C510 HEXACHLOROCYCLOPENTADI		ND	UG/A	3000.0
C515 2,4,6-TRICHLOROPHENOL		ND	UG/A	3000.0
C520 2,4,5-TRICHLOROPHENOL		ND	UG/A	3000.0
ISOSAFROLE (#2)		ND	UG/A	6000.0
C525 2-CHLORONAPHTHALENE		ND	UG/A	3000.0
1-CHLORONAPHTHALENE		ND	UG/A	1500.0
C530 2-NITROANALINE		ND	UG/A	15000.0
1,4-NAPHTHOQUINONE		ND	UG/A	3000.0
C535 DIMETHYLPHTHALATE		ND	UG/A	3000.0
1,3-DINITROBENZENE		ND	UG/A	3000.0
C540 ACENAPHTHYLENE		ND	UG/A	3000.0
C543 2,6-DINITROTOLUENE		ND	UG/A	3000.0
C545 3-NITROANILINE		ND	UG/A	15000.0
C550 ACENAPHTHENE		ND	UG/A	3000.0
C555 2,4-DINITROPHENOL		ND	UG/A	15000.0
C565 DIBENZOFURAN		ND	UG/A	3000.0
C560 4-NITROPHENOL		ND	UG/A	15000.0
PENTACHLOROBENZENE		ND	UG/A	3000.0
C570 2,4-DINITROTOLUENE		ND	UG/A	3000.0
1-NAPHTHYLAMINE		ND	UG/A	3000.0
2-NAPHTHYLAMINE		ND	UG/A	3000.0
2,3,4,6-TETRACHLOROPHENOL		ND	UG/A	6000.0
C580 DIETHYLPHTHALATE		ND	UG/A	3000.0
C590 FLUORENE		ND	UG/A	3000.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068102
Std Id: ST16980831

Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 19:31
Analyst: DAT Instrument ID: F16 Run Factor: 300.
Quan List Threshold: 0.95 Surrogate Vol.: 100.

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C585 4-CHLOROPHENYL-PHENYLE		ND	UG/A	3000.0
5-NITRO-O-TOLUIDINE		ND	UG/A	3000.0
C595 4-NITROANALINE		ND	UG/A	15000.0
C610 4,6-DINITRO-2-METHYLPH		ND	UG/A	15000.0
C615 N-NITROSODIPHENYLAMINE		ND	UG/A	3000.0
C620 AZOBENZENE		ND	UG/A	3000.0
SYM-TRINITROBENZENE		ND	UG/A	3000.0
C625 4-BROMOPHENYL-PHENYLET		ND	UG/A	3000.0
PHENACETIN		ND	UG/A	3000.0
DIALATE	234	ND	UG/A	3000.0
C630 HEXACHLOROBENZENE		ND	UG/A	3000.0
4-AMINOBIIPHENYL		ND	UG/A	15000.0
C635 PENTACHLOROPHENOL		ND	UG/A	3000.0
PRONAMIDE		ND	UG/A	15000.0
PENTACHLORONITROBENZENE		ND	UG/A	3000.0
C640 PHENANTHRENE		ND	UG/A	3000.0
C645 ANTHRACENE		ND	UG/A	3000.0
2SECBUTYL-4,6-DINITROPHENOL		ND	UG/A	3000.0
C647 CARBAZOLE		ND	UG/A	3000.0
C650 DI-N-BUTYLPHTHALATE		ND	UG/A	3000.0
4-NITROQUINOLINE-1-OXIDE		ND	UG/A	3000.0
METHAPYRILENE		ND	UG/A	3000.0
ISODRIN		ND	UG/A	3000.0
C655 FLUORANTHENE		ND	UG/A	3000.0
CHLOROBENZILATE		ND	UG/A	30000.0
C710 BENZIDINE		ND	UG/A	3000.0
C715 PYRENE		ND	UG/A	3000.0
ARAMITE (#1)		ND	UG/A	3000.0
ARAMITE (#2)		ND	UG/A	3000.0
P-DIMETHYLAMINOAZOBENZENE		ND	UG/A	3000.0
3,3'-DIMETHYLBENZIDINE		ND	UG/A	15000.0
KEPONE		ND	UG/A	3000.0
C720 BUTYLBENZYLPHTHALATE		ND	UG/A	3000.0
2-ACETYLAMINOFLUORENE		ND	UG/A	3000.0
C730 BENZO(A)ANTHRACENE		ND	UG/A	6000.0
C725 3,3'-DICHLOROBENZIDINE		ND	UG/A	3000.0
C740 CHRYSENE		ND	UG/A	3000.0
C745 BIS(2-ETHYLHEXYL)PHTHA		ND	UG/A	3000.0
3-METHYLCHOLANTHRENE		ND	UG/A	3000.0
C760 DI-N-OCTYL PHTHALATE		ND	UG/A	3000.0
C765 BENZO(B)FLUORANTHENE		ND	UG/A	3000.0
7,12-DIMETHYLBENZANTHRACENE		ND	UG/A	3000.0
C770 BENZO(K)FLUORANTHENE		ND	UG/A	3000.0
HEXACHLOROPHENE		ND	UG/A	3000.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

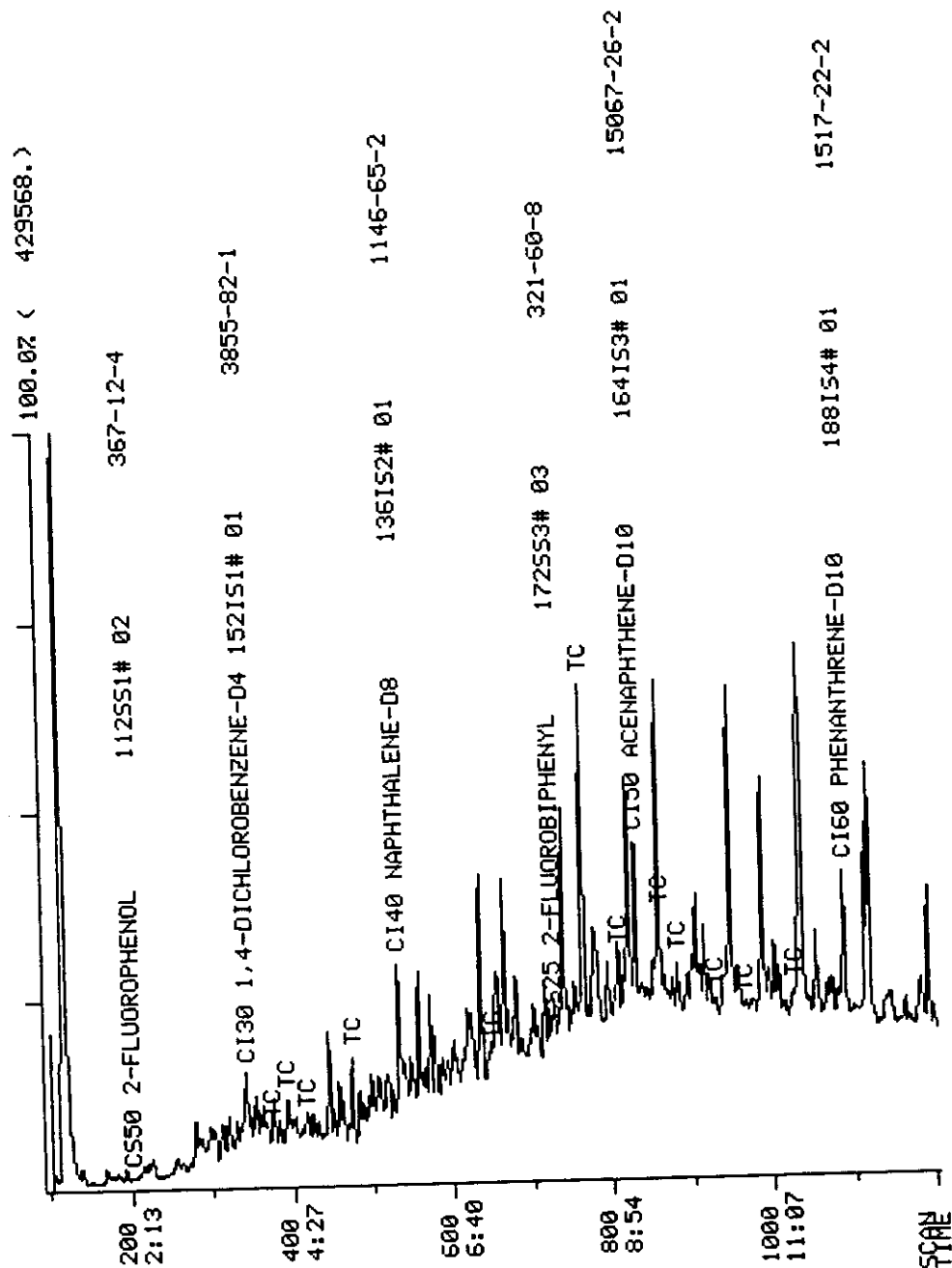
Data File: 30068102
Std Id: ST16980831

Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 19:31
Analyst: DAT Instrument ID: F16 Run Factor: 300.
Quan List Threshold: 0.95 Surrogate Vol.: 100.

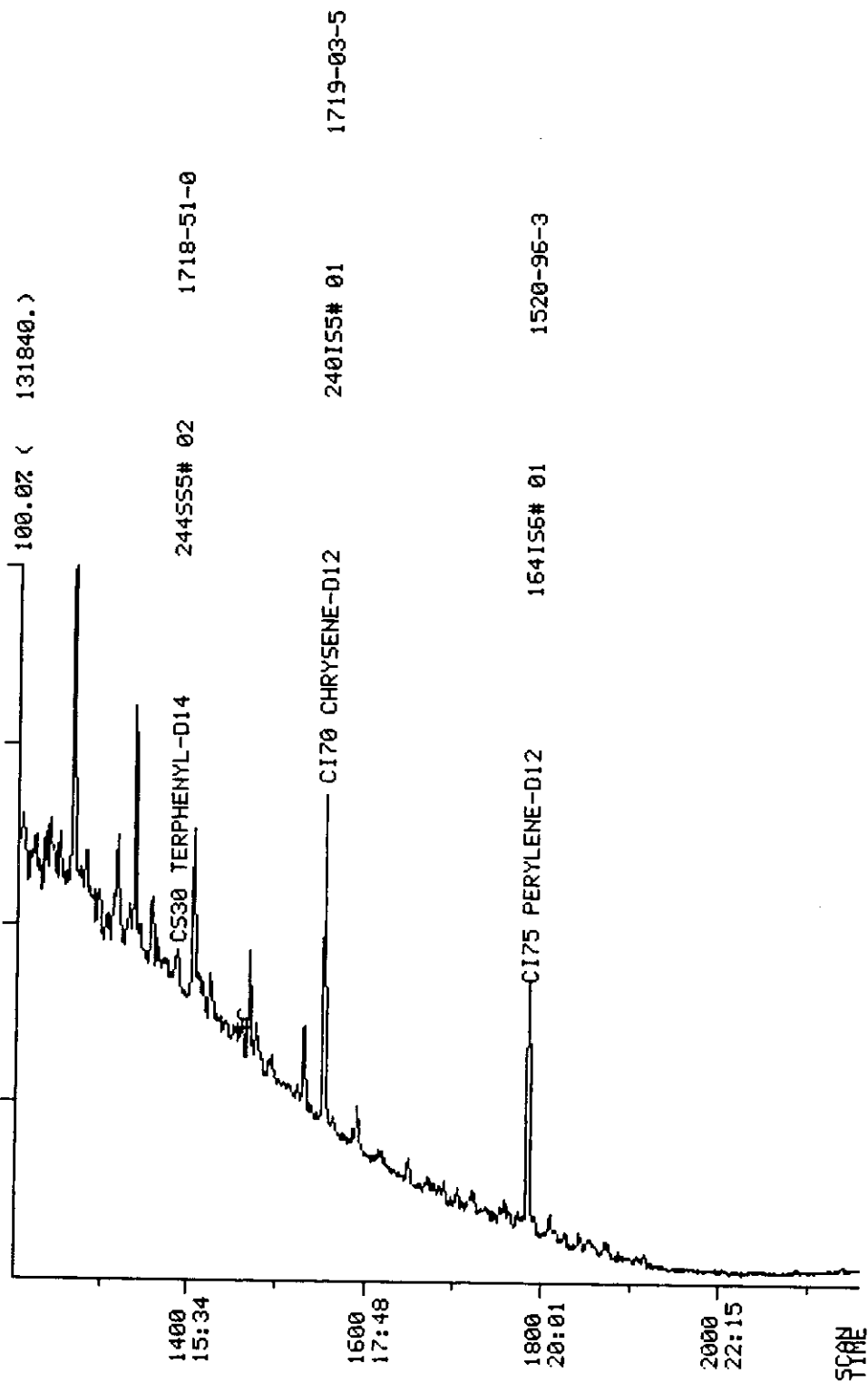
Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C775 BENZO(A)PYRENE		ND	UG/A	3000.0
C780 INDENO(1,2,3-CD)PYRENE		ND	UG/A	3000.0
C785 DIBENZ(A,H)ANTHRACENE		ND	UG/A	3000.0
C790 BENZO(G,H,I)PERYLENE		ND	UG/A	3000.0

DATA FROM FILE: 30068102 SCANS 93 TO 1203 ACQUIRED: 08/31/98 19:31:00
 CALI: 30068102 #3
 SAMPLE: 5-MMS-1B 1/35A/100M INST. ID: F16
 CONDS.: UG/ML *100Z/100Z *(NA/NA >1/35A NA M



DATA FROM FILE: 30068102 SCANS 1283 TO 2158 ACQUIRED: 08/31/98 19:31:00
 SAMPLE: S-MMS-1B 1/35A/100M CALI: 30068102 #3
 CONDS.: UG/ML *100ML *100Z *(NA/NA)/1/35A NA M INST. ID: F16



STANDARD USED FOR THIS REPORT IS ST16980831 REPORT PRINTED: 9/01/98 13:16:54

ADDITIONAL STANDARD USED FOR THIS REPORT IS: ST16980831A

QUANTERRA QUANTITATION SUMMARY

File: 30068102

Sample: S-MM5-18 1/3SA/100M INST. ID: F16
Analyst: DAT Instrument Id: F16 Analyzed: 08/31/98 19:31

Compounds with amounts less than 0.20 reported as NOT FOUND

Library							Units: UG/ML		
No	Entry	Name	Mass	Meth	Scan	Ref	Area	RRF(L)	Amount
1	S1#	1 C130 1,4-DICHLOROBENZENE	152	A	BV	340	1	26359. 1.000	40.000
2	S2#	1 C140 NAPHTHALENE-D8	136	A	BB	530	2	96109. 1.000	40.000
3	S3#	1 C150 ACENAPHTHENE-D10	164	A	BB	829	3	54125. 1.000	40.000
4	S4#	1 C160 PHENANTHRENE-D10	188	A	BB	1088	4	95829. 1.000	40.000
5	S5#	1 C170 CHRYSENE-D12	240	A	BB	1552	5	57388. 1.000	40.000
6	S6#	1 C175 PERYLENE-D12	264	A	BB	1782	6	44375. 1.000	40.000
7	S2#	2 CS20 NITROBENZENE-D5	82	A	BB	425	2	385. 0.470	0.341
8	S3#	3 CS25 2-FLUOROBIPHENYL	172	A	BB	724	3	1292. 1.234	0.774
9	S5#	2 CS30 TERPHENYL-D14	244	A	BB	1384	5	1156. 0.995	0.810
10	S1#	3 CS45 PHENOL-D5	99			1	NOT FOUND		
11	S1#	2 CS50 2-FLUOROPHENOL	112	A	BV	198	1	1262. 1.368	1.400
12	S3#	2 CS55 2,4,6-TRIBROMOPHENO	330	A	BB	971	3	242. 0.236	0.759
13	S4#	4 HEXACHLOROBENZENE-C13	294			4	NOT FOUND		
14	S1#	15 C310 N-NITROSODIMETHYLAM	74			1	NOT FOUND		
15	S1#	20 PYRIDINE	79			1	NOT FOUND		
16	S1#	30 2-PICOLINE	93			1	NOT FOUND		
17	S1#	40 N-NITROSOMETHYLETHYLAMIN	42			1	NOT FOUND		
18	S1#	60 METHYLMETHANESULFONATE	80			1	NOT FOUND		
19	S1#	70 N-NITROSODIETHYLAMINE	102			1	NOT FOUND		
20	S1#	85 ETHYLMETHANE SULFONATE	79			1	NOT FOUND		
21	S1#	95 PENTACHLOROETHANE	117	A	BB	306	1	210. 0.602	0.530
22	S1#	100 C320 ANILINE	93			1	NOT FOUND		
23	S1#	105 C315 PHENOL	94			1	NOT FOUND		
24	S1#	110 C325 BIS(2-CHLOROETHYL)E	93			1	NOT FOUND		
25	S1#	115 C330 2-CHLOROPHENOL	128			1	NOT FOUND		
26	S1#	125 C335 1,3-DICHLOROBENZENE	146			1	NOT FOUND		
27	S1#	130 C340 1,4-DICHLOROBENZENE	146			1	NOT FOUND		
28	S1#	145 C345 BENZYL ALCOHOL	108	A	VV	372	1	546. 0.854	0.970
29	S1#	150 C350 1,2-DICHLOROBENZENE	146			1	NOT FOUND		
30	S1#	160 C355 2-METHYLPHENOL	108	A	BB	392	1	1048. 1.175	1.354
31	S1#	165 C360 2,2'-OXYBIS(1-CLPRO	45			1	NOT FOUND		
32	S1#	170 C361 ACETOPHENONE	105			1	NOT FOUND		
33	S1#	175 N-NITROSOPYRROLIDINE	100			1	NOT FOUND		
34	S1#	180 N-NITROSOMORPHOLINE	56			1	NOT FOUND		
35	S1#	182 3-METHYL PHENOL	108			1	NOT FOUND		
36	S1#	185 C365 4-METHYLPHENOL	108	A	BB	418	1	1628. 1.276	1.936
37	S1#	190 C370 N-NITROSO-DI-N-PROP	70	A	BB	414	1	618. 0.906	1.035
38	S1#	195 O-TOLUIDINE	106			1	NOT FOUND		
39	S1#	200 C375 HEXACHLOROETHANE	117			1	NOT FOUND		
40	S2#	10 C410 NITROBENZENE	77	A	BB	422	2	1481. 0.445	1.385
41	S2#	15 N-NITROSOPIPERIDINE	42			2	NOT FOUND		
42	S2#	20 C415 ISOPHORONE	82	A	BB	473	2	4288. 0.741	2.407
43	S2#	25 C420 2-NITROPHENOL	139			2	NOT FOUND		
44	S2#	30 C425 2,4-DIMETHYLPHENOL	107	A	BB	500	2	621. 0.330	0.784
45	S2#	35 C435 BIS(2-CHLOROETHOXY)	93			2	NOT FOUND		

752

46	S2# 40	C440 2,4-DICHLOROPHENOL	162 A 88	520 2	306. 0.308	0.414
47	S2# 45	C430 BENZOIC ACID	122	2	NOT FOUND	
48	S2# 50	AA-DIMETHYLPHENETHYLAMIN	58	2	NOT FOUND	
49	S2# 55	C445 1,2,4-TRICHLOROBENZ	180	2	NOT FOUND	

50 S2# 60 C450 NAPHTHALENE	128 A BV	533 2	17654. 1.022	7.188
51 S2# 80 C455 4-CHLOROANALINE	127 A BB	556 2	496. 0.454	0.455
52 S2# 85 2,6-DICHLOROPHENOL	162	2	NOT FOUND	
53 S2# 90 HEXACHLOROPROPENE	213	2	NOT FOUND	
54 S2# 95 C460 HEXACHLOROBUTADIENE	225	2	NOT FOUND	
55 S2#115 P-PHENYLENE DIAMINE	108	2	NOT FOUND	
56 S2#120 N-NITROSODI-N-BUTYLAMINE	84	2	NOT FOUND	
57 S2#130 C465 4-CHLORO-3-METHYLPH	107	2	NOT FOUND	
58 S2#140 SAFROLE	162 A BB	648 2	1275. 0.248	2.142
59 S2#145 C470 2-METHYLNAPHTHALENE	142 A BB	652 2	28264. 0.751	15.658
60 S3# 10 1,2,4,5-TETRACHLOROBENZE	214	3	NOT FOUND	
61 S3# 15 ISOSAFROLE (#1)	162	3	NOT FOUND	
62 S3# 20 C510 HEXACHLOROCYCLOPENT	237	3	NOT FOUND	
63 S3# 25 C515 2,4,6-TRICHLOROPHEN	196	3	NOT FOUND	
64 S3# 30 C520 2,4,5-TRICHLOROPHEN	196	3	NOT FOUND	
65 S3# 35 ISOSAFROLE (#2)	104 A BB	738 3	390. 0.195	1.482
66 S3# 40 C525 2-CHLORONAPHTHALENE	162	3	NOT FOUND	
67 S3# 42 1-CHLORONAPHTHALENE	162	3	NOT FOUND	
68 S3# 45 C530 2-NITROANALINE	65 A BB	764 3	655. 0.502	0.965
69 S3# 50 1,4-NAPHTHOQUINONE	158	3	NOT FOUND	
70 S3# 55 C535 DIMETHYLPHTHALATE	163	3	NOT FOUND	
71 S3# 60 1,3-DINITROBENZENE	168 A BB	807 3	258. 0.197	0.966
72 S3# 65 C540 ACENAPHTHYLENE	152 A BB	808 3	622. 1.838	0.250
73 S3# 70 C543 2,6-DINITROTOLUENE	165 A BB	817 3	391. 0.324	0.892
74 S3# 75 C545 3-NITROANILINE	138	3	NOT FOUND	
75 S3# 80 C550 ACENAPHTHENE	153 A BV	835 3	1132. 1.159	0.722
76 S3# 85 C555 2,4-DINITROPHENOL	184 A BV	860 3	308. 0.189	1.202
77 S3# 90 C565 DIBENZOFURAN	168 A BB	867 3	1063. 1.669	0.471
78 S3# 95 C560 4-NITROPHENOL	109	3	NOT FOUND	
79 S3#100 PENTACHLOROBENZENE	250	3	NOT FOUND	
80 S3#105 C570 2,4-DINITROTOLUENE	165 A BB	881 3	551. 0.397	1.025
81 S3#110 1-NAPHTHYLAMINE	143 A BB	882 3	280. 0.837	0.247
82 S3#115 2-NAPHTHYLAMINE	143	3	NOT FOUND	
83 S3#120 2,3,4,6-TETRACHLOROPHENO	232	3	NOT FOUND	
84 S3#130 C580 DIETHYLPHTHALATE	149	3	NOT FOUND	
85 S3#135 C590 FLUORENE	166 A BB	926 3	3536. 1.298	2.014
86 S3#140 C585 4-CHLOROPHENYL-PHEN	204	3	NOT FOUND	
87 S3#145 5-NITRO-O-TOLUIDINE	152 A VB	942 3	1876. 0.356	3.899
88 S3#150 C595 4-NITROANALINE	138	3	NOT FOUND	
89 S4# 10 C610 4,6-DINITRO-2-METHY	198	4	NOT FOUND	
90 S4# 15 C615 N-NITROSODIPHENYLAM	169 A BB	964 4	3648. 0.551	2.765
91 S4# 20 C620 AZOBENZENE	77	4	NOT FOUND	
92 S4# 25 SYM-TRINITROBENZENE	75 A BV	1025 4	464. 0.162	1.198
93 S4# 30 C625 4-BROMOPHENYL-PHENY	248	4	NOT FOUND	
94 S4# 35 PHENACETIN	108	4	NOT FOUND	
95 S4# 37 DIALLATE	234	4	NOT FOUND	
96 S4# 40 C630 HEXACHLOROBENZENE	284	4	NOT FOUND	
97 S4# 45 4-AMINOBIIPHENYL	169	4	NOT FOUND	
98 S4# 50 C635 PENTACHLOROPHENOL	266	4	NOT FOUND	
99 S4# 55 PRONAMIDE	173	4	NOT FOUND	
100 S4# 60 PENTACHLORONITROBENZENE	237 A BB	1066 4	106. 0.080	0.551
101 S4# 65 C640 PHENANTHRENE	178 A BB	1091 4	11034. 1.033	4.460
102 S4# 70 C645 ANTHRACENE	178	4	NOT FOUND	
103 S4# 75 ZSECBUTYL-4,6-DINITROPHE	211	4	NOT FOUND	
104 S4# 80 C647 CARBAZOLE	167 A BB	1149 4	538. 0.764	0.294
105 S4# 85 C650 DI-N-BUTYLPHTHALATE	149	4	NOT FOUND	
106 S4#100 4-NITROQUINOLINE-1-OXIDE	190	4	NOT FOUND	
107 S4#105 METHAPYRILENE	58	4	NOT FOUND	
108 S4#106 ISODRIN	193 A BV	1267 4	201. 0.127	0.662
109 S4#110 C655 FLUORANTHENE	202	4	NOT FOUND	
110 S4#120 CHLOROBENZILATE	139	4	NOT FOUND	

111 S5# 10 C710 BENZIDINE	184	5	NOT FOUND		
112 S5# 15 C715 PYRENE	202 A BB 1335	5	1044.	1.323	0.550
113 S5# 20 ARAMITE (#1)	185	5	NOT FOUND		
114 S5# 25 ARAMITE (#2)	185	5	NOT FOUND		
115 S5# 30 P-DIMETHYLAMINOAZOBENZEN	120	5	NOT FOUND		
116 S5# 35 3,3'-DIMETHYLBENZIDINE	212	5	NOT FOUND		
117 S5# 37 KEPONE	272 A BB 1461	5	278.	0.082	2.364
118 S5# 40 C720 BUTYLBENZYLPHTHALAT	149	5	NOT FOUND		
119 S5# 45 2-ACETYLAMINOFLUORENE	181	5	NOT FOUND		
120 S5# 50 C730 BENZO(A)ANTHRACENE	228	5	NOT FOUND		
121 S5# 55 C725 3,3'-DICHLOROBENZID	252	5	NOT FOUND		
122 S5# 60 C740 CHRYSENE	228 A BB 1555	5	1337.	1.014	0.919
123 S5# 65 C745 BIS(2-ETHYLHEXYL)PH	149	5	NOT FOUND		
124 S5# 85 3-METHYLCHOLANTHRENE	268	5	NOT FOUND		
125 S6# 10 C760 DI-N-OCTYL PHTHALAT	149	6	NOT FOUND		
126 S6# 15 C765 BENZO(B)FLUORANTHEN	252	6	NOT FOUND		
127 S6# 20 7,12-DIMETHYLBENZANTHRAC	256	6	NOT FOUND		
128 S6# 25 C770 BENZO(K)FLUORANTHEN	252	6	NOT FOUND		
129 S6# 30 HEXACHLOROPHENE	196	6	NOT FOUND		
130 S6# 35 C775 BENZO(A)PYRENE	252	6	NOT FOUND		
131 S6# 55 C780 INDENO(1,2,3-CD)PYR	276	6	NOT FOUND		
132 S6# 60 C785 DIBENZ(A,H)ANTHRACE	278	6	NOT FOUND		
133 S6# 65 C790 BENZO(G,H,I)PERYLEN	276	6	NOT FOUND		

TARGET COMPOUND COMPARISON

COMPOUND: C470 2-METHYLNAPHTHALENE 142 S2#145

91-57-6

BASE M/Z: 142 RIC: 114048.

RAW DATA: 30068102 #652

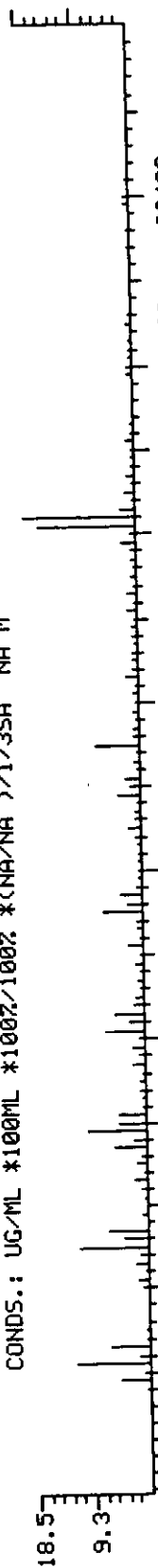
08/31/98 19:31

SAMPLE: S-MM5-1B 1/35A/100M

INST. ID: F16

CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA) >1/35A NA M

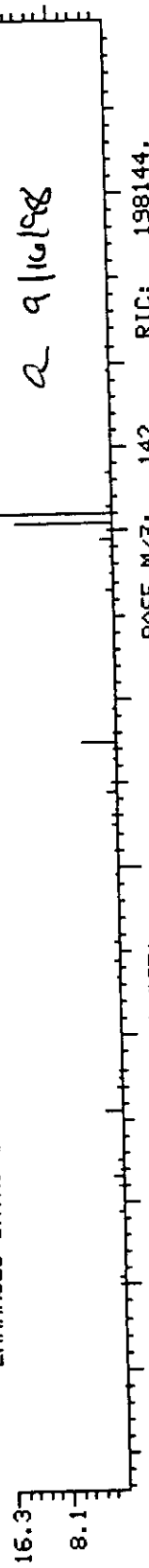
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BASE M/Z: 142 RIC: 29408.

8112.

ENHANCED DATA: 30068102 #652

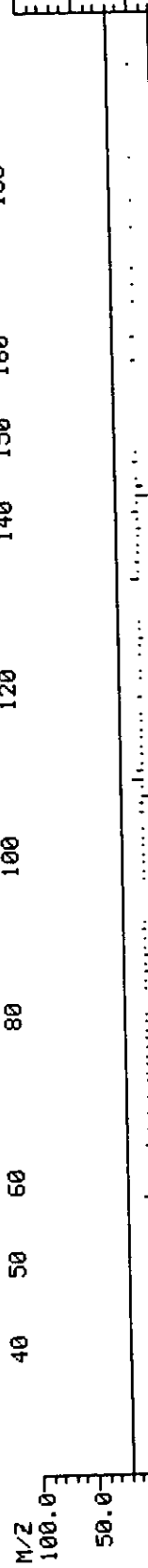
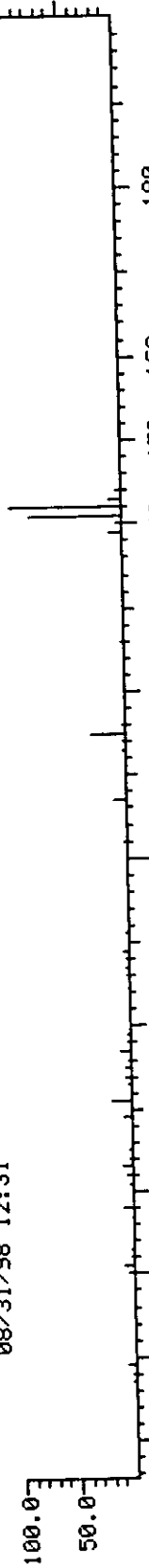


BASE M/Z: 142 RIC: 198144.

49792.

STANDARD FILE: ST16980831 #651

08/31/98 12:31



QUANTERRA Internal Standard Check

Standard Filename: ST16980831

Analyzed: 08/31/98 12:31

Standard	IS# 1			IS# 2			IS# 3		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	28649.		3.78	100432.		5.90	54347.		9.21
Upper Limit	57297.=+200		4.28	200864.=+200		6.40	108695.=+200		9.71
Lower Limit	14324.=+ 50		3.28	50216.=+ 50		5.40	27174.=+ 50		8.71
Filename									
1 30068102	26359.= 92		3.78	96109.= 95		5.90	54125.= 99		9.22

IS# 1 = C130 1,4-DICHLOROBENZENE-D4

IS# 2 = C140 NAPHTHALENE-D8

IS# 3 = C150 ACENAPHTHENE-D10

* - indicates an I.S. is outside QC limit(s)

Standard	IS# 4			IS# 5			IS# 6		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	98520.		12.08	62289.		17.26	50750.		19.83
Upper Limit	197039.=+200		12.58	124578.=+200		17.76	101500.=+200		20.33
Lower Limit	49260.=+ 50		11.58	31145.=+ 50		16.76	25375.=+ 50		19.33
Filename									
1 30068102	95829.= 97		12.10	57388.= 92		17.26	44375.= 87		19.82

IS# 4 = C160 PHENANTHRENE-D10

IS# 5 = C170 CHRYSENE-D12

IS# 6 = C175 PERYLENE-D12

* - indicates an I.S. is outside QC limit(s)

Data Reduced by: *FW* Date: *9/8/98*
Data Reviewed by: *FW* Date: *9/11/98*

Data File: 30068102

QUANTERRA GC/MS TIC REPORT (Part 1)

Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
Analyst: DAT Date Analyzed: 08/31/98 19:31
Run Factor: 300.

# SCAN	Concentration in Sample (UG/A)	CAS #
1 444	7800.	62108-23-0
DECANE, 2,5,6-TRIMETHYL-		

57 46400
alkane

X 473	7100.	74962-98-4
2-TRIDECEN-1-OL, (E)-		

unknown 5700

3 556	7800.	112-40-3
DODECANE <i>or isomer</i>		

X 570	6700.	17301-23-4
UNDECANE, 2,6-DIMETHYL-		

5 617	15000.	00-00-0
UNKNOWN		

5200

6 633	17000.	50871-03-9
1-DECENE, 3,4-DIMETHYL-		

57904300
unkn. hydrocarbon

7 654	10000.	34303-81-6
3-HEXADECENE, (Z)- <i>or isomer</i>		

8 663	11000.	1120-21-4
UNDECANE		

X 678	7400.	00-00-0
UNKNOWN		

5700

10 738	20000.	54105-67-8
HEPTADECANE, 2,6-DIMETHYL-		

5746400
alkane

750

11 764 26000. ~~62108-21-8~~ alkane 5746400
~~DECANE, 6-ETHYL-2-METHYL-~~

12 777 14000. 00-00-0 5700
UNKNOWN

13 820 22000. 18344-37-1 or isomer
HEPTADECANE, 2,6,10,14-TETRAMETHYL-

14 858 22000. 17312-57-1
DODECANE, 3-METHYL-

15 904 9900. ~~52785-43-4~~ alkane 5746400
~~NONADECANOL~~

16 948 26000. 629-92-5 or isomer
NONADECANE

17 988 18000. 629-92-5
NONADECANE

18 1036 40000. 54105-67-8
HEPTADECANE, 2,6-DIMETHYL-

X 1054 7000. 7206-19-1
3-OCTADECENE, (E)-

20 1114 12000. ~~29812-79-1~~ alkane 5646400
~~HYDROXYLAMINE, O-DECYL-~~

21 1119 17000. 629-92-5
NONADECANE or isomer

22 1191 8100. 629-92-5
NONADECANE

~~23~~ 1265 6900. ~~54105-67-8~~
~~HEPTADECANE, 2,6-DIMETHYL~~ alkane

24 1335 9800.
~~2-METHYLOCTADECANE~~ or is nonadecane or isomer
C 9/18/98

25 1402 7800. 629-92-5
NONADECANE



DATA FILE: 30068102

QUANTERRA GC/MS TIC REPORT (Part 2)

CONCENTRATION = AREA(TIC)*CONC(1S)/AREA(1S)

			INT.		AMOUNT					
#	FIT	PURITY	STD.	RT	RRT	AREA	HEIGHT	(UG/ML)	LIB	LIB #
1	941	658	2	4:56	0.536	197944.	59360.	26.115	NB	19015.
2	874	669	2	5:15	0.571	180512.	46417.	23.815	NB	22482.
3	943	842	2	6:11	0.671	196010.	69931.	25.860	NB	15343.
4	972	853	2	6:20	0.688	168832.	54235.	22.274	NB	19054.
5	901	595	2	6:51	0.744	367872.	39360.	48.534	UK	1.
6	948	721	2	7:02	0.764	429472.	115616.	56.661	NB	14799.
7	930	700	2	7:16	0.789	262912.	45696.	34.687	NB	28768.
8	960	846	2	7:22	0.800	268288.	94208.	35.396	NB	11607.
9	944	401	2	7:32	0.818	186752.	43392.	24.639	UK	1.
10	950	826	2	8:12	0.890	495104.	127744.	65.320	NB	37462.
11	953	716	2	8:29	0.922	666007.	189300.	87.868	NB	19006.
12	911	347	2	8:38	0.937	363392.	50048.	47.943	UK	1.
13	972	826	2	9:07	0.989	552366.	122822.	72.875	NB	42196.
14	964	773	2	9:32	1.035	565839.	173725.	74.652	NB	19007.
15	938	661	2	10:03	1.090	250112.	56832.	32.998	NB	40233.
16	958	847	2	10:32	1.144	648704.	177408.	85.585	NB	37465.
17	943	830	3	10:59	0.908	500650.	128483.	58.597	NB	37465.
18	970	865	3	11:31	0.952	1137280.	201472.	133.109	NB	37462.
19	906	703	3	11:43	0.969	200704.	46336.	23.491	NB	34410.
20	957	799	3	12:23	1.024	336000.	104960.	39.326	NB	15969.
21	964	842	3	12:26	1.028	471552.	141056.	55.191	NB	37465.
22	971	792	3	13:14	1.095	231936.	77312.	27.146	NB	37465.
23	924	763	3	14:04	1.163	195456.	57749.	22.876	NB	37462.
24	933	740	4	14:50	0.860	167008.	42891.	32.701	NB	37456.
25	965	690	4	15:35	0.903	132416.	30848.	25.927	NB	37465.

Library Search Data: 30068102 # 444 Base m/z: 43
 08/31/98 19:31:00 + 4:56 Cali: 30068102 # 3 RIC: 49216.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 897 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 19015 DECANE, 2,5,6-TRIMETHYL-
 2 11607 UNDECANE
 3 15969 HYDROXYLAMINE, O-DECYL-
 4 19523 1-DECANOL, 2-ETHYL-
 5 11602 OCTANE, 2,4,6-TRIMETHYL-
 6 14793 1-UNDECENE, 4-METHYL-
 7 8104 OCTANE, 3,5-DIMETHYL-
 8 5151 HEPTANE, 3,4-DIMETHYL-
 9 12074 1-HEPTANOL, 2-PROPYL-

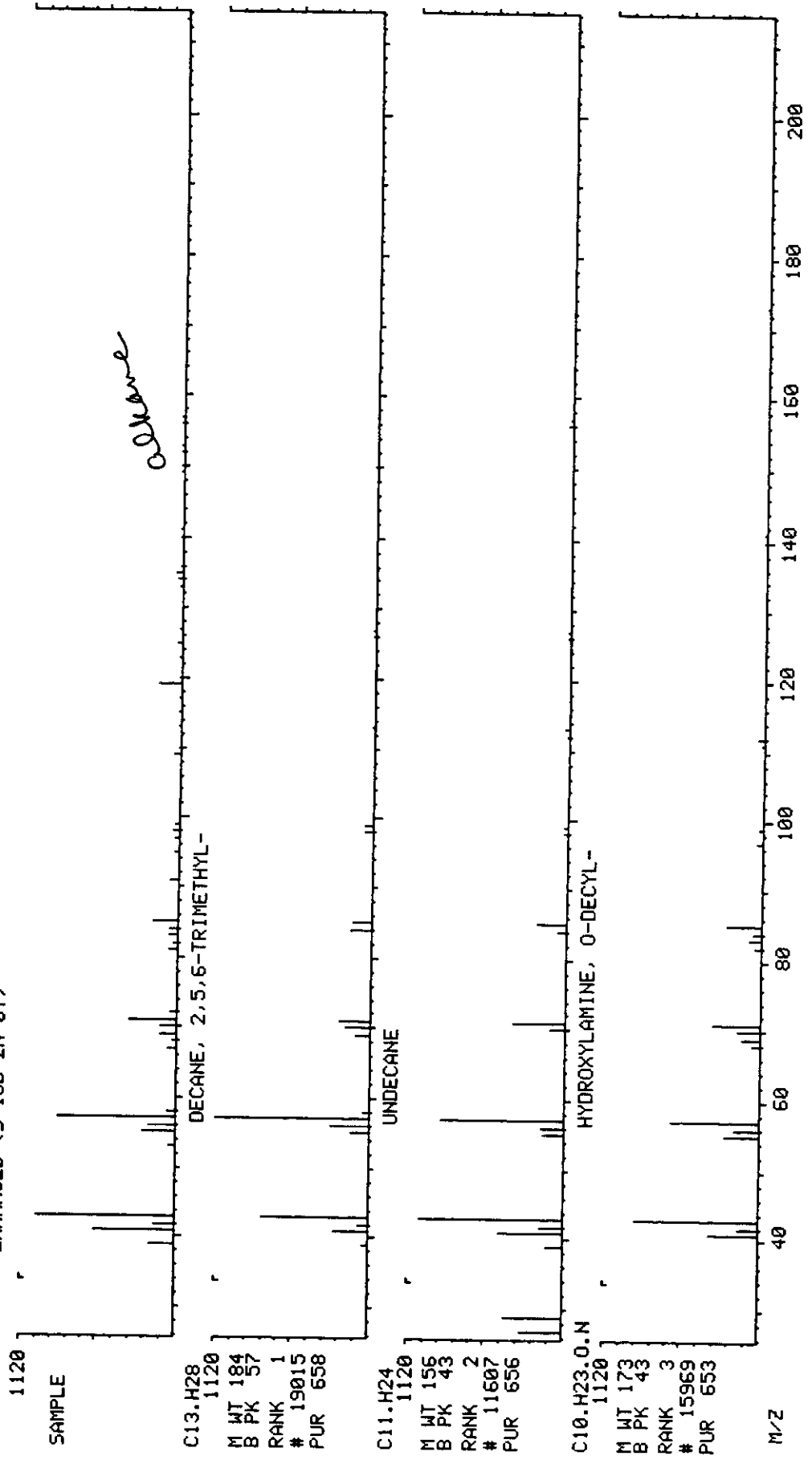
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H28	184	57	658	941	660
2	C11.H24	156	43	656	979	664
3	C10.H23.O.N	173	43	653	948	675
4	C12.H26.O	186	57	641	915	670
5	C11.H24	156	57	638	949	638
6	C12.H24	168	43	638	915	672
7	C10.H22	142	57	637	932	649
8	C9.H20	128	43	632	920	634
9	C10.H22.O	158	43	629	922	648

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	62108-23-0
2	—	—	—	—	1120-21-4
3	—	—	—	—	29812-79-1
4	—	—	—	—	21078-65-9
5	—	—	—	—	62016-37-9
6	—	—	—	—	74630-39-0
7	—	—	—	—	15869-93-9
8	—	—	—	—	922-28-1
9	—	—	—	—	10042-59-8

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 4:56
 SAMPLE: S-MMS-1B 1/35A/100M
 CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068102 # 444
 CALI: 30068102 # 3

BASE M/Z: 43
 RIC: 49215.



Library Search Data: 30068102 # 473 Base m/z: 43
 08/31/98 19:31:00 + 5:16 Cali: 30068102 # 3 RIC: 41024.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 459 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 22482 2-TRIDECEN-1-OL, (E)-
 2 25971 OXIRANE, DODECYL-
 3 37444 OXIRANE, HEXADECYL-
 4 34800 (R)-(-)-(Z)-14-METHYL-8-HEXADECEN-1-OL
 5 39230 9-EICOSYNE
 6 37449 OCTADECANAL
 7 29222 (Z)6-PENTADECEN-1-OL
 8 39224 1-EICOSYNE
 9 29226 PENTADECANAL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H26.O	198	57	669	874	675
2	C14.H28.O	212	41	660	933	660
3	C18.H36.O	268	57	656	845	679
4	C17.H34.O	254	55	651	852	661
5	C20.H38	278	81	649	902	666
6	C18.H36.O	268	43	642	843	685
7	C15.H30.O	226	67	624	867	642
8	C20.H38	278	82	622	841	648
9	C15.H30.O	226	82	606	890	606

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	74962-98-4
2	—	—	—	—	3234-28-4
3	—	—	—	—	7390-81-0
4	—	—	—	—	30689-78-2
5	—	—	—	—	71899-38-2
6	—	—	—	—	638-66-4
7	—	—	—	—	68797-95-5
8	—	—	—	—	765-27-5
9	—	—	—	—	2765-11-9

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 19:31:00 + 5:16

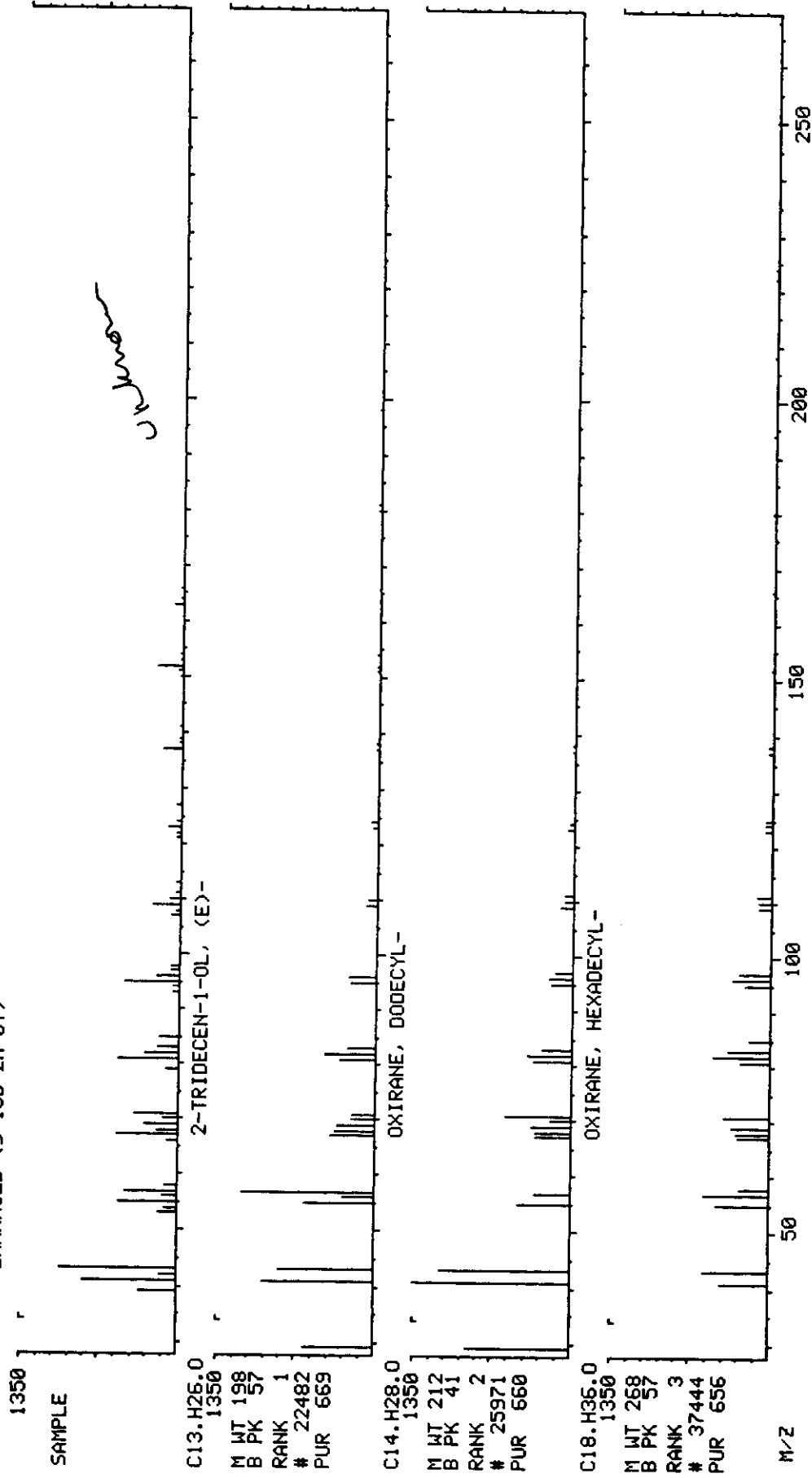
SAMPLE: S-MM5-1B 1/35A/100M

CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA) 1/35A NA M
ENHANCED (S 158 2N 0T)

DATA: 30068102 # 473
CALI: 30068102 # 3

BASE M/Z: 43
RIC: 41024.

INST. ID: F16



Library Search Data: 30068102 # 556 Base m/z: 43
 08/31/98 19:31:00 + 6:11 Cali: 30068102 # 3 RIC: 62784.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

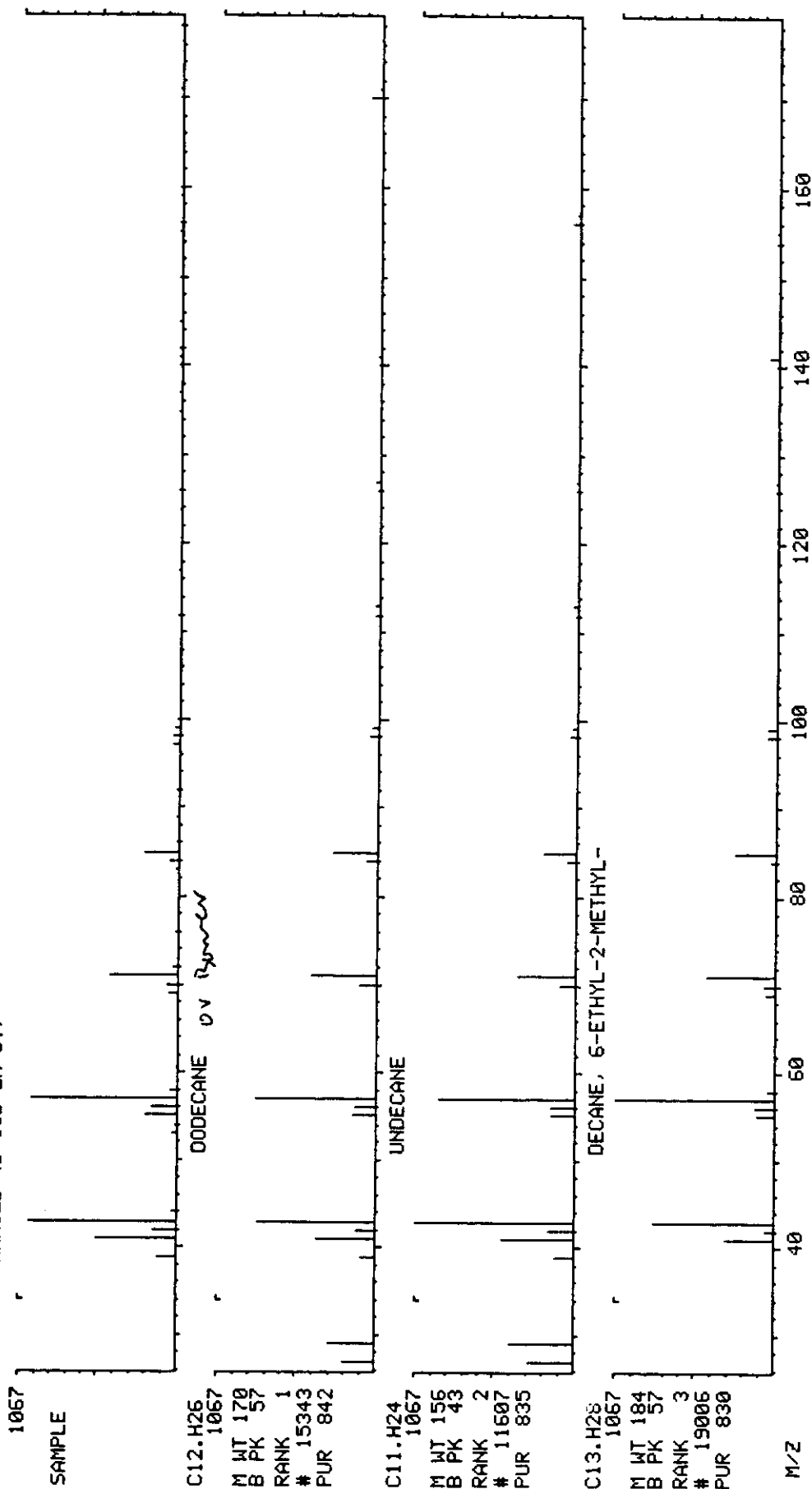
62231 spectra in LIBRARYNB searched for maximum PURITY
 990 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 15343 DODECANE
 2 11607 UNDECANE
 3 19006 DECANE, 6-ETHYL-2-METHYL-
 4 11602 OCTANE, 2,4,6-TRIMETHYL-
 5 19016 UNDECANE, 4,7-DIMETHYL-
 6 19015 DECANE, 2,5,6-TRIMETHYL-
 7 19028 DECANE, 2,4,6-TRIMETHYL-
 8 22530 TETRADECANE
 9 15359 UNDECANE, 5-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H26	170	57	842	943	868
2	C11.H24	156	43	835	955	856
3	C13.H28	184	57	830	945	842
4	C11.H24	156	57	825	957	836
5	C13.H28	184	43	819	933	845
6	C13.H28	184	57	817	917	833
7	C13.H28	184	43	812	916	852
8	C14.H30	198	43	810	904	883
9	C12.H26	170	43	808	929	829

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	112-40-3
2	—	—	—	—	1120-21-4
3	—	—	—	—	62108-21-8
4	—	—	—	—	62016-37-9
5	—	—	—	—	17301-32-5
6	—	—	—	—	62108-23-0
7	—	—	—	—	62108-27-4
8	—	—	—	—	629-59-4
9	—	—	—	—	1632-70-8

MID LIBRARY SEARCH <LIBRARYNB>
 08/31/98 19:31:00 + 5:11
 SAMPLE: S-MM5-1B 1/3SA/100M
 CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/3SA NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068102 # 555
 CALI: 30068102 # 3
 BASE M/Z: 43
 RIC: 62784.



Library Search Data: 30068102 # 570 Base m/z: 57
 08/31/98 19:31:00 + 6:20 Cali: 30068102 # 3 RIC: 47680.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYN8 searched for maximum PURITY
 876 matched at least 7 of the 16 largest peaks in the unknown

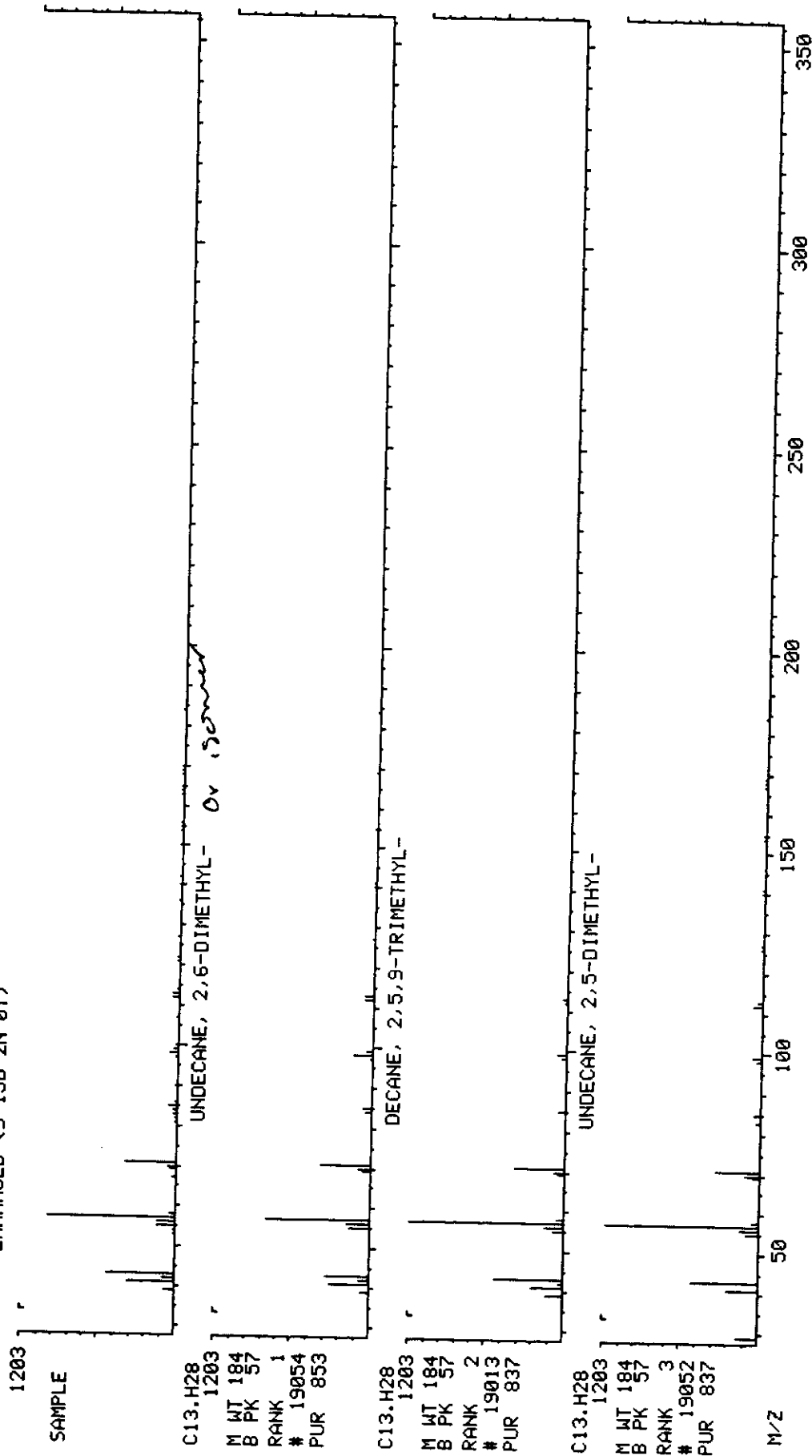
Rank In. Name
 1 19054 UNDECANE, 2,6-DIMETHYL-
 2 19013 DECANE, 2,5,9-TRIMETHYL-
 3 19052 UNDECANE, 2,5-DIMETHYL-
 4 19002 DODECANE, 6-METHYL-
 5 19026 DECANE, 2,6,8-TRIMETHYL-
 6 8104 OCTANE, 3,5-DIMETHYL-
 7 18996 UNDECANE, 3,6-DIMETHYL-
 8 19015 DECANE, 2,5,6-TRIMETHYL-
 9 8071 NONANE, 3-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H28	184	57	853	972	859
2	C13.H28	184	57	837	977	837
3	C13.H28	184	57	837	972	842
4	C13.H28	184	57	834	971	839
5	C13.H28	184	57	833	969	838
6	C10.H22	142	57	820	959	820
7	C13.H28	184	57	815	968	817
8	C13.H28	184	57	796	935	798
9	C10.H22	142	57	793	979	793

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	17301-23-4
2	—	—	—	—	62108-22-9
3	—	—	—	—	17301-22-3
4	—	—	—	—	6044-71-9
5	—	—	—	—	62108-26-3
6	—	—	—	—	15869-93-9
7	—	—	—	—	17301-28-9
8	—	—	—	—	62108-23-0
9	—	—	—	—	5911-04-6

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 6:20
 SAMPLE: S-MM5-1B 1/3SA/100M
 CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/3SA NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068102 # 570
 CALI: 30068102 # 3
 BASE M/Z: 57
 RIC: 47580.



Library Search Data: 30068102 # 617 Base m/z: 41
 08/31/98 19:31:00 + 6:52 Cali: 30068102 # 3 RIC: 34752.
 Sample: S-MMS-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 207 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 24994 1,1'-BICYCLOHEXYL, 2-(1-METHYLETHYL)-, TRANS-
 2 33979 BORINIC ACID, DIETHYL-, 1-CYCLOODECEN-1-YL ESTER
 3 24993 1,1'-BICYCLOHEXYL, 2-(1-METHYLETHYL)-, CIS-
 4 40193 CIS-9,10-EPOXYOCTADECAN-1-OL
 5 39227 NAPHTHALENE, DECAHYDRO-2,6-DIMETHYL-3-OCTYL-
 6 31486 ACETAMIDE, N-METHYL-N-[4-[4-METHOXY-1-HEXAHYDROPYRIDYL]-2-BUTYNYL]-
 7 24990 MUUROLANE-B
 8 39519 CYCLOHEXANE, 1-(1,5-DIMETHYLHEXYL)-4-(4-METHYLPENTYL)-
 9 24987 AMORPHANE-A

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C15.H28	208	83	595	901	619
2	C16.H31.O.8	250	83	584	913	612
3	C15.H28	208	83	580	877	621
4	C18.H36.O2	284	55	530	912	555
5	C20.H38	278	109	521	806	586
6	C13.H22.O2.N2	238	43	519	736	597
7	C15.H28	208	109	516	817	566
8	C20.H40	280	43	515	737	619
9	C15.H28	208	109	512	811	539

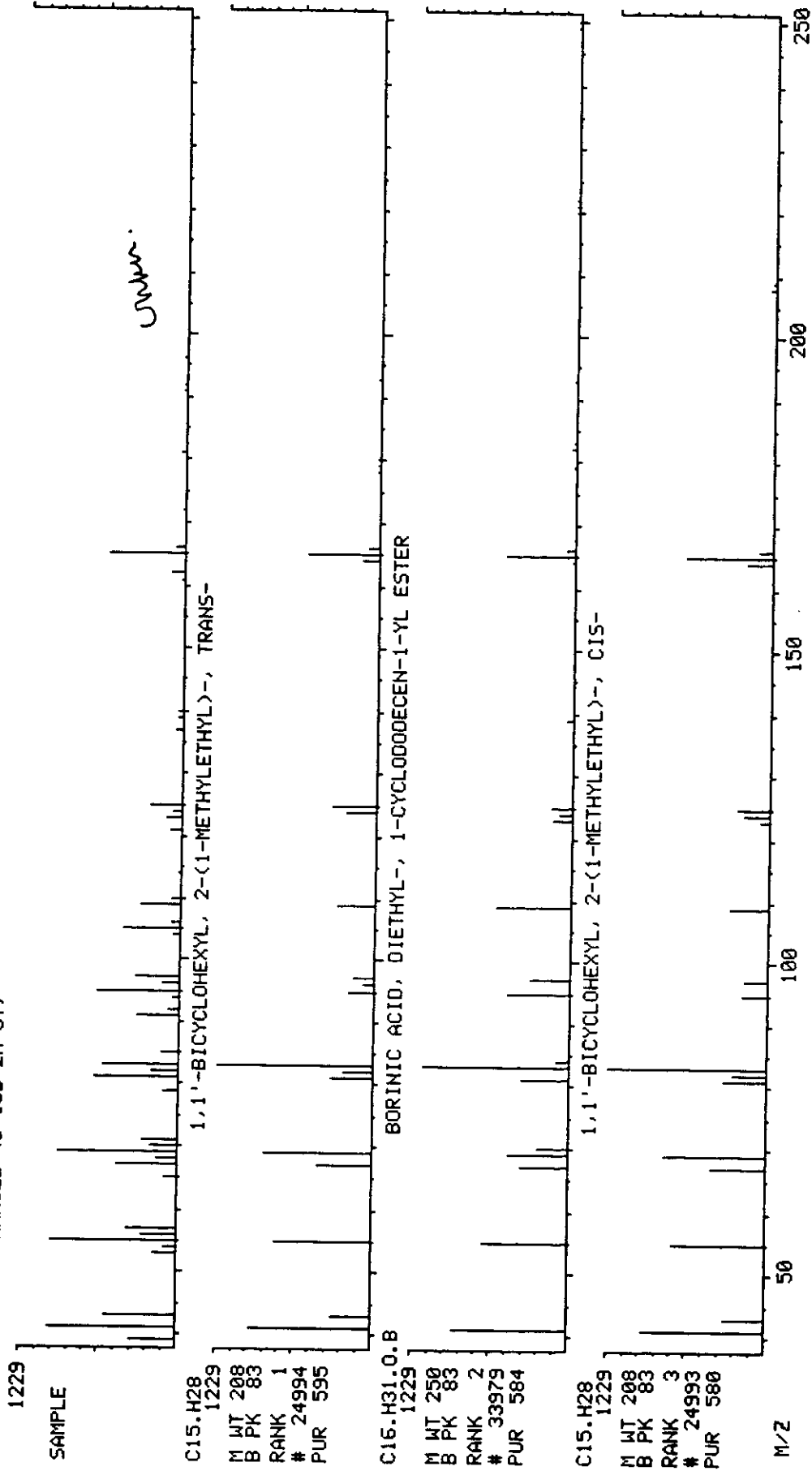
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	50991-16-7
2	—	—	—	—	61142-73-2
3	—	—	—	—	50991-15-6
4	—	—	—	—	13980-12-6
5	—	—	—	—	54964-85-1
6	—	—	—	—	- .
7	—	—	—	—	- .
8	—	—	—	—	56009-20-2
9	—	—	—	—	- .

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 6:52
 SAMPLE: S-MM5-1B 1/35A/100M
 CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA) 1/35A NA M
 ENHANCED (S 15B 2N 0T)

BASE M/Z: 41
 RIC: 34752.

DATA: 30068102 # 617
 CALI: 30068102 # 3

INST. ID: F16
 (NA/NA) 1/35A NA M



Library Search Data: 30068102 # 633 Base m/z: 57
 08/31/98 19:31:00 + 7:02 Cali: 30068102 # 3 RIC: 94336.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 735 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 14799 1-DECENE, 3,4-DIMETHYL-
 2 19523 1-DECANOL, 2-ETHYL-
 3 26408 2-ETHYL-1-DODECANOL
 4 37462 HEPTADECANE, 2,6-DIMETHYL-
 5 26409 2-HEXYL-1-OCTANOL
 6 32418 2-HEXYL-1-DECANOL
 7 22534 TRIDECANE, 7-METHYL-
 8 14751 2-UNDECENE, 5-METHYL-
 9 8539 1-PENTANOL, 4-METHYL-2-PROPYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H24	168	57	721	948	721
2	C12.H26.O	186	57	721	936	731
3	C14.H30.O	214	57	714	919	719
4	C19.H40	268	57	713	920	743
5	C14.H30.O	214	57	712	921	726
6	C16.H34.O	242	57	706	915	719
7	C14.H30	198	57	704	910	714
8	C12.H24	168	57	700	960	700
9	C9.H20.O	144	57	699	982	699

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	50871-03-9
2	—	—	—	—	21078-65-9
3	—	—	—	—	- -
4	—	—	—	—	54105-67-8
5	—	—	—	—	- -
6	—	—	—	—	- -
7	—	—	—	—	26730-14-3
8	—	—	—	—	56851-34-4
9	—	—	—	—	54004-41-0

Library Search Data: 30068102 # 654 Base m/z: 41
 08/31/98 19:31:00 + 7:16 Cali: 30068102 # 3 RIC: 44224.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 687 matched at least 7 of the 16 largest peaks in the unknown

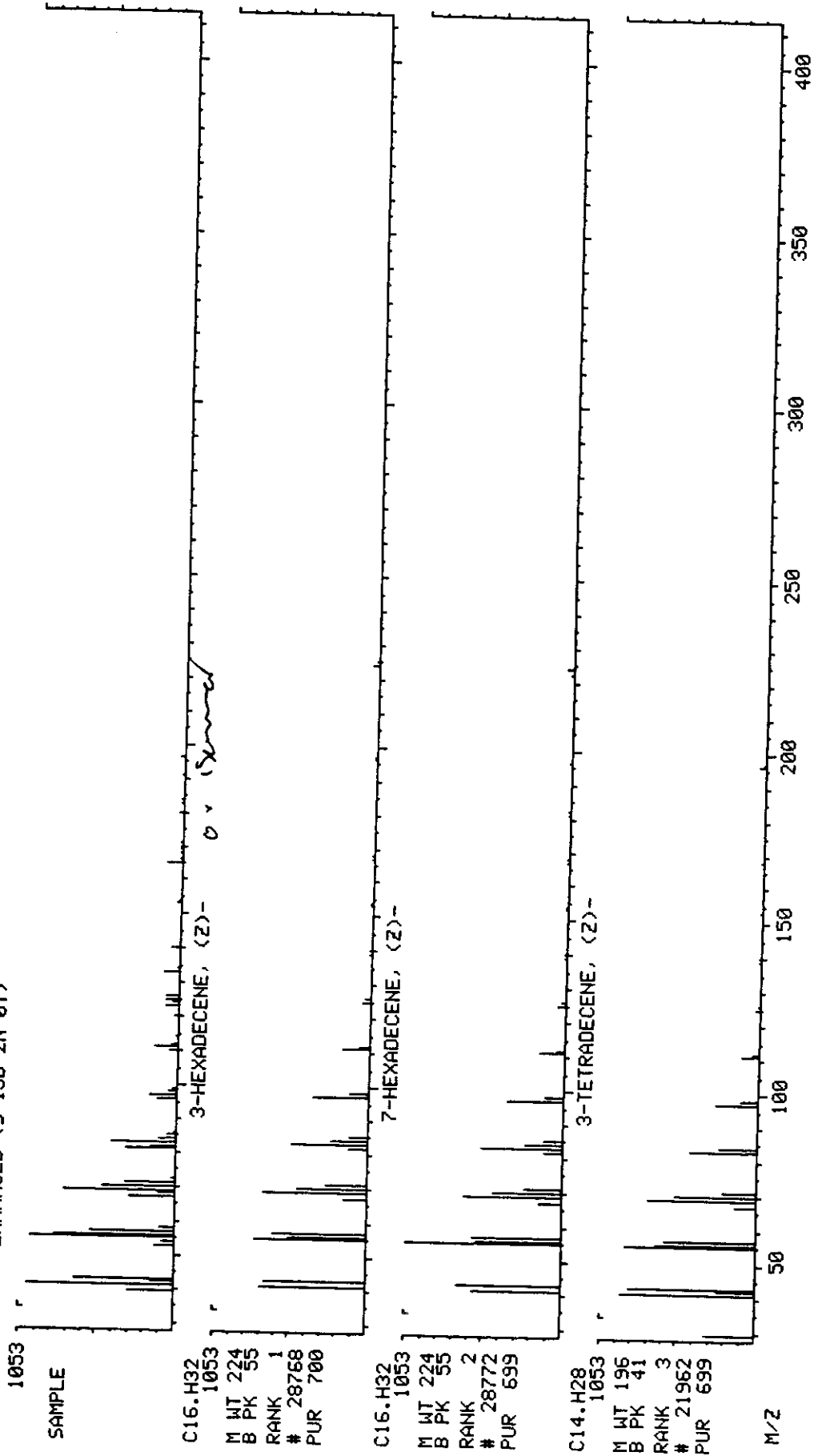
Rank In. Name
 1 28768 3-HEXADECENE, (Z)-
 2 28772 7-HEXADECENE, (Z)-
 3 21962 3-TETRADECENE, (Z)-
 4 21963 3-TETRADECENE, (E)-
 5 34410 3-OCTADECENE, (E)-
 6 32420 1-HEXADECANOL
 7 34411 5-OCTADECENE, (E)-
 8 21973 7-TETRADECENE, (Z)-
 9 19526 1-DODECANOL

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C16.H32	224	55	700	930	721
2	C16.H32	224	55	699	929	719
3	C14.H28	196	41	699	956	702
4	C14.H28	196	41	697	953	702
5	C18.H36	252	69	696	914	716
6	C16.H34.O	242	55	696	927	703
7	C18.H36	252	55	695	924	715
8	C14.H28	196	55	694	949	696
9	C12.H26.O	186	43	693	962	693

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	34303-81-6
2	—	—	—	—	35507-09-6
3	—	—	—	—	41446-67-7
4	—	—	—	—	41446-68-8
5	—	—	—	—	7206-19-1
6	—	—	—	—	36653-82-4
7	—	—	—	—	7206-21-5
8	—	—	—	—	41446-60-0
9	—	—	—	—	112-53-8

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 7:16
 SAMPLE: S-MM5-1B 1/35A/100M
 CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068102 # 654
 CALI: 30068102 # 3
 BASE M/Z: 41
 RIC: 44224.



Library Search Data: 30068102 # 663 Base m/z: 57
 08/31/98 19:31:00 + 7:22 Cali: 30068102 # 3 RIC: 86528.
 Sample: S-MM5-18 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYN8 searched for maximum PURITY
 787 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 11607 UNDECANE
 2 19013 DECANE, 2,5,9-TRIMETHYL-
 3 8104 OCTANE, 3,5-DIMETHYL-
 4 5159 NONANE
 5 11602 OCTANE, 2,4,6-TRIMETHYL-
 6 8073 DECANE
 7 15356 DECANE, 2,4-DIMETHYL-
 8 19016 UNDECANE, 4,7-DIMETHYL-
 9 19026 DECANE, 2,6,8-TRIMETHYL-

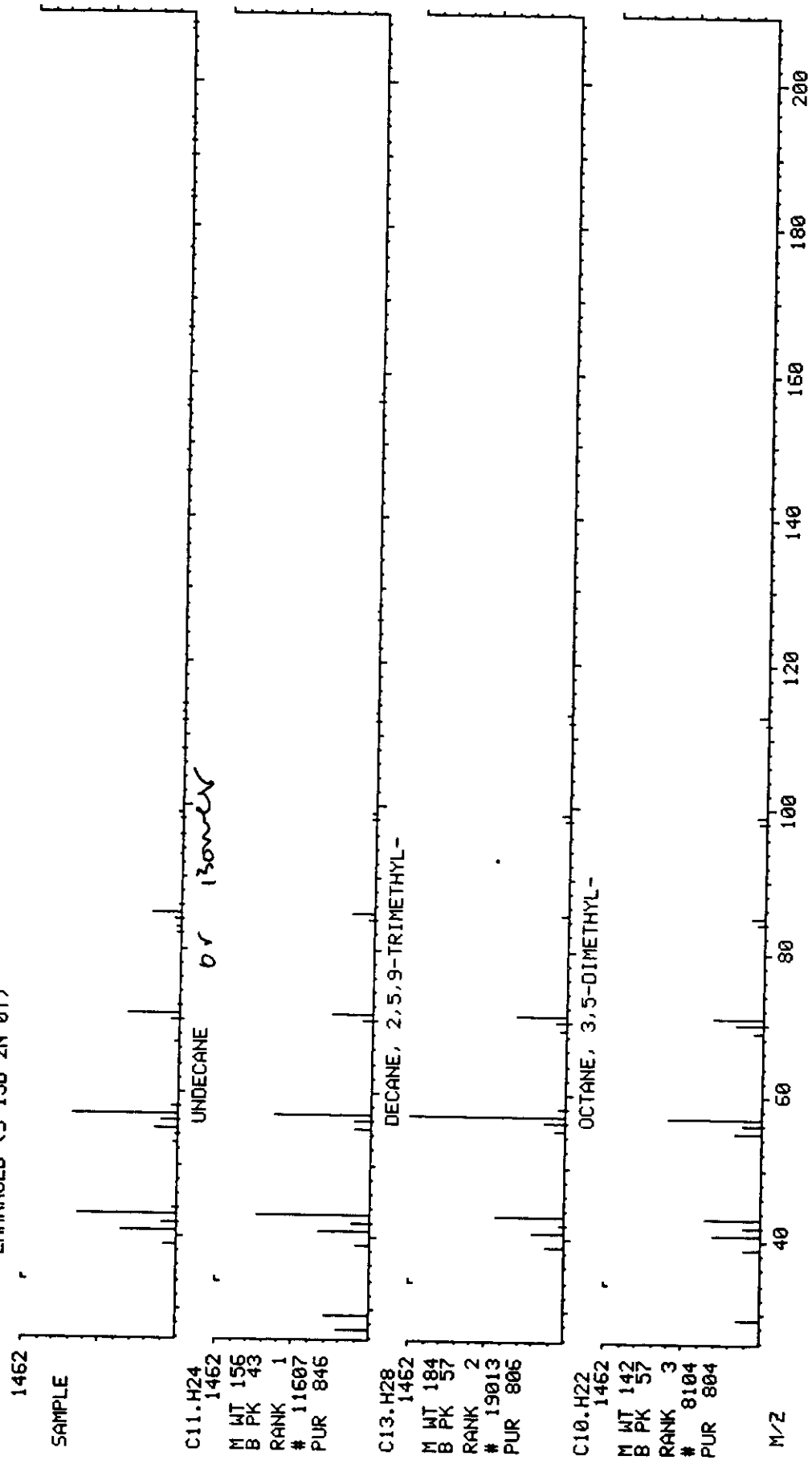
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C11.H24	156	43	846	960	858
2	C13.H28	184	57	806	921	820
3	C10.H22	142	57	804	913	832
4	C9.H20	128	43	803	929	842
5	C11.H24	156	57	799	938	831
6	C10.H22	142	43	797	922	850
7	C12.H26	170	43	794	930	823
8	C13.H28	184	43	793	904	846
9	C13.H28	184	57	792	907	826

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	1120-21-4
2	—	—	—	—	62108-22-9
3	—	—	—	—	15869-93-9
4	—	—	—	—	111-84-2
5	—	—	—	—	62016-37-9
6	—	—	—	—	124-18-5
7	—	—	—	—	2801-84-5
8	—	—	—	—	17301-32-5
9	—	—	—	—	62108-26-3

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 7:22
 SAMPLE: S-MMS-1B 1/35A/100M
 CONDS.: UG/ML *100ML *100%/(NA/NA) 1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068102 # 663
 CALI: 30068102 # 3
 BASE M/Z: 57
 RIC: 86528.

INST. ID: F16
 *(NA/NA) 1/35A NA M



Library Search Data: 30068102 # 678 Base m/z: 57
 08/31/98 19:31:00 + 7:32 Cali: 30068102 # 3 RIC: 45056.
 Sample: S-MM5-18 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

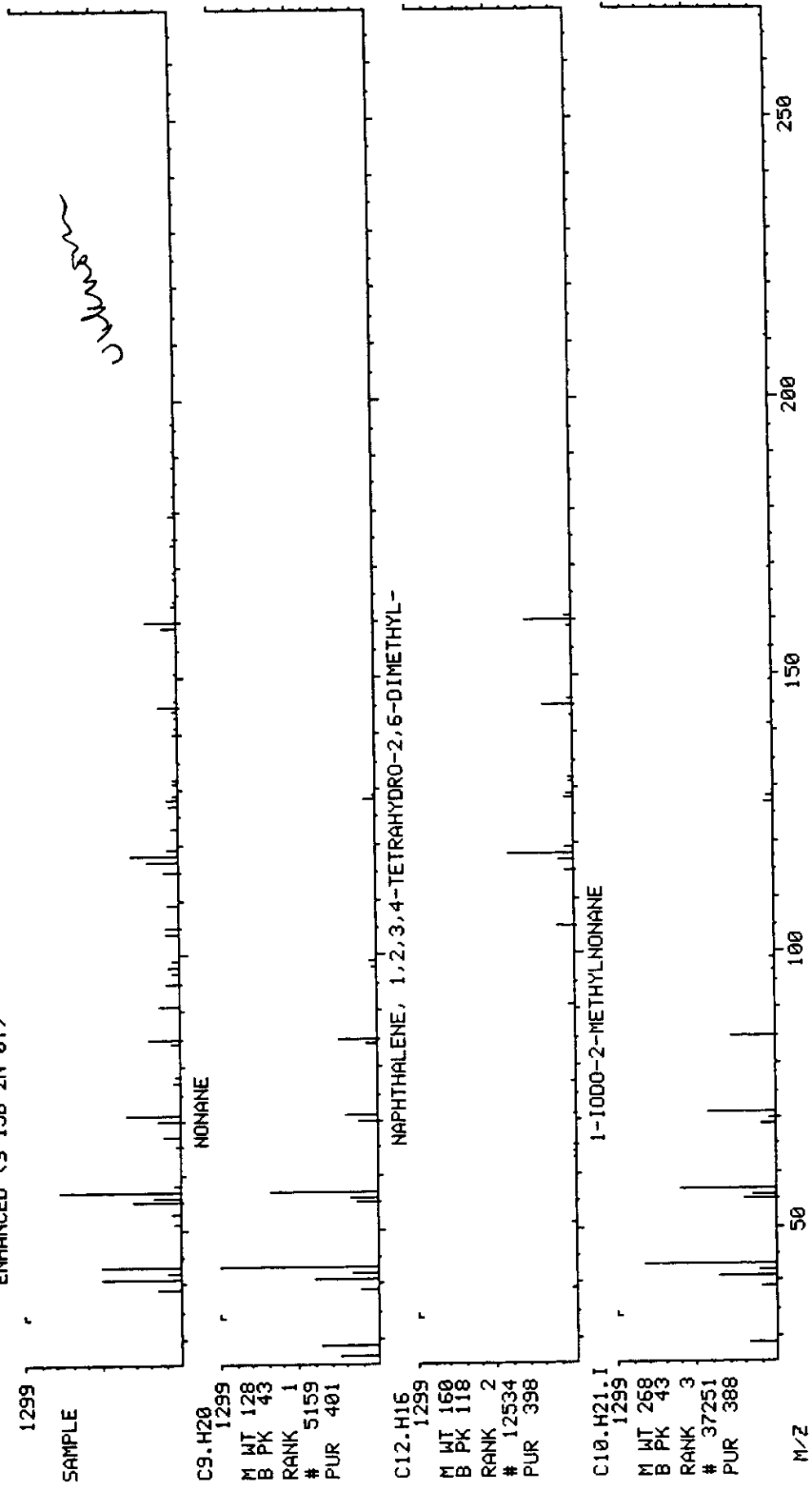
62231 spectra in LIBRARYNB searched for maximum PURITY
 672 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 5159 NONANE
 2 12534 NAPHTHALENE, 1,2,3,4-TETRAHYDRO-2,6-DIMETHYL-
 3 37251 1-iodo-2-METHYLNONANE
 4 11597 NONANE, 3,7-DIMETHYL-
 5 11602 OCTANE, 2,4,6-TRIMETHYL-
 6 8089 NONANE, 2-METHYL-
 7 15343 DODECANE
 8 11607 UNDECANE
 9 5041 3-HEXANONE, 2,4-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C9.H20	128	43	401	944	401
2	C12.H16	160	118	398	832	413
3	C10.H21.I	268	43	388	918	404
4	C11.H24	156	57	381	903	390
5	C11.H24	156	57	377	912	386
6	C10.H22	142	43	376	894	388
7	C12.H26	170	57	373	884	409
8	C11.H24	156	43	373	906	394
9	C8.H16.O	128	57	372	916	375

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	111-84-2
2	—	—	—	—	7524-63-2
3	—	—	—	—	—
4	—	—	—	—	17302-32-8
5	—	—	—	—	62016-37-9
6	—	—	—	—	871-83-0
7	—	—	—	—	112-40-3
8	—	—	—	—	1120-21-4
9	—	—	—	—	18641-70-8

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 7:32
 SAMPLE: 5-MM5-1B 1/35A/100M
 COND5.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068102 # 678
 CALI: 30068102 # 3
 BASE M/Z: 57
 RIC: 45056.



Library Search Data: 30068102 # 738 Base m/z: 57
 08/31/98 19:31:00 + 8:13 Cali: 30068102 # 3 RIC: 119808.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (\$ 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 712 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 25991 DODECANE, 2,6,10-TRIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 22535 DODECANE, 4,6-DIMETHYL-
 5 26001 DODECANE, 2,7,10-TRIMETHYL-
 6 25997 PENTADECANE
 7 25994 DODECANE, 2,6,11-TRIMETHYL-
 8 18998 UNDECANE, 3,7-DIMETHYL-
 9 15352 UNDECANE, 2-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	826	950	847
2	C15.H32	212	57	821	975	826
3	C21.H44	296	57	814	933	851
4	C14.H30	198	57	812	956	816
5	C15.H32	212	57	806	979	815
6	C15.H32	212	57	798	934	822
7	C15.H32	212	57	796	968	807
8	C13.H28	184	43	795	942	795
9	C12.H26	170	43	795	946	799

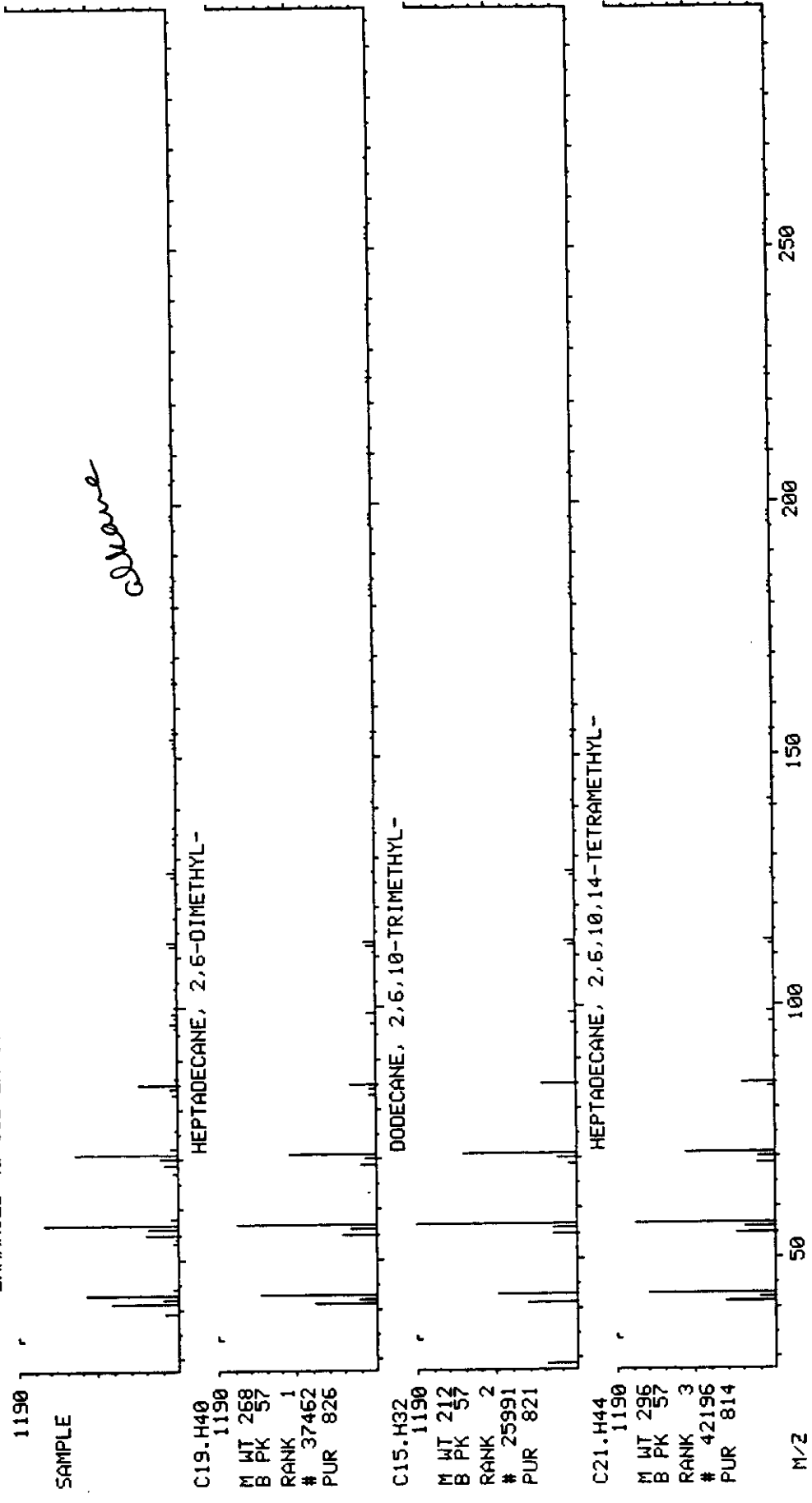
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	3891-98-3
3	—	—	—	—	18344-37-1
4	—	—	—	—	61141-72-8
5	—	—	—	—	74645-98-0
6	—	—	—	—	629-62-9
7	—	—	—	—	31295-56-4
8	—	—	—	—	17301-29-0
9	—	—	—	—	7045-71-8

BASE M/Z: 57
RIC: 119808.

DATA: 30068102 # 738
CALI: 30068102 # 3

MID LIBRARY SEARCH <LIBRARYNB>
08/31/98 19:31:00 + 8:13
SAMPLE: S-MM5-1B 1/35A/100M
CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA) 1/35A NA M
ENHANCED <S 15B 2N 0T>

INST. ID: F16



Library Search Data: 30068102 # 764 Base m/z: 43
 08/31/98 19:31:00 + 8:30 Cali: 30068102 # 3 RIC: 162304.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 672 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 19006 DECANE, 6-ETHYL-2-METHYL-
 2 22530 TETRADECANE
 3 19007 DODECANE, 3-METHYL-
 4 18990 UNDECANE, 2,9-DIMETHYL-
 5 11612 NONANE, 2,5-DIMETHYL-
 6 39681 1-iodoundecane
 7 37252 DECANE, 1-iodo-
 8 11607 UNDECANE
 9 22532 TRIDECANE, 2-METHYL-

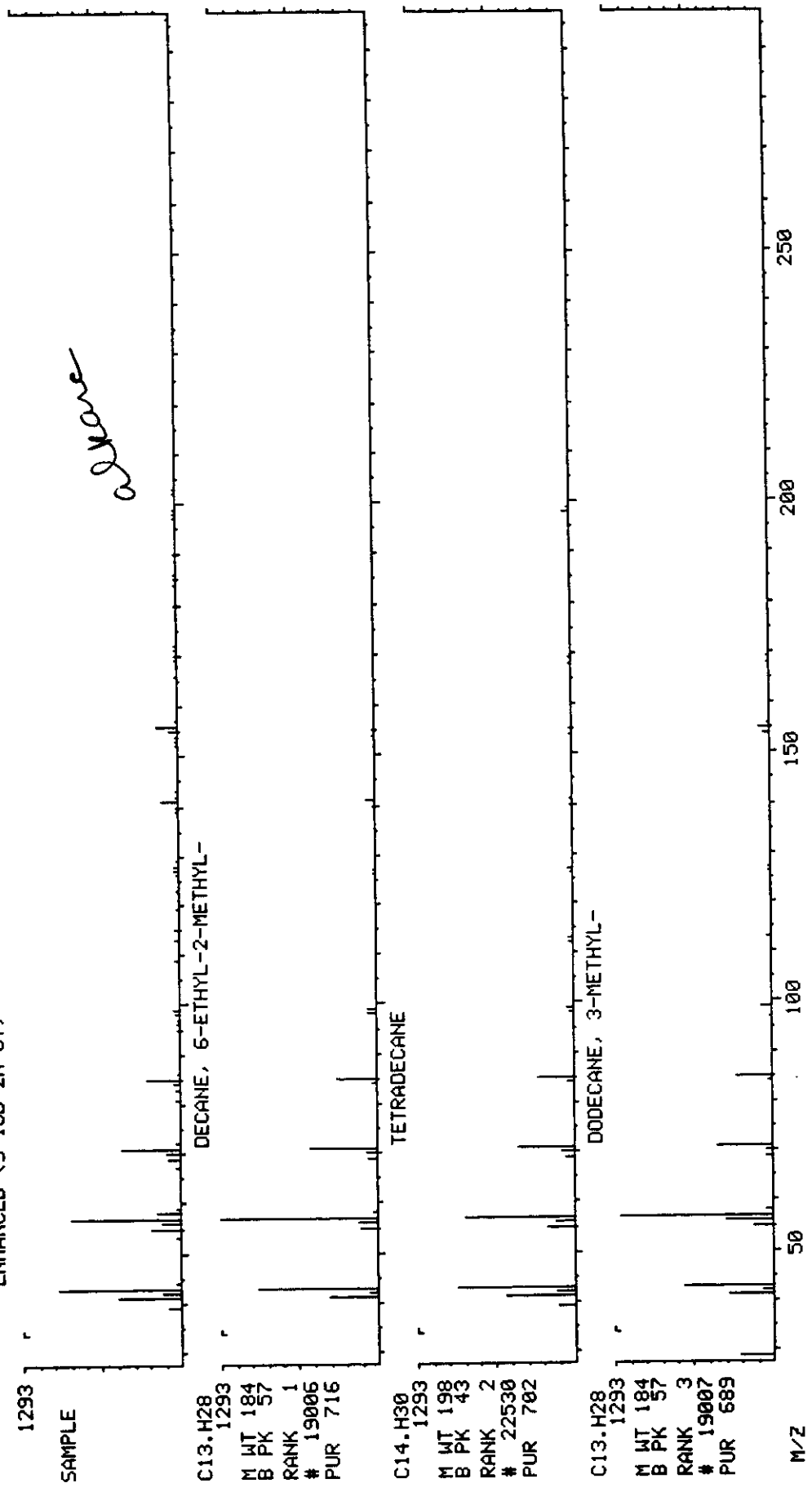
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H28	184	57	716	953	716
2	C14.H30	198	43	702	936	712
3	C13.H28	184	57	689	893	689
4	C13.H28	184	57	683	862	683
5	C11.H24	156	57	683	848	688
6	C11.H23.I	282	57	681	853	687
7	C10.H21.I	268	57	678	876	684
8	C11.H24	156	43	677	956	683
9	C14.H30	198	57	670	859	695

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	62108-21-8
2	—	—	—	—	629-59-4
3	—	—	—	—	17312-57-1
4	—	—	—	—	17301-26-7
5	—	—	—	—	17302-27-1
6	—	—	—	—	4282-44-4
7	—	—	—	—	2050-77-3
8	—	—	—	—	1120-21-4
9	—	—	—	—	1560-96-9

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 8:30
 SAMPLE: S-MMS-1B 1/35A/100M
 CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068102 # 764
 CALI: 30068102 # 3

BASE M/Z: 43
 RIC: 162304.



Library Search Data: 30068102 # 777 Base m/z: 41
 08/31/98 19:31:00 + 8:39 Cal: 30068102 # 3 RIC: 34368.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYN8 searched for maximum PURITY
 154 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 32052 OXIRANE, TETRADECYL-
 2 29187 BUTANOIC ACID, 3,7-DIMETHYL-6-OCTENYL ESTER
 3 29986 7-HYDROXY-3-(1,1-DIMETHYLPROP-2-ENYL)COUMARIN
 4 11518 CYCLOHEXANOL, 5-METHYL-2-(1-METHYLETHYL)-, [1S-(1.ALPHA.,2.BETA.,5.*
 5 35931 HEXADECANE, 1-CHLORO-
 6 50016 CETYLPIRIDINIUM CHLORIDE
 7 46252 UNDECANE, 1,2-DIBROMO-2-METHYL-
 8 32792 FALCARINOL (Z)-(-)-1,9-HEPTADECADIENE-4,6-DIYNE-3-OL
 9 40866 OCTADECANE, 1-CHLORO-

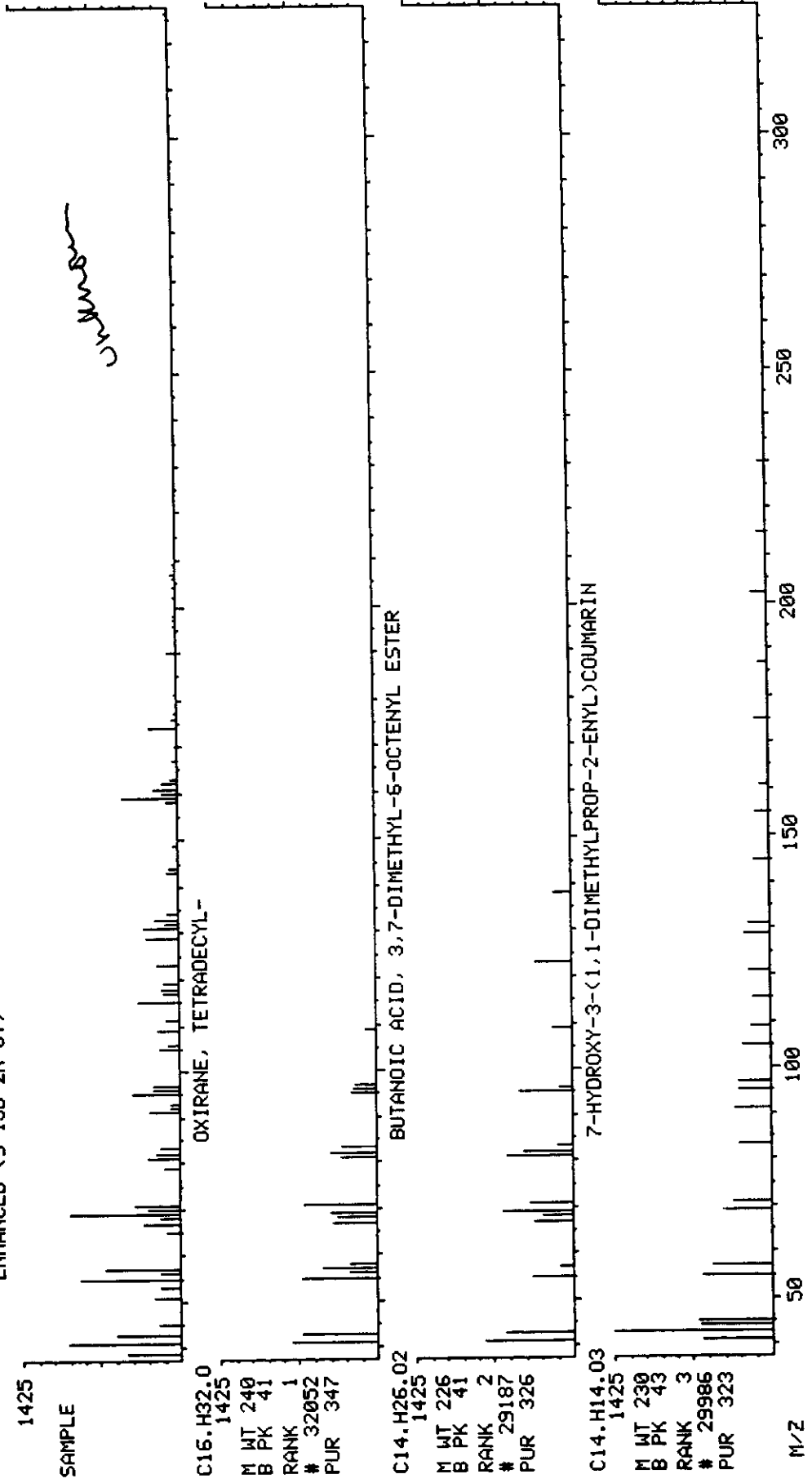
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C16.H32.O	240	41	347	911	359
2	C14.H26.O2	226	41	326	870	343
3	C14.H14.O3	230	43	323	696	420
4	C10.H20.O	156	71	313	833	333
5	C16.H33.CL	260	57	311	813	344
6	C21.H40.O.N.CL	357	55	305	824	334
7	C12.H24.BR2	326	41	294	808	349
8	C17.H24.O	244	55	288	706	388
9	C18.H37.CL	288	57	288	749	338

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	7320-37-8
2	—	—	—	—	141-16-2
3	—	—	—	—	56881-08-4
4	—	—	—	—	23283-97-8
5	—	—	—	—	4860-03-1
6	—	—	—	—	6004-24-6
7	—	—	—	—	55334-43-5
8	—	—	—	—	- -
9	—	—	—	—	3386-33-2

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 8:39
 SAMPLE: 5-MMS-1B 1/35A/100M
 CONDS.: UG/ML *100ML *100% *100% *1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068102 # 777
 CALI: 30068102 # 3

BASE M/Z: 41
 RIC: 34368.



Library Search Data: 30068102 # 820 Base m/z: 57
 08/31/98 19:31:00 + 9:07 Cali: 30068102 # 3 RIC: 117760.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 855 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 37465 NONADECANE
 4 25997 PENTADECANE
 5 18987 UNDECANE, 2,8-DIMETHYL-
 6 19016 UNDECANE, 4,7-DIMETHYL-
 7 25991 DODECANE, 2,6,10-TRIMETHYL-
 8 26001 DODECANE, 2,7,10-TRIMETHYL-
 9 15353 2,6-DIMETHYLDECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C21.H44	296	57	826	972	841
2	C19.H40	268	57	825	959	838
3	C19.H40	268	57	800	943	832
4	C15.H32	212	57	777	943	800
5	C13.H28	184	43	776	964	779
6	C13.H28	184	43	775	951	781
7	C15.H32	212	57	774	958	783
8	C15.H32	212	57	773	956	785
9	C12.H26	170	43	773	924	788

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	18344-37-1
2	—	—	—	—	54105-67-8
3	—	—	—	—	629-92-5
4	—	—	—	—	629-62-9
5	—	—	—	—	17301-25-6
6	—	—	—	—	17301-32-5
7	—	—	—	—	3891-98-3
8	—	—	—	—	74645-98-0
9	—	—	—	—	13150-81-7

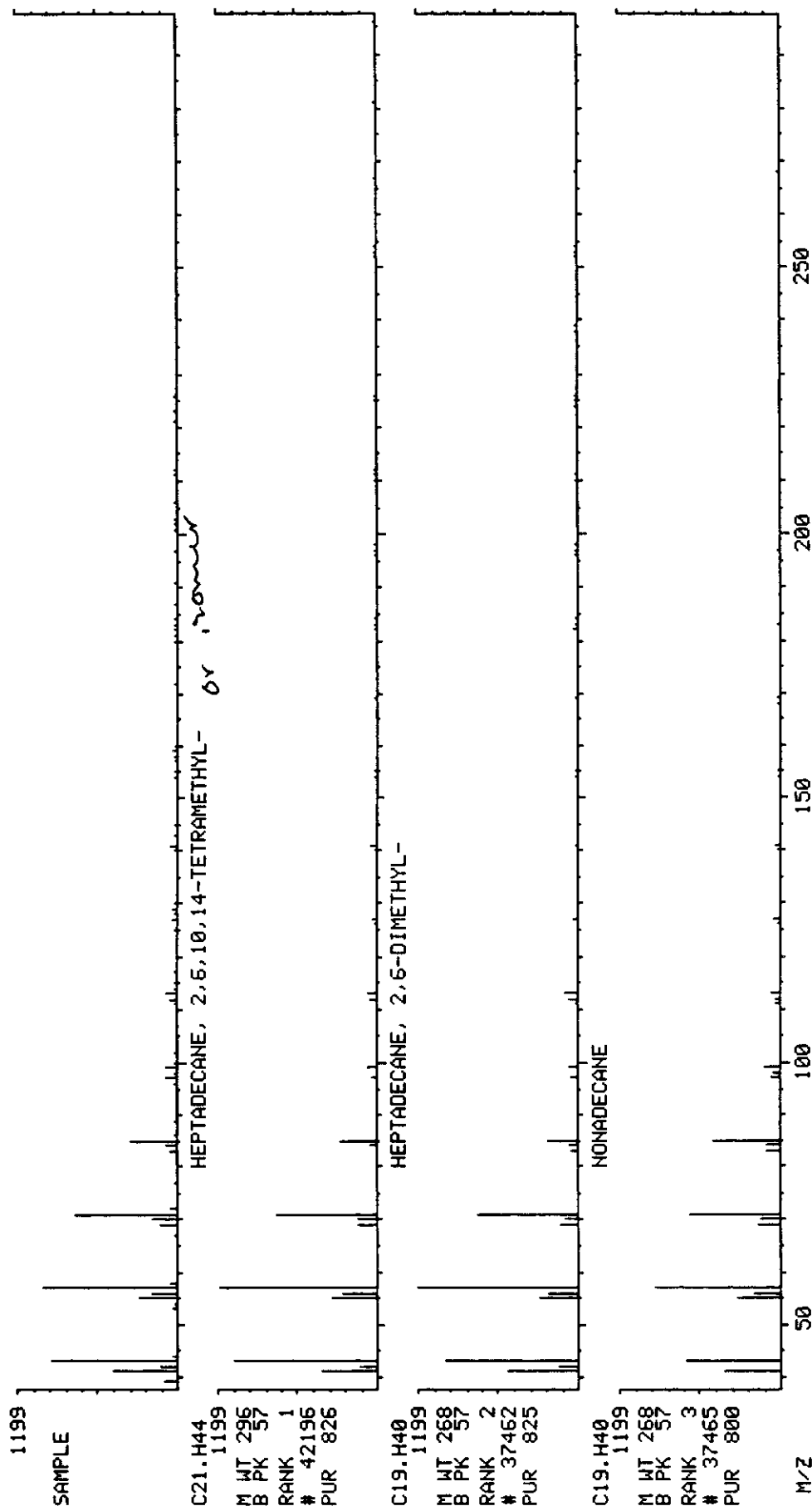
MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 9:07
 SAMPLE: 5-MMS-1B 1/35A/100M
 CONDS.: UG/ML *100ML *100% *100%
 ENHANCED (5 15B 2N 0T)

DATA: 30068102 # 820
 CALI: 30068102 # 3

BASE M/Z: 57
 RIC: 117760.

INST. ID: F16

(NA/NA) > 1/35A NA M



Library Search Data: 30068102 # 858 Base m/z: 57
 08/31/98 19:31:00 + 9:33 Cali: 30068102 # 3 RIC: 162816.
 Sample: S-MM5-18 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 835 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 19007 DODECANE, 3-METHYL-
 2 15352 UNDECANE, 2-METHYL-
 3 15353 2,6-DIMETHYLDECANE
 4 22535 DODECANE, 4,6-DIMETHYL-
 5 18998 UNDECANE, 3,7-DIMETHYL-
 6 22530 TETRADECANE
 7 18990 UNDECANE, 2,9-DIMETHYL-
 8 39681 1-IODOUNDECANE
 9 19006 DECANE, 6-ETHYL-2-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H28	184	57	773	964	774
2	C12.H26	170	43	771	935	771
3	C12.H26	170	43	753	908	757
4	C14.H30	198	57	753	946	755
5	C13.H28	184	43	752	938	752
6	C14.H30	198	43	751	928	762
7	C13.H28	184	57	750	958	751
8	C11.H23.I	282	57	749	965	756
9	C13.H28	184	57	746	949	747

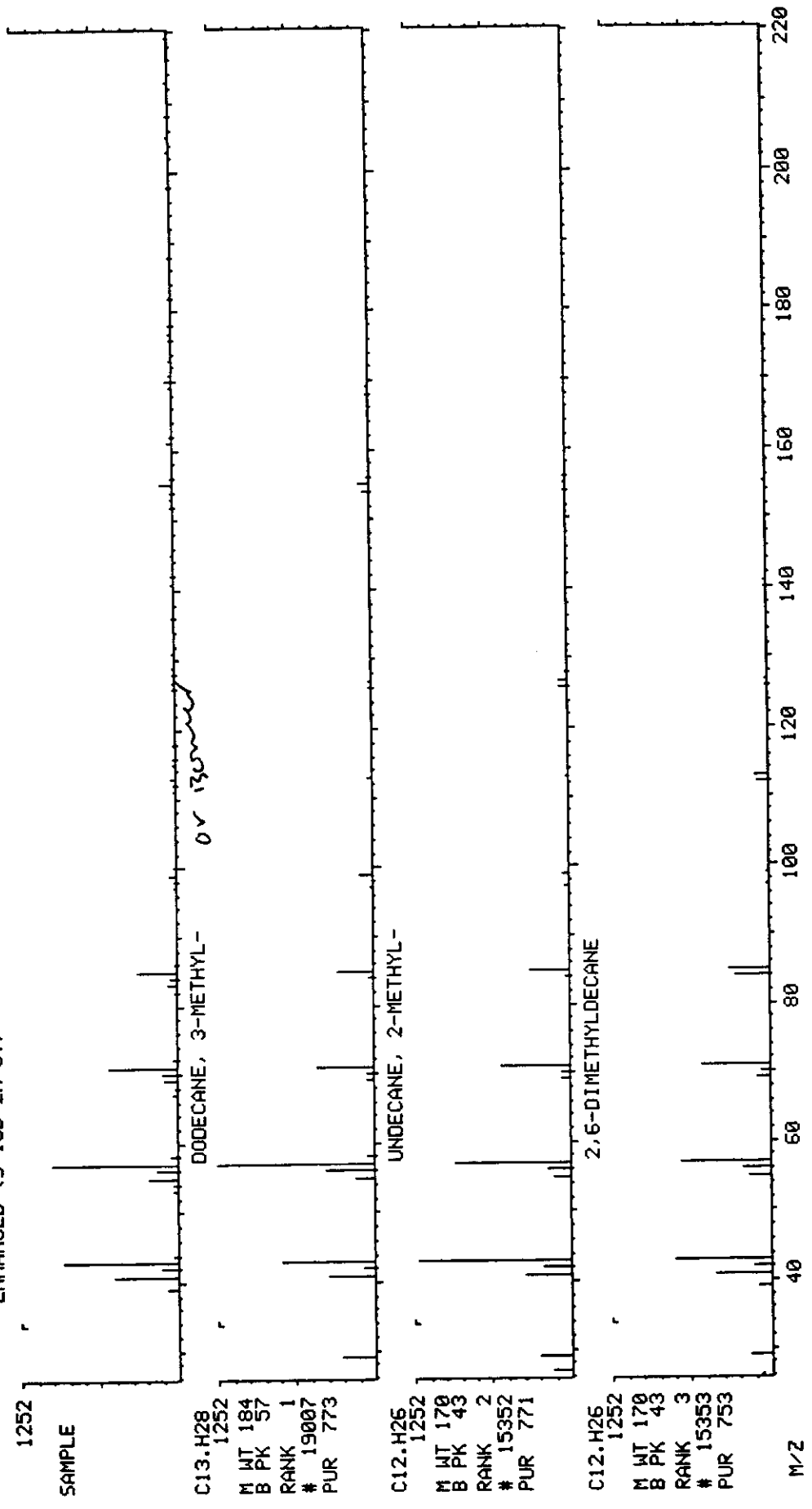
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	17312-57-1
2	---	---	---	---	7045-71-8
3	---	---	---	---	13150-81-7
4	---	---	---	---	61141-72-8
5	---	---	---	---	17301-29-0
6	---	---	---	---	629-59-4
7	---	---	---	---	17301-26-7
8	---	---	---	---	4282-44-4
9	---	---	---	---	62108-21-8

DATA: 30068102 # 858
CALI: 30068102 # 3

BASE M/Z: 57
RIC: 162816.

MTD LIBRARY SEARCH (LIBRARYNB)
08/31/98 19:31:00 + 9:33
SAMPLE: S-MMS-1B 1/35A/100M
CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 158 2N 0T)

INST. ID: F16



Library Search Data: 30068102 # 904 Base m/z: 43
 08/31/98 19:31:00 + 10:03 Cali: 30068102 # 3 RIC: 58816.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 650 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 40233 NONADECANOL
 2 42521 1-EICOSANOL
 3 42180 OCTADECANE, 1-(ETHENYLOXY)-
 4 37444 OXIRANE, HEXADECYL-
 5 25971 OXIRANE, DODECYL-
 6 58380 DECANEDIOIC ACID, DIDECYL ESTER
 7 37449 OCTADECANAL
 8 28772 7-HEXADECENE, (Z)-
 9 19525 2-METHYL-1-UNDECANOL

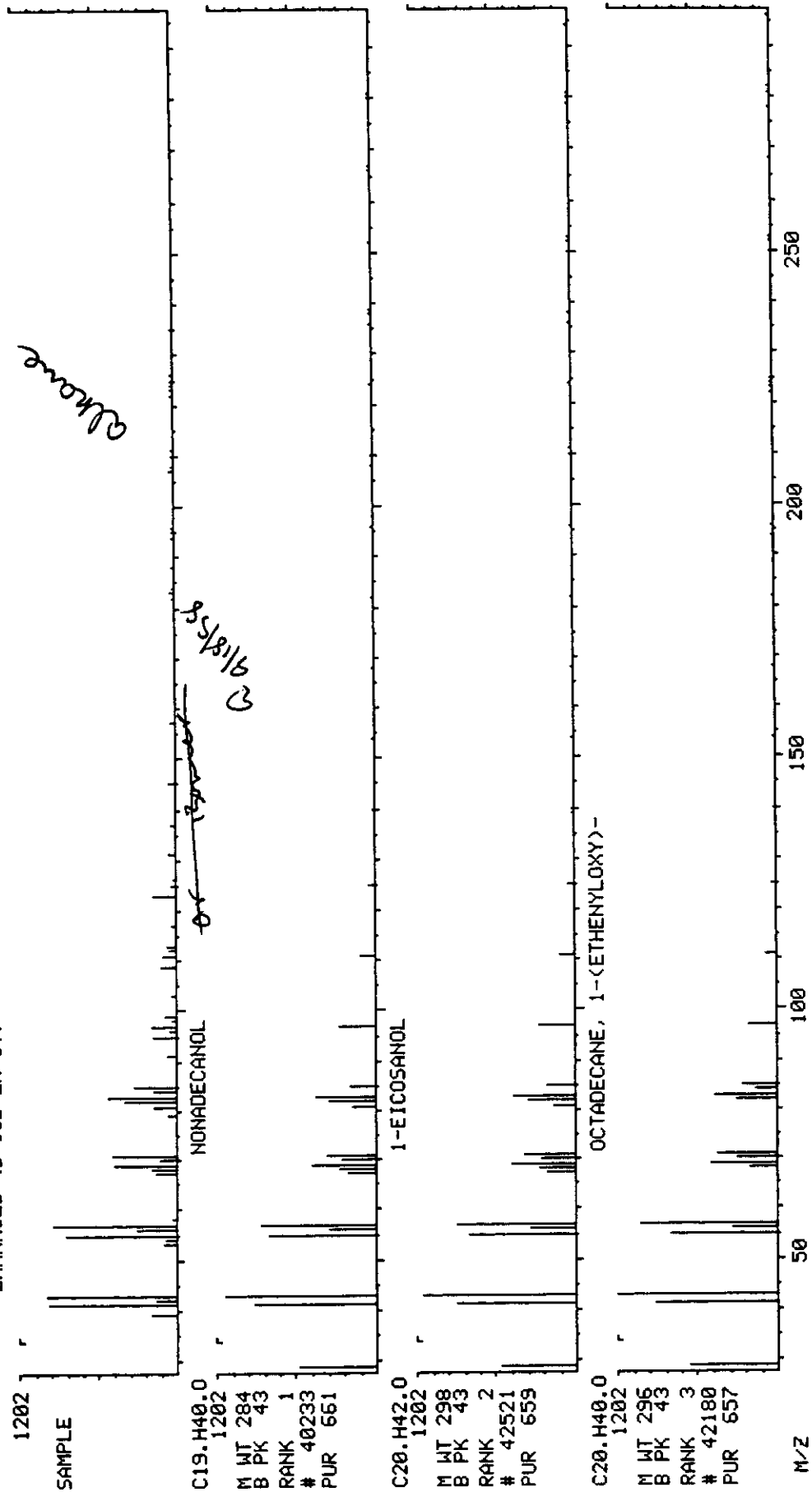
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40.O	284	43	661	938	693
2	C20.H42.O	298	43	659	935	695
3	C20.H40.O	296	43	657	948	684
4	C18.H36.O	268	57	648	853	693
5	C14.H28.O	212	41	630	889	656
6	C30.H58.O4	482	57	629	917	643
7	C18.H36.O	268	43	626	850	688
8	C16.H32	224	55	625	894	634
9	C12.H26.O	186	57	624	935	634

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	52783-43-4
2	—	—	—	—	629-96-9
3	—	—	—	—	930-02-9
4	—	—	—	—	7390-81-0
5	—	—	—	—	3234-28-4
6	—	—	—	—	2432-89-5
7	—	—	—	—	638-66-4
8	—	—	—	—	35507-09-6
9	—	—	—	—	10522-26-6

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 10:03
 SAMPLE: S-MMS-1B 1/35A/100M
 CONDS.: UG/ML *100Z/100Z *(NA/NA) >1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068102 # 904
 CALI: 30068102 # 3

BASE M/Z: 43
 RIC: 58816.



Library Search Data: 30068102 # 948 Base m/z: 57
 08/31/98 19:31:00 + 10:33 Cali: 30068102 # 3 RIC: 154112.
 Sample: S-MM5-18 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

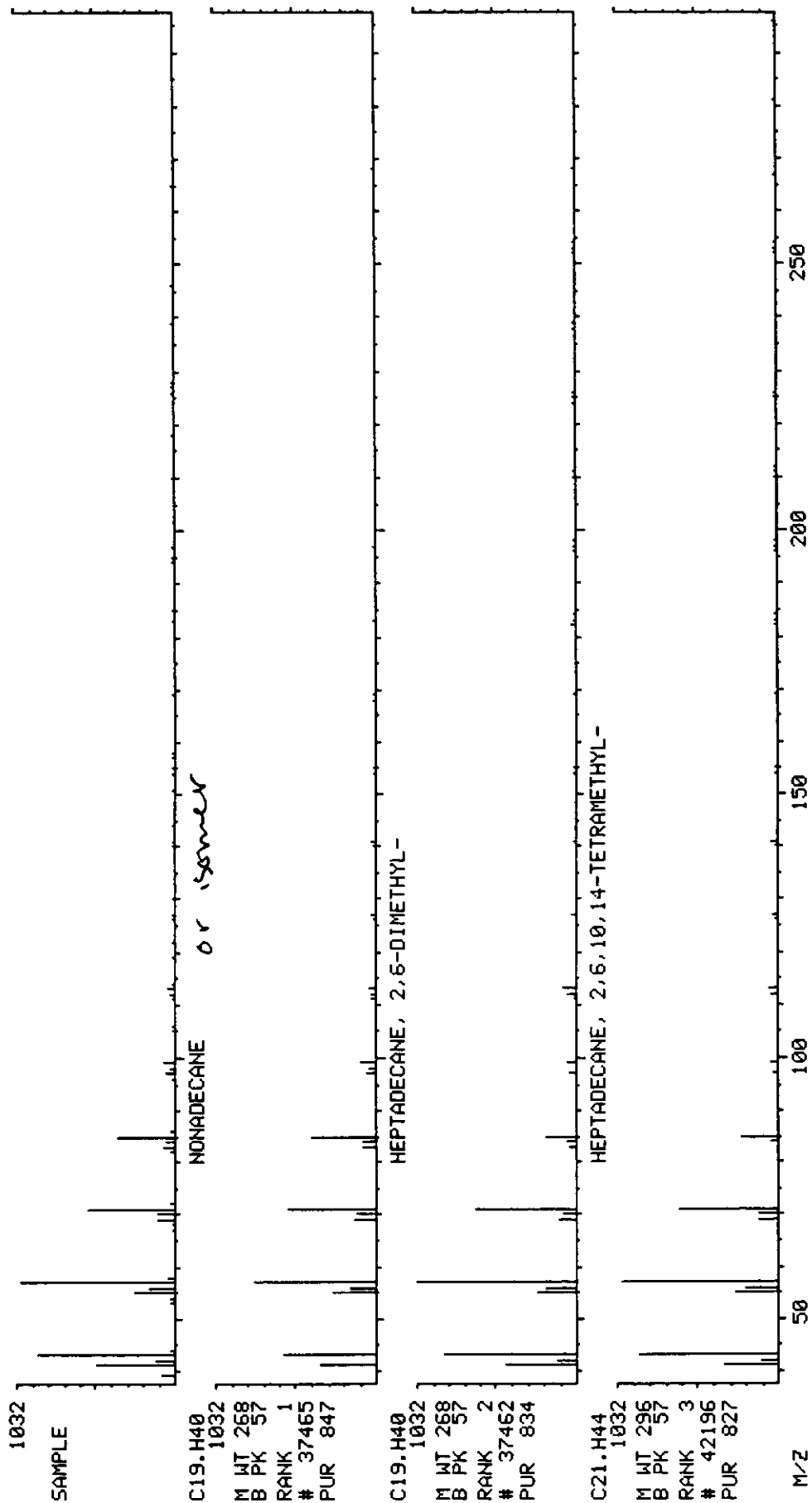
62231 spectra in LIBRARYNB searched for maximum PURITY
 360 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 15969 HYDROXYLAMINE, O-DECYL-
 5 25997 PENTADECANE
 6 19016 UNDECANE, 4,7-DIMETHYL-
 7 22535 DODECANE, 4,6-DIMETHYL-
 8 25994 DODECANE, 2,6,11-TRIMETHYL-
 9 15353 2,6-DIMETHYLDECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	847	958	866
2	C19.H40	268	57	834	942	858
3	C21.H44	296	57	827	950	846
4	C10.H23.O.N	173	43	825	967	827
5	C15.H32	212	57	817	954	842
6	C13.H28	184	43	812	964	813
7	C14.H30	198	57	801	951	805
8	C15.H32	212	57	799	965	802
9	C12.H26	170	43	797	927	807

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	629-92-5
2	---	---	---	---	54105-67-8
3	---	---	---	---	18344-37-1
4	---	---	---	---	29812-79-1
5	---	---	---	---	629-62-9
6	---	---	---	---	17301-32-5
7	---	---	---	---	61141-72-8
8	---	---	---	---	31295-56-4
9	---	---	---	---	13150-81-7

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 19:31:00 + 10:33
SAMPLE: S-MMS-1B 1/35A/100M
CONDS.: UG/ML *100ML *1007/1007
ENHANCED (S 15B 2N 0T)



Library Search Data: 30068102 # 988 Base m/z: 57
 08/31/98 19:31:00 + 10:59 Cali: 30068102 # 3 RIC: 114816.
 Sample: S-MMS-18 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYN8 searched for maximum PURITY
 644 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 25997 PENTADECANE
 3 19054 UNDECANE, 2,6-DIMETHYL-
 4 22530 TETRADECANE
 5 37462 HEPTADECANE, 2,6-DIMETHYL-
 6 25996 TRIDECANE, 4,8-DIMETHYL-
 7 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 8 29263 HEXADECANE
 9 32059 HEPTADECANE

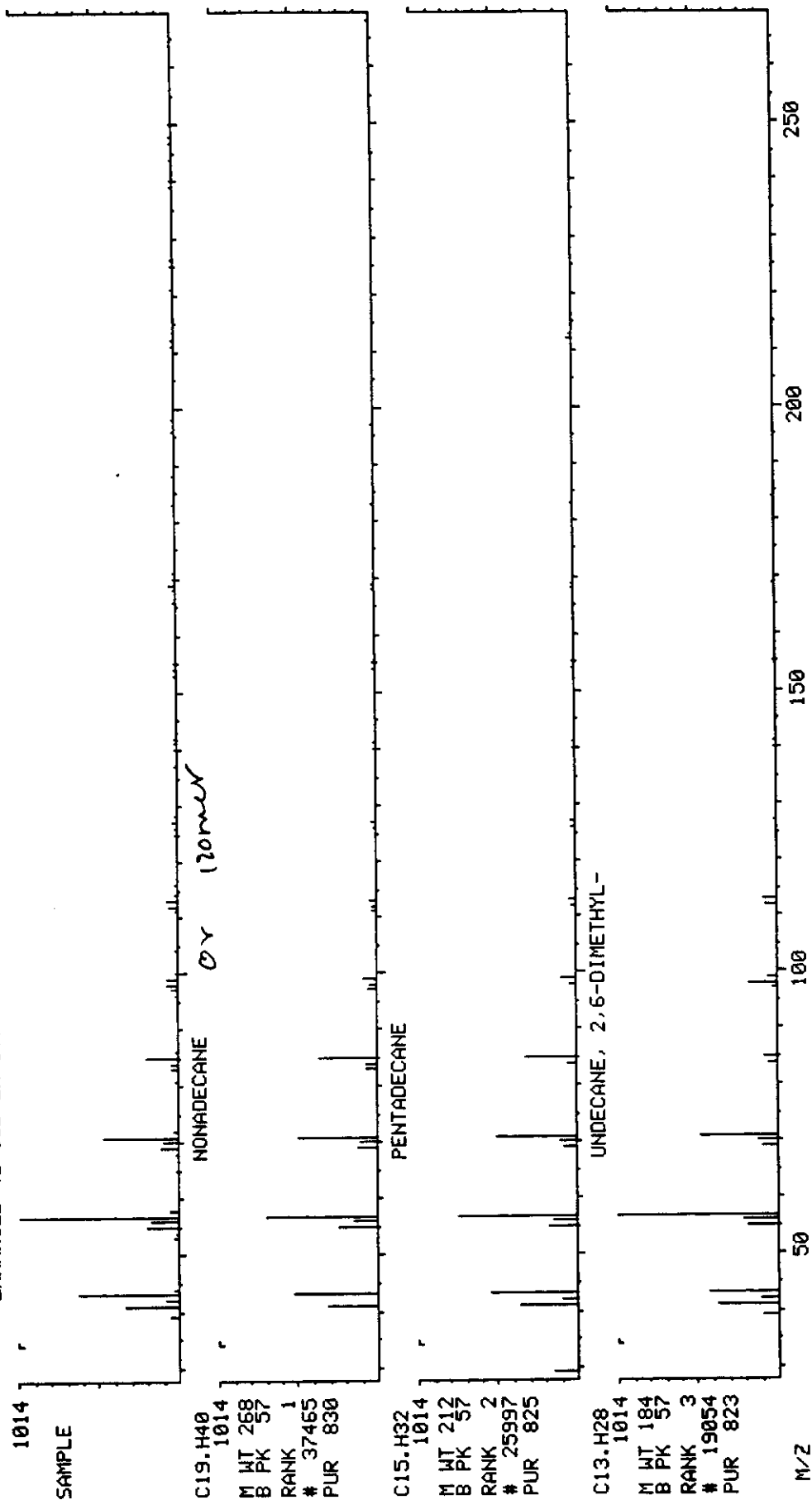
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	830	943	849
2	C15.H32	212	57	825	959	828
3	C13.H28	184	57	823	951	830
4	C14.H30	198	43	820	947	825
5	C19.H40	268	57	819	945	841
6	C15.H32	212	57	818	944	826
7	C21.H44	296	57	816	950	830
8	C16.H34	226	57	809	933	821
9	C17.H36	240	57	807	929	821

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	629-62-9
3	—	—	—	—	17301-23-4
4	—	—	—	—	629-59-4
5	—	—	—	—	54105-67-8
6	—	—	—	—	55030-62-1
7	—	—	—	—	18344-37-1
8	—	—	—	—	544-76-3
9	—	—	—	—	629-78-7

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 10:59
 SAMPLE: S-MMS-1B 1/35A/100M
 CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA) >1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068102 # 988
 CALI: 30068102 # 3

BASE M/Z: 57
 RIC: 114816.



Library Search Data: 30068102 #1036 Base m/z: 57
 08/31/98 19:31:00 + 11:31 Cali: 30068102 # 3 RIC: 176640.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 840 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 3 22535 DODECANE, 4,6-DIMETHYL-
 4 26001 DODECANE, 2,7,10-TRIMETHYL-
 5 25991 DODECANE, 2,6,10-TRIMETHYL-
 6 37465 NONADECANE
 7 37466 PENTADECANE, 2,6,10,14-TETRAMETHYL-
 8 25994 DODECANE, 2,6,11-TRIMETHYL-
 9 29264 TRIDECANE, 5-PROPYL-

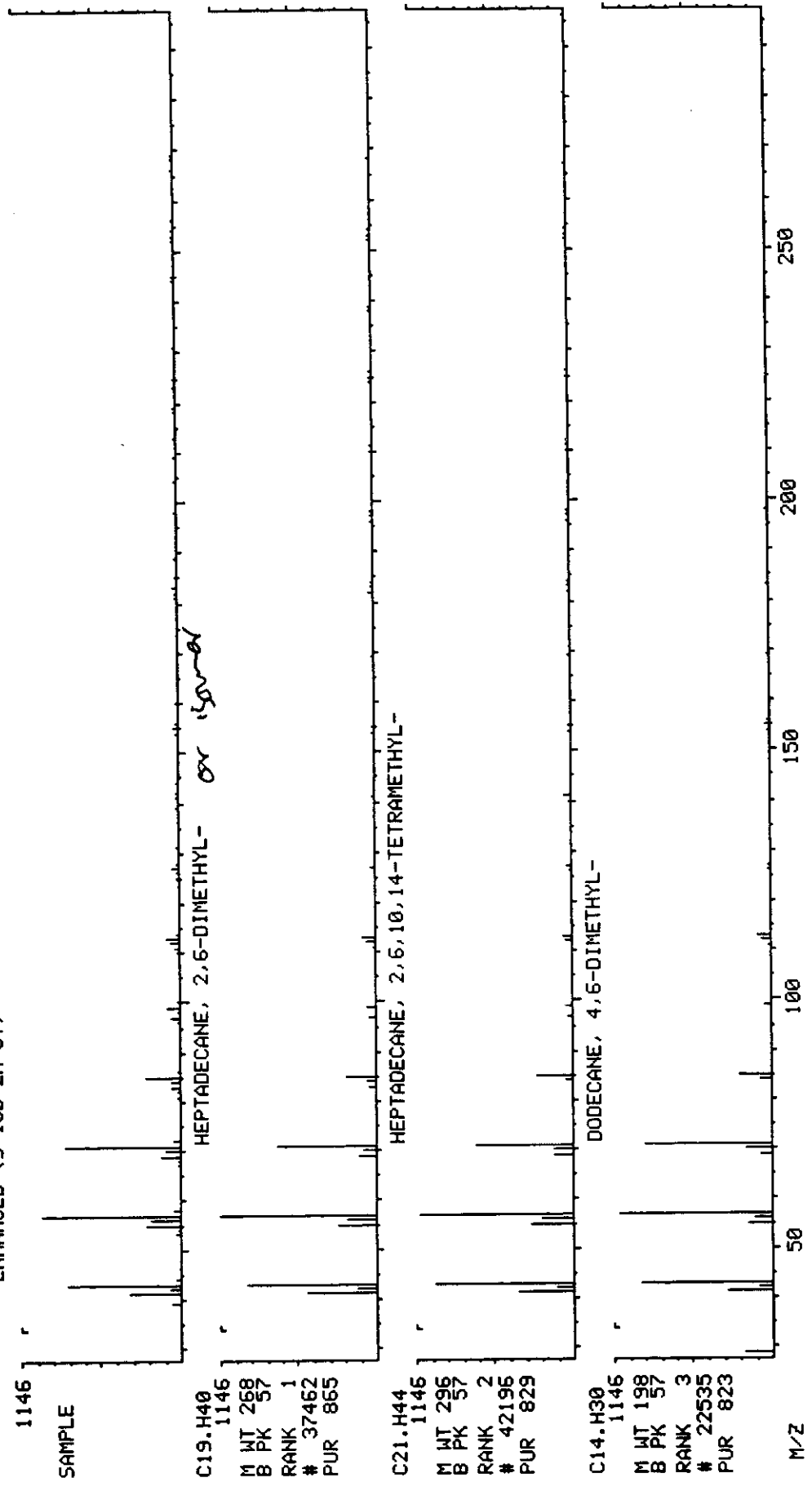
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	865	970	875
2	C21.H44	296	57	829	948	853
3	C14.H30	198	57	823	969	824
4	C15.H32	212	57	822	970	828
5	C15.H32	212	57	820	972	824
6	C19.H40	268	57	815	930	838
7	C19.H40	268	71	806	937	826
8	C15.H32	212	57	803	948	814
9	C16.H34	226	57	791	934	833

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	18344-37-1
3	—	—	—	—	61141-72-8
4	—	—	—	—	74645-98-0
5	—	—	—	—	3891-98-3
6	—	—	—	—	629-92-5
7	—	—	—	—	1921-70-6
8	—	—	—	—	31295-56-4
9	—	—	—	—	55045-11-9

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 11:31
 SAMPLE: S-MM5-1B 1/35A/100M
 CONDS.: UG/ML *100ML *100% *100% (NA/NA) 1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068102 #1036
 CALI: 30068102 # 3

BASE M/Z: 57
 RIC: 176640.



Library Search Data: 30068102 #1054 Base m/z: 55
 08/31/98 19:31:00 + 11:43 Cali: 30068102 # 3 RIC: 48576.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 648 matched at least 7 of the 16 largest peaks in the unknown

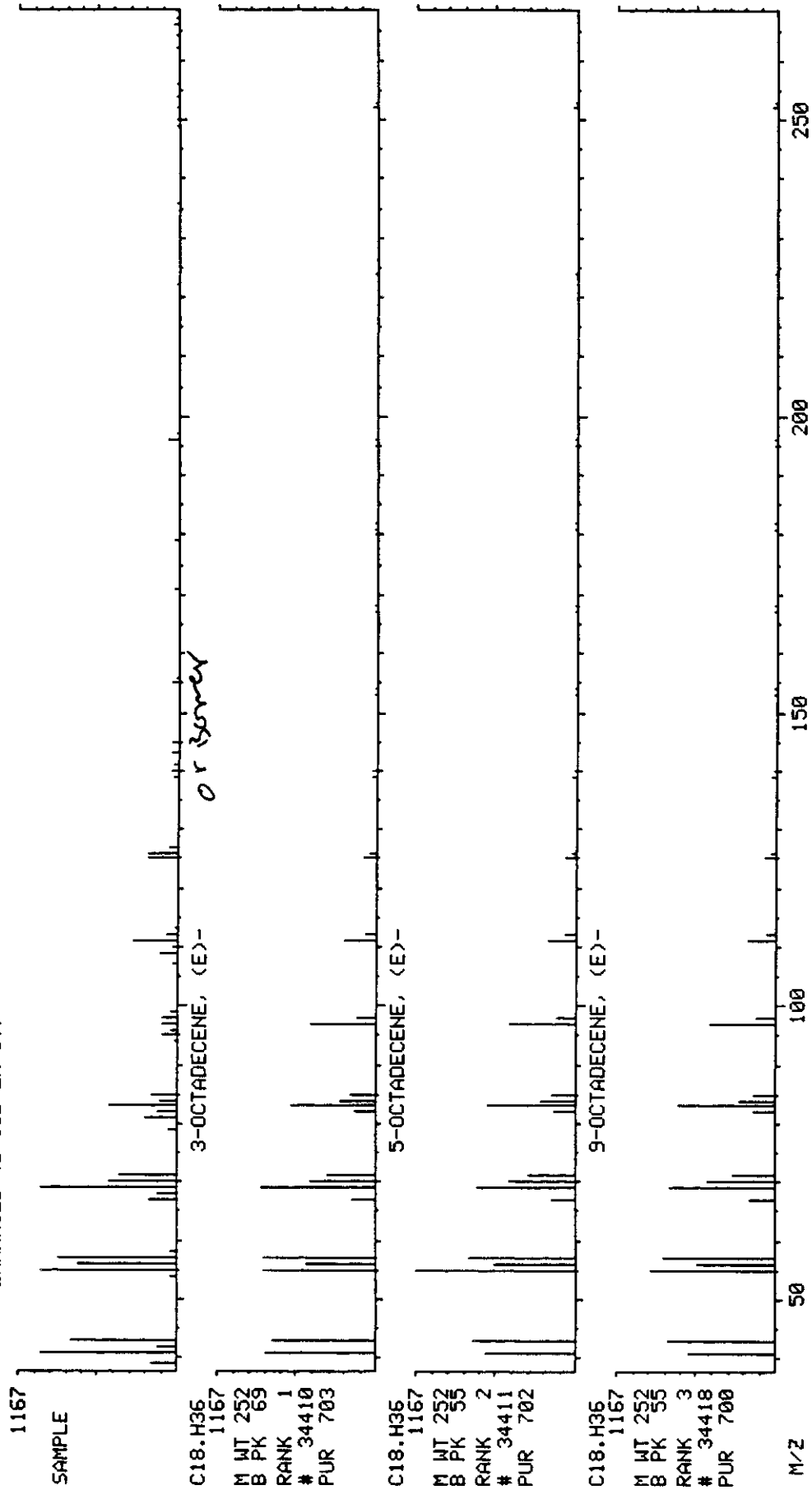
Rank In. Name
 1 34410 3-OCTADECENE, (E)-
 2 34411 5-OCTADECENE, (E)-
 3 34418 9-OCTADECENE, (E)-
 4 28768 3-HEXADECENE, (Z)-
 5 32425 1-DECANOL, 2-HEXYL-
 6 26409 2-HEXYL-1-OCTANOL
 7 28772 7-HEXADECENE, (Z)-
 8 21963 3-TETRADECENE, (E)-
 9 32418 2-HEXYL-1-DECANOL

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C18.H36	252	69	703	906	717
2	C18.H36	252	55	702	904	713
3	C18.H36	252	55	700	901	711
4	C16.H32	224	55	700	903	722
5	C16.H34.O	242	57	698	911	720
6	C14.H30.O	214	57	696	924	703
7	C16.H32	224	55	694	895	713
8	C14.H28	196	41	693	925	696
9	C16.H34.O	242	57	692	916	704

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	7206-19-1
2	—	—	—	—	7206-21-5
3	—	—	—	—	7206-25-9
4	—	—	—	—	34303-81-6
5	—	—	—	—	2425-77-6
6	—	—	—	—	- -
7	—	—	—	—	35507-09-6
8	—	—	—	—	41446-68-8
9	—	—	—	—	- -

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 11:43
 SAMPLE: S-MMS-1B 1/35A/100M
 CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA) >1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068102 #1054
 CALI: 30068102 # 3
 BASE M/Z: 55
 RIC: 48576.



Library Search Data: 30068102 #1114 Base m/z: 57
 08/31/98 19:31:00 + 12:23 Cali: 30068102 # 3 RIC: 80128.
 Sample: S-MM5-18 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 360 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 15969 HYDROXYLAMINE, O-DECYL-
 2 19016 UNDECANE, 4,7-DIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 37462 HEPTADECANE, 2,6-DIMETHYL-
 5 37465 NONADECANE
 6 11602 OCTANE, 2,4,6-TRIMETHYL-
 7 11612 NONANE, 2,5-DIMETHYL-
 8 25997 PENTADECANE
 9 15353 2,6-DIMETHYLDECANE

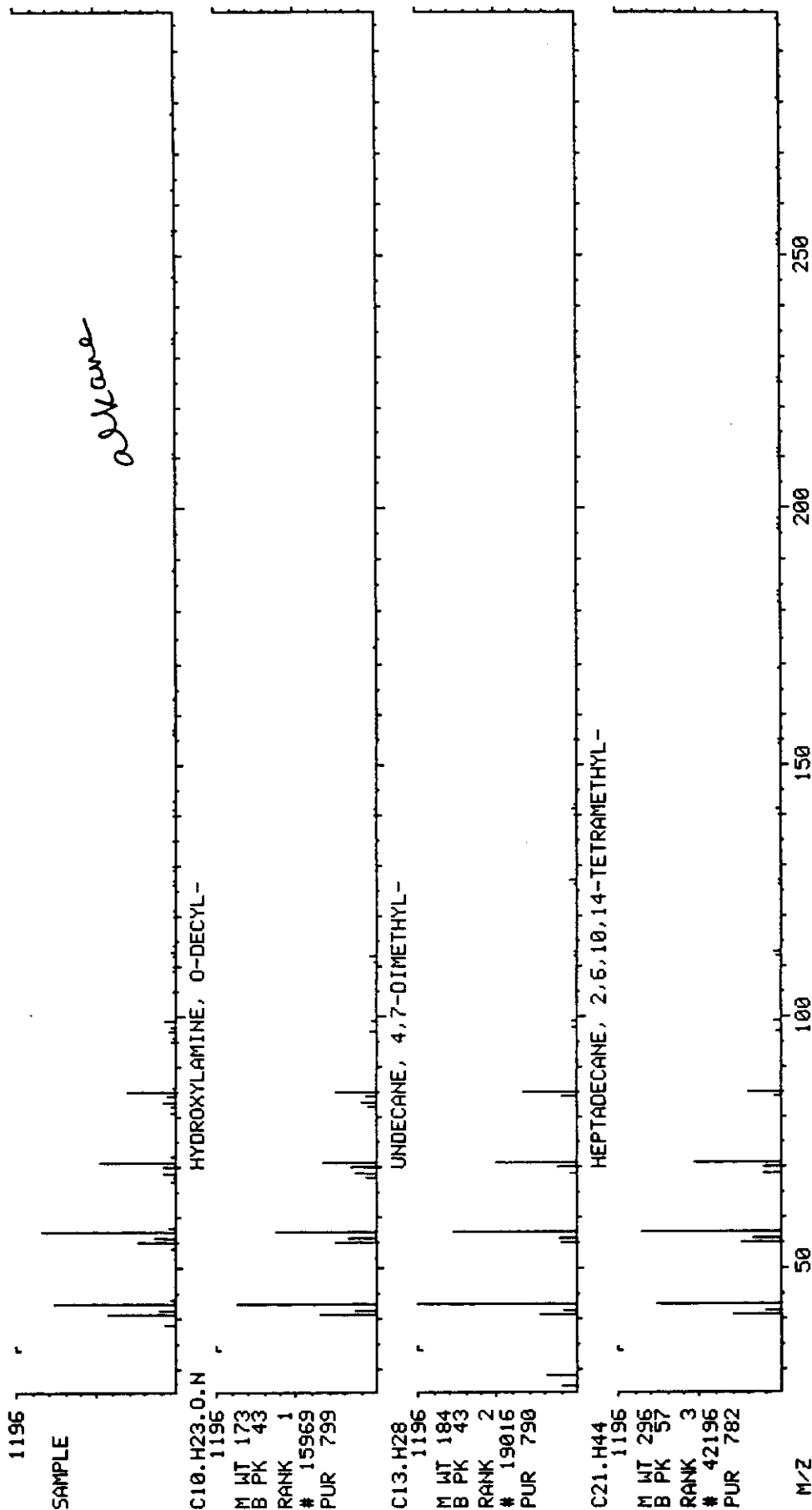
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C10.H23.O.N	173	43	799	957	808
2	C13.H28	184	43	790	966	792
3	C21.H44	296	57	782	940	814
4	C19.H40	268	57	777	917	823
5	C19.H40	268	57	773	919	829
6	C11.H24	156	57	765	968	765
7	C11.H24	156	57	762	913	762
8	C15.H32	212	57	761	930	807
9	C12.H26	170	43	760	916	775

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	29812-79-1
2	---	---	---	---	17301-32-5
3	---	---	---	---	18344-37-1
4	---	---	---	---	54105-67-8
5	---	---	---	---	629-92-5
6	---	---	---	---	62016-37-9
7	---	---	---	---	17302-27-1
8	---	---	---	---	629-62-9
9	---	---	---	---	13150-81-7

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 12:23
 SAMPLE: S-MM5-1B 1/35A/100M
 COND.: UG/ML *100ML *100Z/100Z *(NA/NA) >1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068102 #1114
 CALI: 30068102 # 3

BASE M/Z: 57
 RIC: 80128.



Library Search Data: 30068102 #1119 Base m/z: 57
 08/31/98 19:31:00 + 12:27 Cali: 30068102 # 3 RIC: 112896.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N QT)

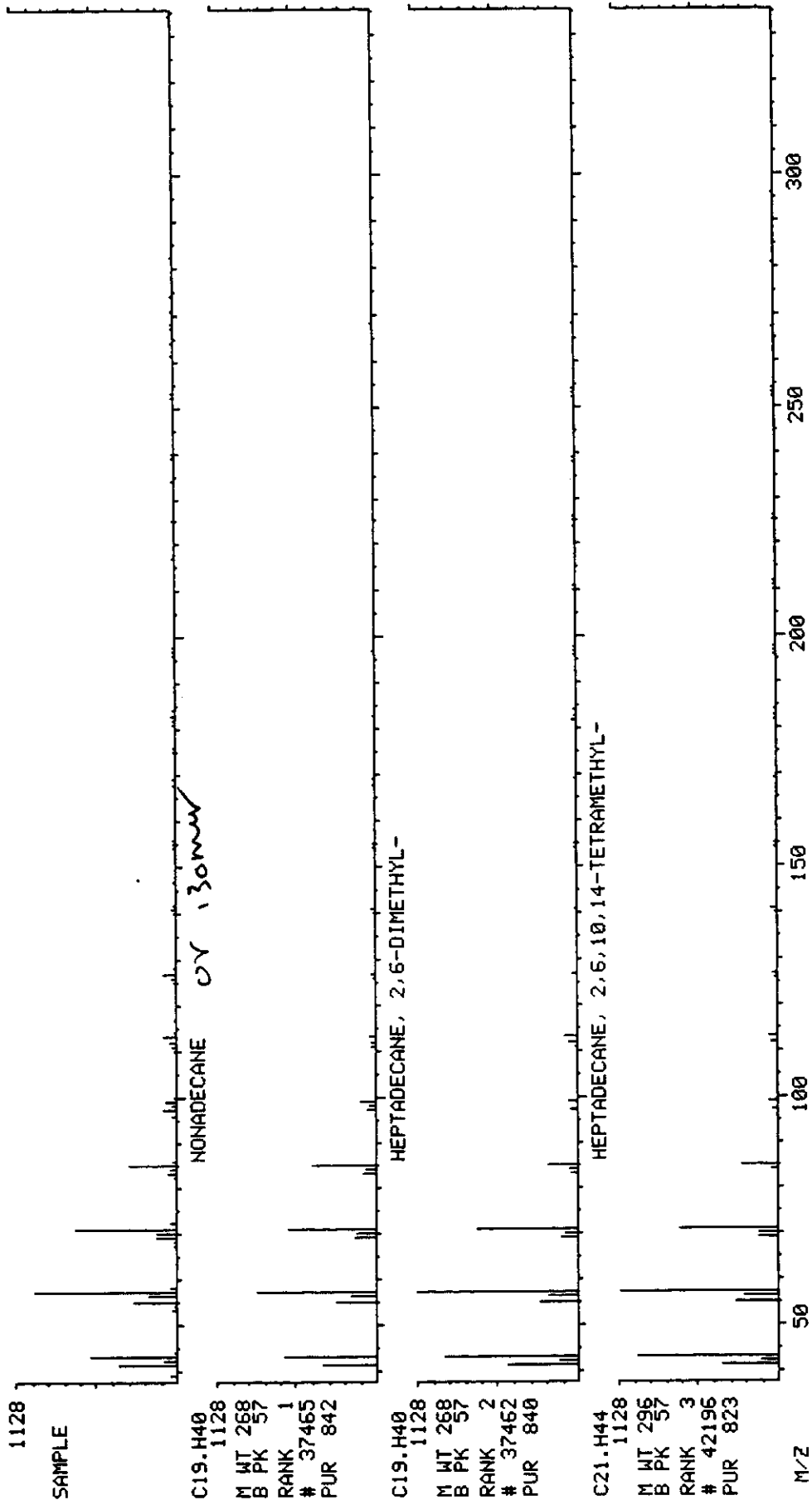
62231 spectra in LIBRARYNB searched for maximum PURITY
 840 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 46161 TRICOSANE
 5 32059 HEPTADECANE
 6 25991 DODECANE, 2,6,10-TRIMETHYL-
 7 26001 DODECANE, 2,7,10-TRIMETHYL-
 8 42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-
 9 37466 PENTADECANE, 2,6,10,14-TETRAMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	842	964	842
2	C19.H40	268	57	840	953	844
3	C21.H44	296	57	823	943	835
4	C23.H48	324	43	819	924	856
5	C17.H36	240	57	809	940	816
6	C15.H32	212	57	807	973	808
7	C15.H32	212	57	803	972	803
8	C21.H44	296	57	802	953	828
9	C19.H40	268	71	800	935	807

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	18344-37-1
4	—	—	—	—	638-67-5
5	—	—	—	—	629-78-7
6	—	—	—	—	3891-98-3
7	—	—	—	—	74645-98-0
8	—	—	—	—	54833-48-6
9	—	—	—	—	1921-70-6

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 12:27
 SAMPLE: 5-MM5-1B 1/35A/100M
 CONDS.: UC/ML *100ML *100% *100% *(NA/NA) /1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068102 #1119
 CALI: 30068102 # 3
 BASE M/Z: 57
 RIC: 112896.



Library Search Data: 30068102 #1191 Base m/z: 57
 08/31/98 19:31:00 + 13:15 Cali: 30068102 # 3 RIC: 71424.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

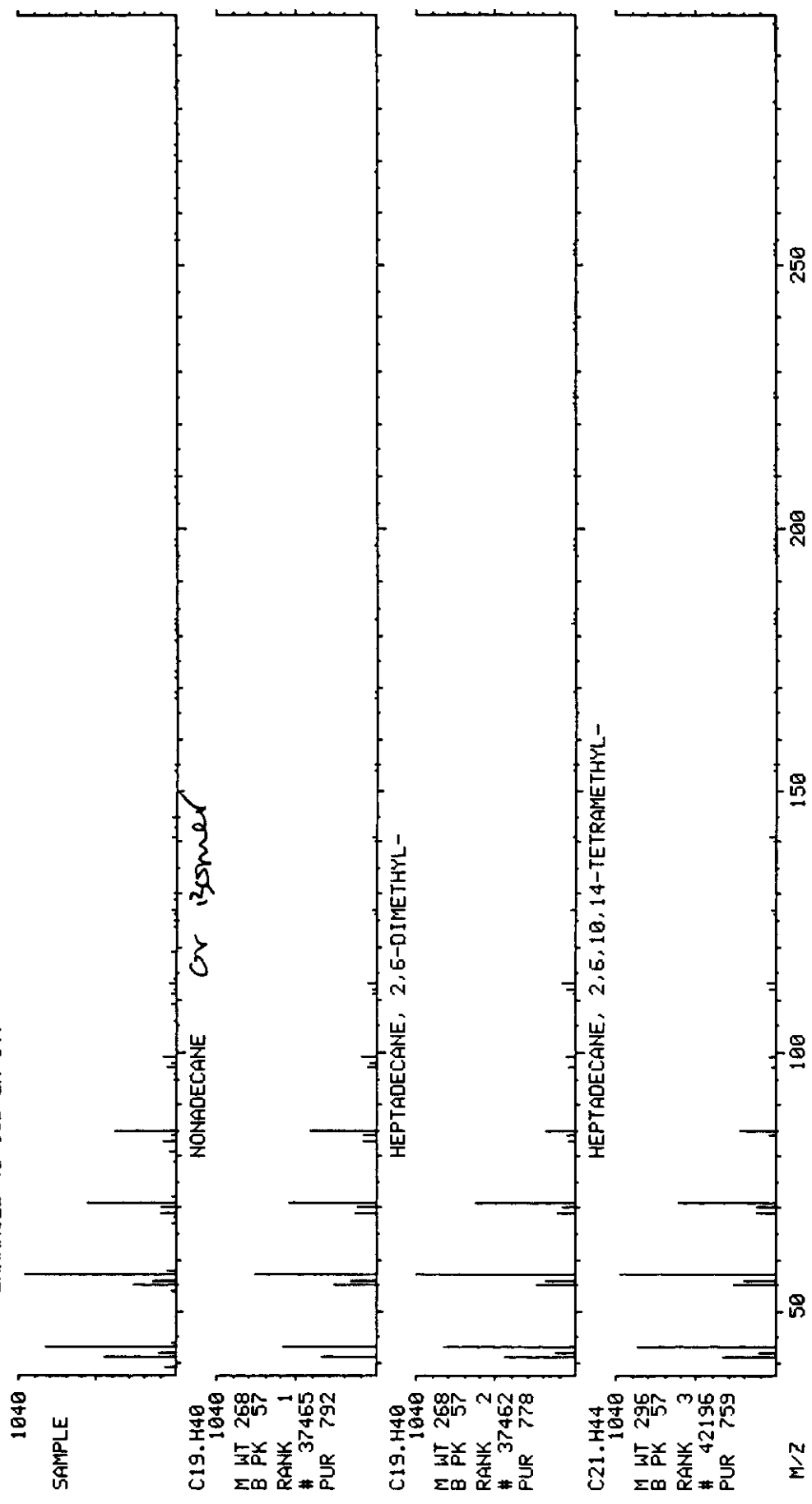
62231 spectra in LIBRARYNB searched for maximum PURITY
 879 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 46161 TRICOSANE
 5 32058 HEXADECANE, 3-METHYL-
 6 22530 TETRADECANE
 7 22536 TRIDECANE, 3-METHYL-
 8 25997 PENTADECANE
 9 42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	Rfit
1	C19.H40	268	57	792	971	798
2	C19.H40	268	57	778	942	788
3	C21.H44	296	57	759	955	773
4	C23.H48	324	43	749	911	808
5	C17.H36	240	57	748	932	758
6	C14.H30	198	43	743	948	752
7	C14.H30	198	57	740	928	748
8	C15.H32	212	57	736	948	764
9	C21.H44	296	57	736	933	764

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	18344-37-1
4	—	—	—	—	638-67-5
5	—	—	—	—	6418-43-5
6	—	—	—	—	629-59-4
7	—	—	—	—	6418-41-3
8	—	—	—	—	629-62-9
9	—	—	—	—	54833-48-6

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 19:31:00 + 13:15
 SAMPLE: S-MM5-1B 1/35A/100M
 COND5.: UG/ML *100ML *100Z/100Z *(NA/NA >1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30068102 #1191
 CALI: 30068102 # 3
 BASE M/Z: 57
 RIC: 71424.



Library Search Data: 30068102 #1265 Base m/z: 57
 08/31/98 19:31:00 + 14:04 Cali: 30068102 # 3 RIC: 53184.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 604 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-
 3 25997 PENTADECANE
 4 37465 NONADECANE
 5 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 6 25994 DODECANE, 2,6,11-TRIMETHYL-
 7 26001 DODECANE, 2,7,10-TRIMETHYL-
 8 32058 HEXADECANE, 3-METHYL-
 9 22535 DODECANE, 4,6-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	763	924	790
2	C21.H44	296	57	759	923	793
3	C15.H32	212	57	749	949	771
4	C19.H40	268	57	747	919	796
5	C21.H44	296	57	746	926	779
6	C15.H32	212	57	745	960	750
7	C15.H32	212	57	745	959	749
8	C17.H36	240	57	739	917	765
9	C14.H30	198	57	736	939	739

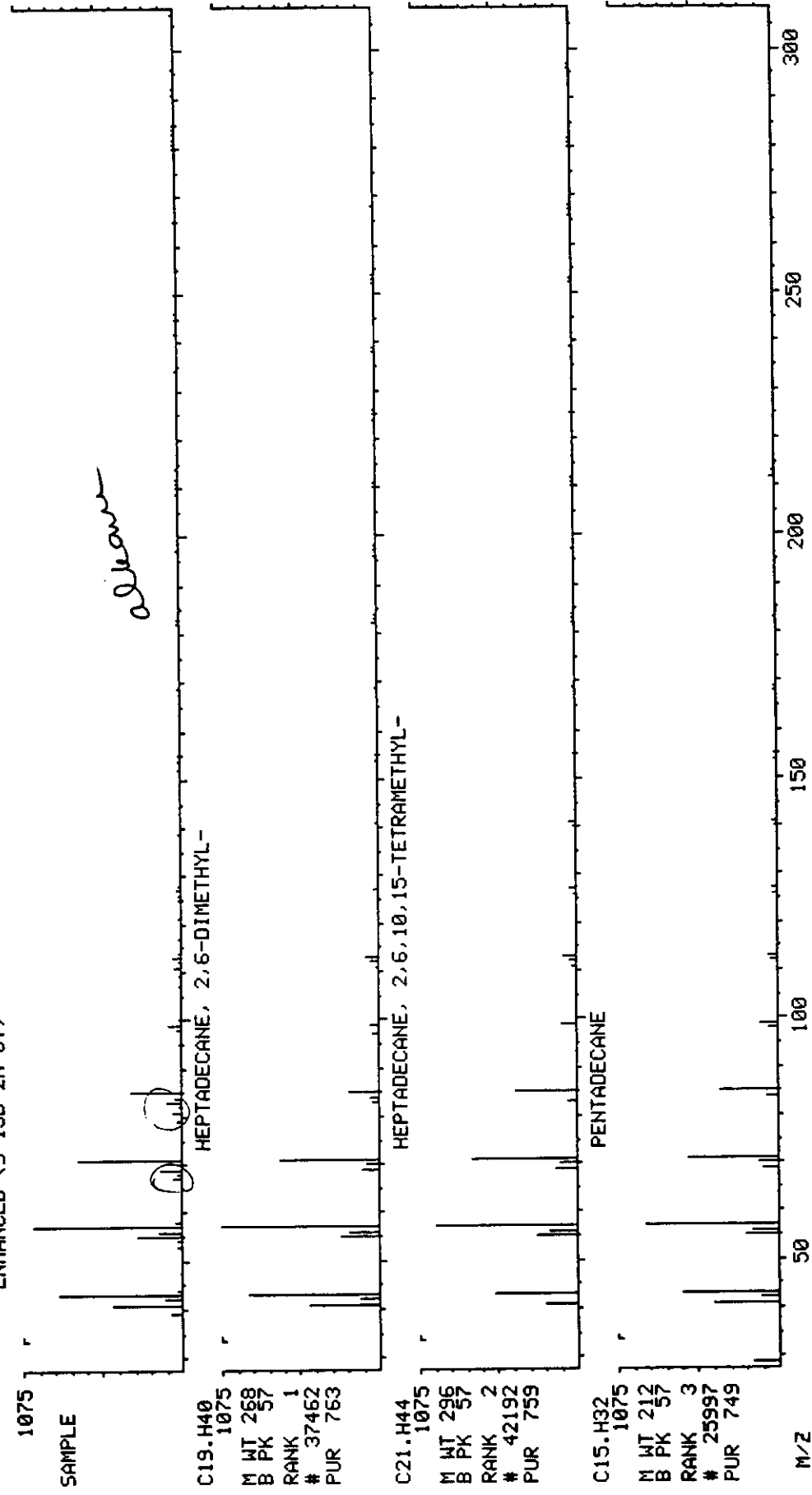
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	54833-48-6
3	—	—	—	—	629-62-9
4	—	—	—	—	629-92-5
5	—	—	—	—	18344-37-1
6	—	—	—	—	31295-56-4
7	—	—	—	—	74645-98-0
8	—	—	—	—	6418-43-5
9	—	—	—	—	61141-72-8

BASE M/Z: 57
RIC: 53184.

DATA: 30068102 #1265
CALI: 30068102 # 3

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 19:31:00 + 14:04
SAMPLE: S-MMS-1B 1/35A/100M
CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA) 1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068102 #1335 Base m/z: 57
 08/31/98 19:31:00 + 14:51 Cali: 30068102 # 3 RIC: 40704.
 Sample: S-MM5-1B 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 715 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37456 2-METHYLOCTADECANE
 2 37465 NONADECANE
 3 50998 HENEICOSANE, 11-(1-ETHYLPROPYL)-
 4 37462 HEPTADECANE, 2,6-DIMETHYL-
 5 32058 HEXADECANE, 3-METHYL-
 6 49555 PENTACOSANE
 7 46161 TRICOSANE
 8 29263 HEXADECANE
 9 34808 HEPTADECANE, 3-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	43	740	933	758
2	C19.H40	268	57	712	942	748
3	C26.H54	366	43	709	885	738
4	C19.H40	268	57	708	872	755
5	C17.H36	240	57	707	928	723
6	C25.H52	352	43	703	889	782
7	C23.H48	324	43	701	901	770
8	C16.H34	226	57	697	919	737
9	C18.H38	254	57	695	894	714

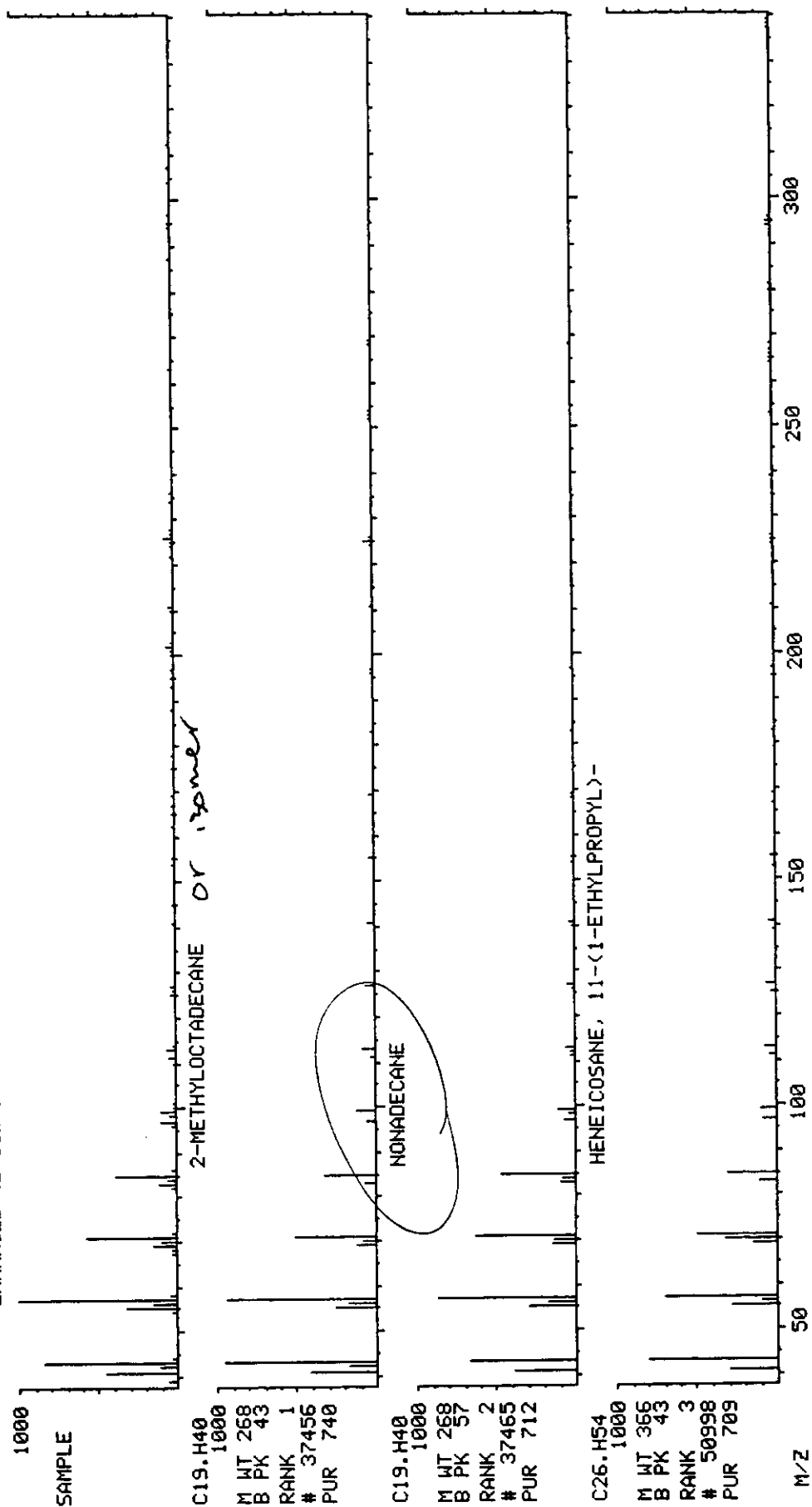
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	55282-11-6
3	—	—	—	—	54105-67-8
4	—	—	—	—	6418-43-5
5	—	—	—	—	629-99-2
6	—	—	—	—	638-67-5
7	—	—	—	—	544-76-3
8	—	—	—	—	6418-44-6
9	—	—	—	—	

BASE M/Z: 57
RIC: 40704.

DATA: 30058102 #1335
CALI: 30058102 # 3

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 19:31:00 + 14:51
SAMPLE: S-MM5-1B 1/35A/100M
CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA) 1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



801

Library Search Data: 30068102 #1402 Base m/z: 57
 08/31/98 19:31:00 + 15:36 Cali: 30068102 # 3 RIC: 29856.
 Sample: S-MM5-18 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 714 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 46161 TRICOSANE
 4 49555 PENTACOSANE
 5 32059 HEPTADECANE
 6 42193 EICOSANE, 10-METHYL-
 7 37456 2-METHYLOCTADECANE
 8 44314 DOCOSANE
 9 42197 HENEICOSANE

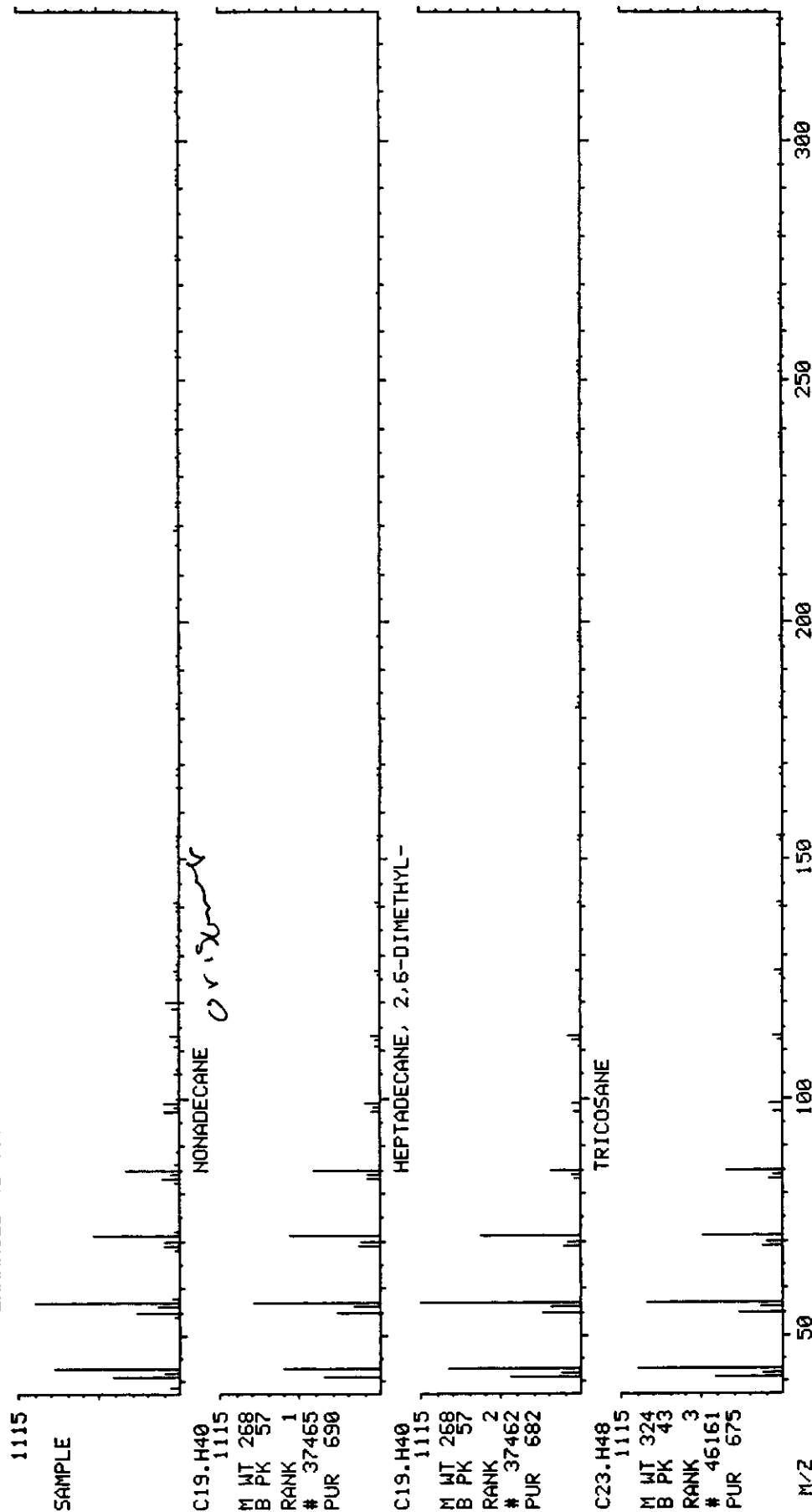
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	690	965	706
2	C19.H40	268	57	682	938	692
3	C23.H48	324	43	675	923	727
4	C25.H52	352	43	664	898	734
5	C17.H36	240	57	660	954	671
6	C21.H44	296	57	658	914	679
7	C19.H40	268	43	658	948	677
8	C22.H46	310	57	656	912	701
9	C21.H44	296	57	654	912	692

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	638-67-5
4	—	—	—	—	629-99-2
5	—	—	—	—	629-78-7
6	—	—	—	—	54833-23-7
7	—	—	—	—	- -
8	—	—	—	—	629-97-0
9	—	—	—	—	629-94-7

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 19:31:00 + 15:36
SAMPLE: S-MM5-1B 1/35A/100M
CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA) >1/35A NA M
ENHANCED (S 15B 2N 0T)

DATA: 30068102 #1402
CALI: 30068102 # 3

BASE M/Z: 57
RIC: 29856.



CC
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C

TIC SELECTION REPORT

DATA FILE: 30068102

THE FOLLOWING PEAKS WERE REJECTED BECAUSE
AT LEAST 40 % OF THEIR SIZE WAS ACCOUNTED FOR BY
TARGET COMPOUNDS ELUTING WITHIN 4 SCANS OF THE
PEAK TOP.

SCAN	SIZE	AMOUNT
341	188320.	24.846
530	398528.	52.579
829	303185.	40.000
1088	341760.	40.000
1552	204287.	40.000
1783	192640.	40.000

TOTAL NUMBER OF UNIDENTIFIED PEAKS WITH SIZE
GREATER THAN 10 % OF THE CLOSEST INTERNAL STANDARD
THAT DOES NOT HAVE INTERFERENCES = 95

INTERNAL STANDARDS THAT HAVE RIC SIZE LESS THAN
50 % OR GREATER THAN 200 % OF THE ESTIMATED RIC SIZE
ARE CONSIDERED TO HAVE INTERFERENCES AND WILL NOT BE USED
FOR QUANTITATION.

#	INTERNAL STANDARD	RIC SIZE	PERCENT OF ESTIMATED RIC SIZE	SCAN
1	C140 NAPHTHALENE-D8	398528.	260 *	530
2	C150 ACENAPHTHENE-D10	303185.	156	829
3	C160 PHENANTHRENE-D10	341760.	157	1088
4	C170 CHRYSENE-D12	204287.	145	1552
5	C175 PERYLENE-D12	192640.	162	1783

* INDICATES INTERFERENCE

SIZE = AREA

Semivolatile Organics
Method 0010/8270

Client Name: Pacific Environmental Services
Client ID: T-MM5-2-F, FH, XAD, COND, BH
LAB ID: 300681-0003-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 25 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
Phenol	ND	ug/Sample	30	R
bis(2-Chloroethyl)ether	ND	ug/Sample	30	
2-Chlorophenol	ND	ug/Sample	30	
1,3-Dichlorobenzene	ND	ug/Sample	30	
1,4-Dichlorobenzene	ND	ug/Sample	30	
Benzyl alcohol	ND	ug/Sample	30	
1,2-Dichlorobenzene	ND	ug/Sample	30	
2-Methylphenol	ND	ug/Sample	30	
2,2'-Oxybis(1-chloropropane)	ND	ug/Sample	30	
3/4-Methylphenol	ND	ug/Sample	30	
N-Nitroso-di-n-propylamine	ND	ug/Sample	30	
Hexachloroethane	ND	ug/Sample	30	
Nitrobenzene	ND	ug/Sample	30	
Isophorone	ND	ug/Sample	30	
2-Nitrophenol	ND	ug/Sample	30	
2,4-Dimethylphenol	ND	ug/Sample	30	
Benzoic acid	ND	ug/Sample	150	
bis(2-Chloroethoxy)-methane	ND	ug/Sample	30	
2,4-Dichlorophenol	ND	ug/Sample	30	
1,2,4-Trichlorobenzene	ND	ug/Sample	30	
Naphthalene	ND	ug/Sample	30	
4-Chloroaniline	ND	ug/Sample	30	
Hexachlorobutadiene	ND	ug/Sample	30	
4-Chloro-3-methylphenol	ND	ug/Sample	30	
2-Methylnaphthalene	60	ug/Sample	30	
Hexachlorocyclopentadiene	ND	ug/Sample	30	
2,4,6-Trichlorophenol	ND	ug/Sample	30	
2,4,5-Trichlorophenol	ND	ug/Sample	150	
2-Chloronaphthalene	ND	ug/Sample	30	
2-Nitroaniline	ND	ug/Sample	30	
Dimethyl phthalate	ND	ug/Sample	30	
Acenaphthylene	ND	ug/Sample	30	
3-Nitroaniline	ND	ug/Sample	150	
Acenaphthene	ND	ug/Sample	30	
2,4-Dinitrophenol	ND	ug/Sample	150	
4-Nitrophenol	ND	ug/Sample	150	
Dibenzofuran	ND	ug/Sample	30	

Note R = Reporting limit(s) raised due to sample volume limitations.
ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.
Rev 230787

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Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: T-MM5-2-F, FH, XAD, COND, BH
LAB ID: 300681-0003-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 25 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
2,4-Dinitrotoluene	ND	ug/Sample	30	
2,6-Dinitrotoluene	ND	ug/Sample	30	
Diethyl phthalate	ND	ug/Sample	30	
4-Chlorophenyl phenyl ether	ND	ug/Sample	30	
Fluorene	ND	ug/Sample	30	
4-Nitroaniline	ND	ug/Sample	150	
4,6-Dinitro-2-methylphenol	ND	ug/Sample	150	
N-Nitrosodiphenylamine	ND	ug/Sample	30	
4-Bromophenyl phenyl ether	ND	ug/Sample	30	
Hexachlorobenzene	ND	ug/Sample	30	
Pentachlorophenol	ND	ug/Sample	150	
Phenanthrene	ND	ug/Sample	30	
Anthracene	ND	ug/Sample	30	
Di-n-butyl phthalate	ND	ug/Sample	30	
Fluoranthene	ND	ug/Sample	30	
Pyrene	ND	ug/Sample	30	
Butyl benzyl phthalate	ND	ug/Sample	30	
3,3'-Dichlorobenzidine	ND	ug/Sample	60	
Benzo(a)anthracene	ND	ug/Sample	30	
bis(2-Ethylhexyl)-phthalate	ND	ug/Sample	30	
Chrysene	ND	ug/Sample	30	
Di-n-octyl phthalate	ND	ug/Sample	30	
Benzo(b)fluoranthene	ND	ug/Sample	30	
Benzo(k)fluoranthene	ND	ug/Sample	30	
Benzo(a)pyrene	ND	ug/Sample	30	
Indeno(1,2,3-cd)pyrene	ND	ug/Sample	30	
Dibenz(a,h)anthracene	ND	ug/Sample	30	
Benzo(g,h,i)perylene	ND	ug/Sample	30	
Acetophenone	ND	ug/Sample	30	
4-Aminobiphenyl	ND	ug/Sample	150	
Aniline	ND	ug/Sample	30	
Benzidine	ND	ug/Sample	300	
3,3'-Dimethylbenzidine	ND	ug/Sample	60	
N-Nitrosodimethylamine	ND	ug/Sample	30	
N-Nitrosomorpholine	ND	ug/Sample	30	
Pentachloronitrobenzene (PCNB)	ND	ug/Sample	150	
o-Toluidine	ND	ug/Sample	60	

ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

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The cover letter is an integral part of this report.
Rev 230787

Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: T-MM5-2-F, FH, XAD, COND, BH
LAB ID: 300681-0003-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 25 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
2-Methoxybenzenamine	ND	ug/Sample	--	K
Biphenyl	ND	ug/Sample	--	K
Chloroacetophenone	ND	ug/Sample	--	K
Cumene	ND	ug/Sample	--	K
DBCP (1,2-Dibromo-3-chloropropane)	ND	ug/Sample	--	K
Benzo(e)pyrene	ND	ug/Sample	--	K
N-N-Diethylaniline	ND	ug/Sample	--	K
Dimethylaniline	ND	ug/Sample	--	K
3,3'-Dimethoxybenzidine	ND	ug/Sample	--	K
Hydroquinone	ND	ug/Sample	--	K
4,4'-Methyl-bis(2-chloroaniline)	ND	ug/Sample	--	K
4-Nitrodiphenyl	ND	ug/Sample	--	K
Trifluralin	ND	ug/Sample	--	K

Surrogate	Recovery	Acceptable Range	
Nitrobenzene-d5	87 %	45 - 107	
2-Fluorobiphenyl	105 %	62 - 110	
Terphenyl-d14	78 %	58 - 135	
Phenol-d5	67 %	43 - 130	
2-Fluorophenol	62 %	36 - 111	
2,4,6-Tribromophenol	49 %	58 - 131	i

Note i = Surrogate recovery is outside of control limits.

Note K = Identified by mass spectrum only; quantitation based on 1:1 response with internal standard.

ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

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Semivolatiles Library Search (20 Compound TID)
Method 8270

Client Name: Pacific Environmental Services
Client ID: T-MM5-2-F, FH, XAD, COND, BH
LAB ID: 300681-0003-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 25 JUL 98
Prepared: NA

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Result	Units	Reporting Limit	Qualifier
4-Hydroxy-4-methyl-2-pentanone	450	ug/Sample	--	0
n-Nonane	370	ug/Sample	--	0
Unknown alkane	180	ug/Sample	--	
1H-Indene, 1-ethylidene-	240	ug/Sample	--	0
Undecane	180	ug/Sample	--	0
Heptadecane, 2,6,10,14 -tetramethyl-	220	ug/Sample	--	0
Unknown alkane	320	ug/Sample	--	
Unknown	200	ug/Sample	--	
Heptadecane, 2,6,10,14 -tetramethyl-	260	ug/Sample	--	0
Pentadecane	320	ug/Sample	--	0
Nonadecane	290	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	190	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	460	ug/Sample	--	0
Unknown alkane	160	ug/Sample	--	
Nonadecane	200	ug/Sample	--	0
Nonadecane	110	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	130	ug/Sample	--	0
Nonadecane	110	ug/Sample	--	0
3-Octadecene, (E)-	160	ug/Sample	--	0
Unknown	120	ug/Sample	--	

Note 0 = Or structurally similar compound (isomer).
NA = Not Applicable

Reported By: Emily Uebelhoer

Approved By: Karin Yee

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The cover letter is an integral part of this report.
Rev 230787

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068103
Std Id: ST16980831

Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 20:01
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Surrogate Spike Recoveries
8270-G, LIMS 10/10/96

Surrogate	Surrogate Amount (ug)		% Recovery	
	Spiked	Measured	Measured	QC limits
CS20 NITROBENZENE-D5	50.00	43.50	87.0	45 107
CS25 2-FLUOROBIPHENY	50.00	52.50	105.	62 110
CS30 TERPHENYL-D14	50.00	38.80	77.6	58 135
CS45 PHENOL-D5	100.0	67.20	67.2	43 130
CS50 2-FLUOROPHENOL	100.0	61.90	61.9	36 111
CS55 2,4,6-TRIBROMOP	100.0	48.90	48.9	* 58 131

RI confirms

29/10/98

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
HEXACHLOROBENZENE-C13		ND	UG/A	30.0
C310 N-NITROSODIMETHYLAMINE		ND	UG/A	30.0
PYRIDINE		ND	UG/A	60.0
2-PICOLINE		ND	UG/A	30.0
N-NITROSOMETHYLETHYLAMINE		ND	UG/A	30.0
METHYLMETHANESULFONATE		ND	UG/A	30.0
N-NITROSODIETHYLAMINE		ND	UG/A	30.0
ETHYLMETHANE SULFONATE		ND	UG/A	30.0
PENTACHLOROETHANE		ND	UG/A	30.0
C320 ANILINE		ND	UG/A	30.0
C315 PHENOL		ND	UG/A	30.0
C325 BIS(2-CHLOROETHYL)ETHER		ND	UG/A	30.0
C330 2-CHLOROPHENOL		ND	UG/A	30.0
C335 1,3-DICHLOROBENZENE		ND	UG/A	30.0
C340 1,4-DICHLOROBENZENE		ND	UG/A	30.0
C345 BENZYL ALCOHOL		ND	UG/A	30.0
C350 1,2-DICHLOROBENZENE		ND	UG/A	30.0
C355 2-METHYLPHENOL		ND	UG/A	30.0
C360 2,2'-OXYBIS(1-CLPROPAN)		ND	UG/A	30.0
C361 ACETOPHENONE		ND	UG/A	30.0
N-NITROSOPYRROLIDINE		ND	UG/A	30.0
N-NITROSOMORPHOLINE		ND	UG/A	30.0
3-METHYL PHENOL		ND	UG/A	30.0
C365 4-METHYLPHENOL		ND	UG/A	30.0
C370 N-NITROSO-DI-N-PROPYLAM		ND	UG/A	30.0
O-TOLUIDINE		ND	UG/A	30.0
C375 HEXACHLOROETHANE		ND	UG/A	30.0
C410 NITROBENZENE		ND	UG/A	30.0
N-NITROSOPIPERIDINE		ND	UG/A	30.0

Reviewed by:

8/9/17/96

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QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068103
Std Id: ST16980831

Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 20:01
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C415 ISOPHORONE		ND	UG/A	30.0
C420 2-NITROPHENOL		ND	UG/A	30.0
C425 2,4-DIMETHYLPHENOL		ND	UG/A	30.0
C435 BIS(2-CHLOROETHOXY)METH		ND	UG/A	30.0
C440 2,4-DICHLOROPHENOL		ND	UG/A	30.0
C430 BENZOIC ACID		ND	UG/A	150.0
AA-DIMETHYLPHENETHYLAMINE		ND	UG/A	30.0
C445 1,2,4-TRICHLOROBENZENE		ND	UG/A	30.0
C450 NAPHTHALENE		ND	UG/A	30.0
C455 4-CHLOROANALINE		ND	UG/A	30.0
2,6-DICHLOROPHENOL		ND	UG/A	30.0
HEXACHLOROPROPENE		ND	UG/A	30.0
C460 HEXACHLOROBUTADIENE		ND	UG/A	30.0
P-PHENYLENE DIAMINE		ND	UG/A	30.0
N-NITROSODI-N-BUTYLAMINE		ND	UG/A	30.0
C465 4-CHLORO-3-METHYLPHENO		ND	UG/A	30.0
SAFROLE		ND	UG/A	30.0
C470 2-METHYLNAPHTHALENE	653	60	UG/A	30.0
1,2,4,5-TETRACHLOROBENZENE		ND	UG/A	30.0
ISOSAFROLE (#1)		ND	UG/A	60.0
C510 HEXACHLOROCYCLOPENTADI		ND	UG/A	30.0
C515 2,4,6-TRICHLOROPHENOL		ND	UG/A	30.0
C520 2,4,5-TRICHLOROPHENOL		ND	UG/A	30.0
ISOSAFROLE (#2)		ND	UG/A	60.0
C525 2-CHLORONAPHTHALENE		ND	UG/A	30.0
1-CHLORONAPHTHALENE		ND	UG/A	15.0
C530 2-NITROANALINE		ND	UG/A	150.0
1,4-NAPHTHOQUINONE		ND	UG/A	30.0
C535 DIMETHYLPHTHALATE		ND	UG/A	30.0
1,3-DINITROBENZENE		ND	UG/A	30.0
C540 ACENAPHTHYLENE		ND	UG/A	30.0
C543 2,6-DINITROTOLUENE		ND	UG/A	30.0
C545 3-NITROANILINE		ND	UG/A	150.0
C550 ACENAPHTHENE		ND	UG/A	30.0
C555 2,4-DINITROPHENOL		ND	UG/A	150.0
C565 DIBENZOFURAN		ND	UG/A	30.0
C560 4-NITROPHENOL		ND	UG/A	150.0
PENTACHLOROBENZENE		ND	UG/A	30.0
C570 2,4-DINITROTOLUENE		ND	UG/A	30.0
1-NAPHTHYLAMINE		ND	UG/A	30.0
2-NAPHTHYLAMINE		ND	UG/A	30.0
2,3,4,6-TETRACHLOROPHENOL		ND	UG/A	60.0
C580 DIETHYLPHTHALATE		ND	UG/A	30.0
C590 FLUORENE		ND	UG/A	30.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068103
Std Id: ST16980831

Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 20:01
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
=====				
C585 4-CHLOROPHENYL-PHENYLE		ND	UG/A	30.0
5-NITRO-O-TOLUIDINE		ND	UG/A	30.0
C595 4-NITROANALINE		ND	UG/A	150.0
C610 4,6-DINITRO-2-METHYLPH		ND	UG/A	150.0
C615 N-NITROSODIPHENYLAMINE		ND	UG/A	30.0
C620 AZOBENZENE		ND	UG/A	30.0
SYM-TRINITROBENZENE		ND	UG/A	30.0
C625 4-BROMOPHENYL-PHENYLET		ND	UG/A	30.0
PHENACETIN		ND	UG/A	30.0
DIALATE	234	ND	UG/A	30.0
C630 HEXACHLOROBENZENE		ND	UG/A	30.0
4-AMINOBIIPHENYL		ND	UG/A	30.0
C635 PENTACHLOROPHENOL		ND	UG/A	150.0
PRONAMIDE		ND	UG/A	30.0
PENTACHLORONITROBENZENE		ND	UG/A	150.0
C640 PHENANTHRENE		ND	UG/A	30.0
C645 ANTHRACENE		ND	UG/A	30.0
2SECBUTYL-4,6-DINITROPHENOL		ND	UG/A	30.0
C647 CARBAZOLE		ND	UG/A	30.0
C650 DI-N-BUTYLPHTHALATE		ND	UG/A	30.0
4-NITROQUINOLINE-1-OXIDE		ND	UG/A	30.0
METHAPYRILENE		ND	UG/A	30.0
ISODRIN		ND	UG/A	30.0
C655 FLUORANTHENE		ND	UG/A	30.0
CHLOROBENZILATE		ND	UG/A	30.0
C710 BENZIDINE		ND	UG/A	300.0
C715 PYRENE		ND	UG/A	30.0
ARAMITE (#1)		ND	UG/A	30.0
ARAMITE (#2)		ND	UG/A	30.0
P-DIMETHYLAMINOAZOBENZENE		ND	UG/A	30.0
3,3'-DIMETHYLBENZIDINE		ND	UG/A	30.0
KEPONE		ND	UG/A	150.0
C720 BUTYLBENZYLPHTHALATE		ND	UG/A	30.0
2-ACETYLAMINOFLUORENE		ND	UG/A	30.0
C730 BENZO(A)ANTHRACENE		ND	UG/A	30.0
C725 3,3'-DICHLOROBENZIDINE		ND	UG/A	60.0
C740 CHRYSENE		ND	UG/A	30.0
C745 BIS(2-ETHYLHEXYL)PHTHA		ND	UG/A	30.0
3-METHYLCHOLANTHRENE		ND	UG/A	30.0
C760 DI-N-OCTYL PHTHALATE		ND	UG/A	30.0
C765 BENZO(B)FLUORANTHENE		ND	UG/A	30.0
7,12-DIMETHYLBENZANTHRACENE		ND	UG/A	30.0
C770 BENZO(K)FLUORANTHENE		ND	UG/A	30.0
HEXACHLOROPHENE		ND	UG/A	30.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

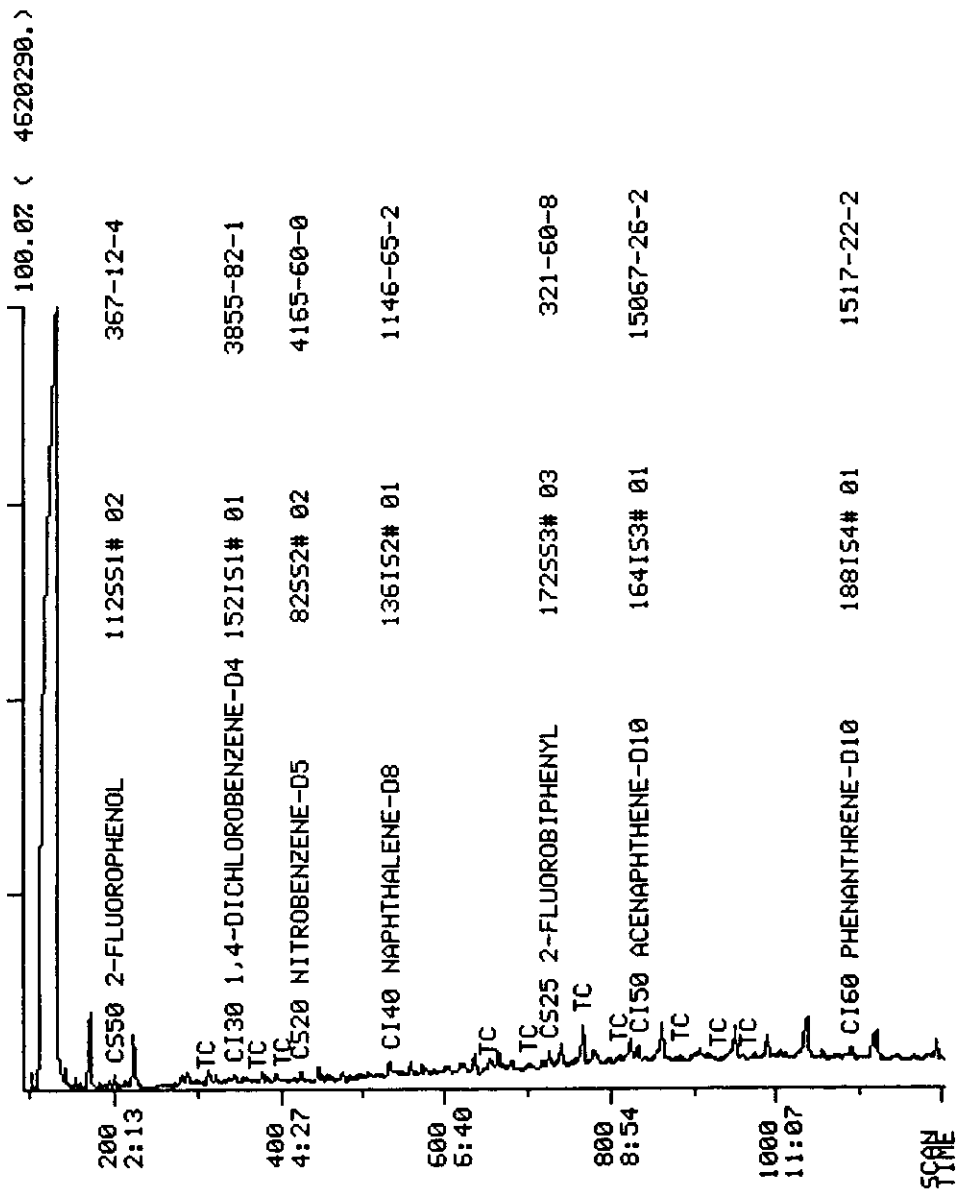
Data File: 30068103
Std Id: ST16980831

Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 20:01
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

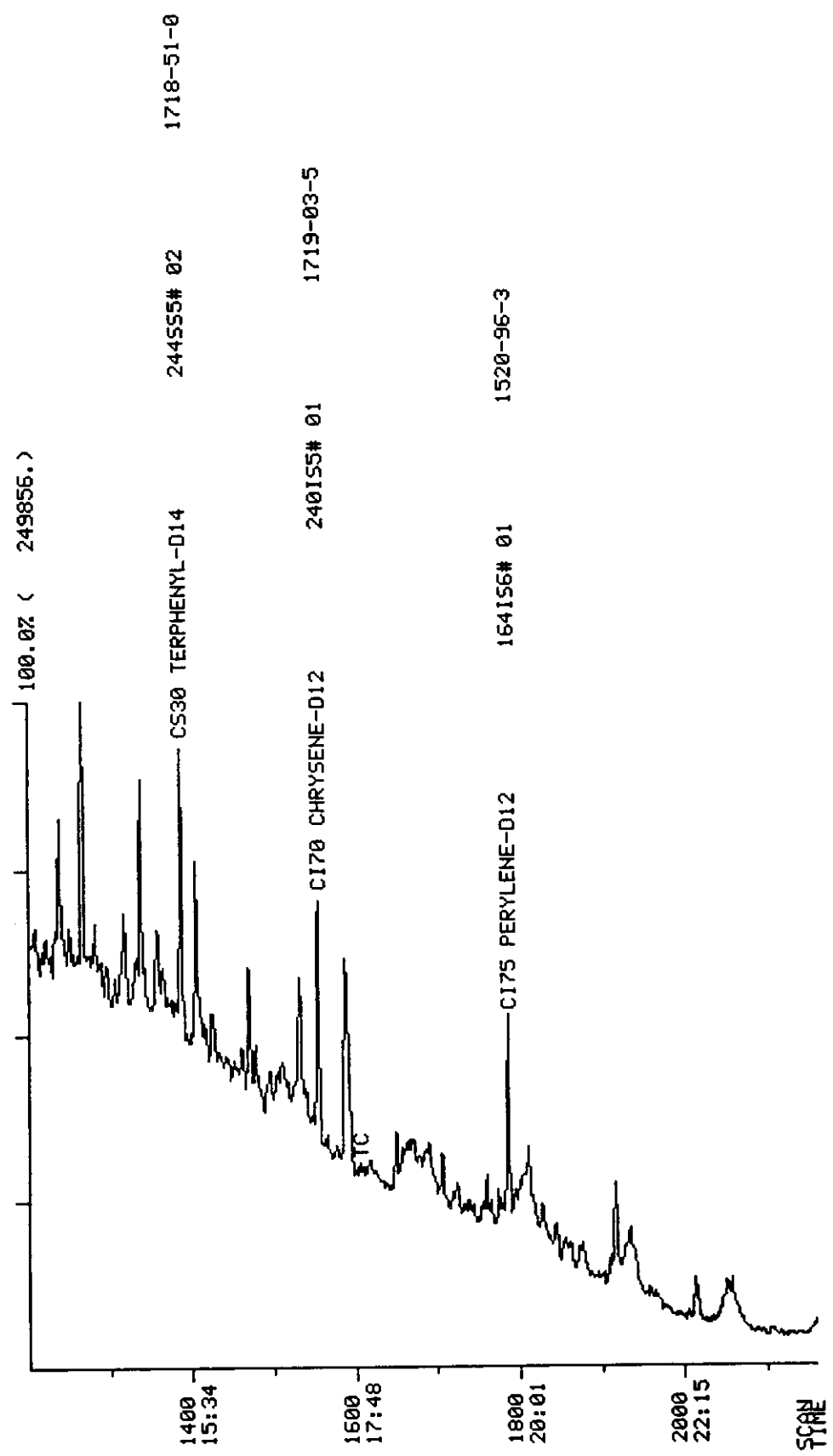
Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C775 BENZO(A)PYRENE		ND	UG/A	30.0
C780 INDENO(1,2,3-CD)PYRENE		ND	UG/A	30.0
C785 DIBENZ(A,H)ANTHRACENE		ND	UG/A	30.0
C790 BENZO(G,H,I)PERYLENE		ND	UG/A	30.0

DATA FROM FILE: 30068103 SCANS 93 TO 1203 ACQUIRED: 08/31/98 20:01:00
 CALI: 30068103 #3
 SAMPLE: T-MM5-2 1/35A/1ML INST. ID: F15
 CONDS.: UG/ML *1ML *100%/100% *CNA/NA >1/35A NA M



DATA FROM FILE: 30068103 SCANS 1203 TO 2158 ACQUIRED: 08/31/98 20:01:00
 CALI: 30068103 #3
 SAMPLE: T-MMS-2 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *1002/1002 *(NA/NA)/1/35A NA M



ADDITIONAL STANDARD USED FOR THIS REPORT IS: ST16980831A

QUANTERRA QUANTITATION SUMMARY

File: 30068103

Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Analyst: DAT Instrument Id: F16 Analyzed: 08/31/98 20:01

Compounds with amounts less than 0.20 reported as NOT FOUND

Library		Units: UG/ML					
No Entry	Name	Mass Meth	Scan	Ref	Area RRF(L)	Amount	
1 S1#	1 C130 1,4-DICHLOROBENZENE	152 A	BV	341 1	31920. 1.000	40.000	
2 S2#	1 C140 NAPHTHALENE-D8	136 A	BB	531 2	111468. 1.000	40.000	
3 S3#	1 C150 ACENAPHTHENE-D10	164 A	BB	830 3	58555. 1.000	40.000	
4 S4#	1 C160 PHENANTHRENE-D10	188 A	BB	1088 4	96270. 1.000	40.000	
5 S5#	1 C170 CHRYSENE-D12	240 A	BB	1554 5	92258. 1.000	40.000	
6 S6#	1 C175 PERYLENE-D12	264 A	BB	1784 6	85395. 1.000	40.000	
7 S2#	2 CS20 NITROBENZENE-D5	82 A	BB	422 2	56985. 0.470	43.550	
8 S3#	3 CS25 2-FLUOROBIPHENYL	172 A	BB	724 3	94839. 1.234	52.489	
9 S5#	2 CS30 TERPHENYL-D14	244 A	BB	1386 5	89070. 0.995	38.803	
10 S1#	3 CS45 PHENOL-D5	99 A	BB	310 1	108633. 2.025	67.210	
11 S1#	2 CS50 2-FLUOROPHENOL	112 A	BB	199 1	67610. 1.368	61.923	
12 S3#	2 CS55 2,4,6-TRIBROMOPHENO	330 A	BB	971 3	16861. 0.236	48.877	
13 S4#	4 HEXACHLOROBENZENE-C13	294		4 NOT FOUND			
14 S1#	15 C310 N-NITROSODIMETHYLAM	74		1 NOT FOUND			
15 S1#	20 PYRIDINE	79		1 NOT FOUND			
16 S1#	30 2-PICOLINE	93		1 NOT FOUND			
17 S1#	40 N-NITROSOMETHYLETHYLAMIN	42		1 NOT FOUND			
18 S1#	60 METHYLMETHANESULFONATE	80		1 NOT FOUND			
19 S1#	70 N-NITROSODIETHYLAMINE	102		1 NOT FOUND			
20 S1#	85 ETHYLMETHANE SULFONATE	79		1 NOT FOUND			
21 S1#	95 PENTACHLOROETHANE	117 A	BB	307 1	821. 0.602	1.710	
22 S1#	100 C320 ANILINE	93		1 NOT FOUND			
23 S1#	105 C315 PHENOL	94 A	BB	311 1	11667. 1.968	7.430	
24 S1#	110 C325 BIS(2-CHLOROETHYL)E	93 A	BB	320 1	634. 1.368	0.581	
25 S1#	115 C330 2-CHLOROPHENOL	128		1 NOT FOUND			
26 S1#	125 C335 1,3-DICHLOROBENZENE	146		1 NOT FOUND			
27 S1#	130 C340 1,4-DICHLOROBENZENE	146 A	BB	343 1	532. 1.617	0.412	
28 S1#	145 C345 BENZYL ALCOHOL	108 A	VB	370 1	5992. 0.854	8.789	
29 S1#	150 C350 1,2-DICHLOROBENZENE	146		1 NOT FOUND			
30 S1#	160 C355 2-METHYLPHENOL	108		1 NOT FOUND			
31 S1#	165 C360 2,2'-OXYBIS(1-CLPRO	45		1 NOT FOUND			
32 S1#	170 C361 ACETOPHENONE	105 A	BB	401 1	11312. 1.934	7.330	
33 S1#	175 N-NITROSOPYRROLIDINE	100 A	VB	402 1	224. 0.553	0.508	
34 S1#	180 N-NITROSOMORPHOLINE	56 A	VV	410 1	3898. 0.821	5.952	
35 S1#	182 3-METHYL PHENOL	108		1 NOT FOUND			
36 S1#	185 C365 4-METHYLPHENOL	108		1 NOT FOUND			
37 S1#	190 C370 N-NITROSO-DI-N-PROP	70 A	VB	404 1	1437. 0.906	1.988	
38 S1#	195 O-TOLUIDINE	106		1 NOT FOUND			
39 S1#	200 C375 HEXACHLOROETHANE	117		1 NOT FOUND			
40 S2#	10 C410 NITROBENZENE	77 A	BB	423 2	1534. 0.445	1.237	
41 S2#	15 N-NITROSOPIPERIDINE	42		2 NOT FOUND			
42 S2#	20 C415 ISOPHORONE	82		2 NOT FOUND			
43 S2#	25 C420 2-NITROPHENOL	139		2 NOT FOUND			
44 S2#	30 C425 2,4-DIMETHYLPHENOL	107 A	BB	500 2	820. 0.330	0.892	
45 S2#	35 C435 BIS(2-CHLOROETHOXY)	93		2 NOT FOUND			

46	S2# 40	C440 2,4-DICHLOROPHENOL	162 A BB	519 2	346. 0.308	0.404
47	S2# 45	C430 BENZOIC ACID	122	2 NOT FOUND		
48	S2# 50	AA-DIMETHYLPHENETHYLAMIN	58	2 NOT FOUND		
49	S2# 55	C445 1,2,4-TRICHLOROBENZ	180 A VV	527 2	220. 0.331	0.239

50	S2# 60	C450 NAPHTHALENE	128	A BV	534	2	27047.	1.022	9.495
51	S2# 80	C455 4-CHLOROANALINE	127	A BB	556	2	596.	0.454	0.471
52	S2# 85	2,6-DICHLOROPHENOL	162		2	NOT FOUND			
53	S2# 90	HEXACHLOROPROPENE	213		2	NOT FOUND			
54	S2# 95	C460 HEXACHLOROBUTADIENE	225		2	NOT FOUND			
55	S2#115	P-PHENYLENE DIAMINE	108		2	NOT FOUND			
56	S2#120	N-NITROSODI-N-BUTYLAMINE	84		2	NOT FOUND			
57	S2#130	C465 4-CHLORO-3-METHYLPH	107		2	NOT FOUND			
58	S2#140	SAFROLE	162	A BB	649	2	1150.	0.248	1.666
59	S2#145	C470 2-METHYLNAPHTHALENE	142	A BB	653	2	42032.	0.751	20.077
60	S3# 10	1,2,4,5-TETRACHLOROBENZE	214		3	NOT FOUND			
61	S3# 15	ISOSAFROLE (#1)	162	A BB	698	3	451.	0.044	7.066
62	S3# 20	C510 HEXACHLOROCYCLOPENT	237		3	NOT FOUND			
63	S3# 25	C515 2,4,6-TRICHLOROPHEN	196		3	NOT FOUND			
64	S3# 30	C520 2,4,5-TRICHLOROPHEN	196		3	NOT FOUND			
65	S3# 35	ISOSAFROLE (#2)	104		3	NOT FOUND			
66	S3# 40	C525 2-CHLORONAPHTHALENE	162		3	NOT FOUND			
67	S3# 42	1-CHLORONAPHTHALENE	162		3	NOT FOUND			
68	S3# 45	C530 2-NITROANALINE	65	A BB	765	3	848.	0.502	1.155
69	S3# 50	1,4-NAPHTHOQUINONE	158		3	NOT FOUND			
70	S3# 55	C535 DIMETHYLPHTHALATE	163		3	NOT FOUND			
71	S3# 60	1,3-DINITROBENZENE	168	A BB	808	3	398.	0.197	1.378
72	S3# 65	C540 ACENAPHTHYLENE	152		3	NOT FOUND			
73	S3# 70	C543 2,6-DINITROTOLUENE	165		3	NOT FOUND			
74	S3# 75	C545 3-NITROANILINE	138		3	NOT FOUND			
75	S3# 80	C550 ACENAPHTHENE	153	A VV	836	3	1939.	1.159	1.143
76	S3# 85	C555 2,4-DINITROPHENOL	184	A BB	858	3	158.	0.189	0.570
77	S3# 90	C565 DIBENZOFURAN	168	A VB	868	3	1797.	1.669	0.735
78	S3# 95	C560 4-NITROPHENOL	109		3	NOT FOUND			
79	S3#100	PENTACHLOROBENZENE	250		3	NOT FOUND			
80	S3#105	C570 2,4-DINITROTOLUENE	165	A VB	882	3	891.	0.397	1.532
81	S3#110	1-NAPHTHYLAMINE	143	A BV	882	3	394.	0.837	0.321
82	S3#115	2-NAPHTHYLAMINE	143	A BB	903	3	418.	0.966	0.296
83	S3#120	2,3,4,6-TETRACHLOROPHENO	232		3	NOT FOUND			
84	S3#130	C580 DIETHYLPHTHALATE	149		3	NOT FOUND			
85	S3#135	C590 FLUORENE	166	A BB	927	3	5465.	1.298	2.877
86	S3#140	C585 4-CHLOROPHENYL-PHEN	204		3	NOT FOUND			
87	S3#145	5-NITRO-O-TOLUIDINE	152	A VB	942	3	1762.	0.356	3.386
88	S3#150	C595 4-NITROANALINE	138	A BB	950	3	223.	0.307	0.495
89	S4# 10	C610 4,6-DINITRO-2-METHY	198		4	NOT FOUND			
90	S4# 15	C615 N-NITROSODIPHENYLAM	169	A BB	965	4	4015.	0.551	3.029
91	S4# 20	C620 AZOBENZENE	77		4	NOT FOUND			
92	S4# 25	SYM-TRINITROBENZENE	75	A BB	1026	4	132.	0.162	0.340
93	S4# 30	C625 4-BROMOPHENYL-PHENY	248		4	NOT FOUND			
94	S4# 35	PHENACETIN	108		4	NOT FOUND			
95	S4# 37	DIALATE	234		4	NOT FOUND			
96	S4# 40	C630 HEXACHLOROBENZENE	284		4	NOT FOUND			
97	S4# 45	4-AMINOBIIPHENYL	169		4	NOT FOUND			
98	S4# 50	C635 PENTACHLOROPHENOL	266		4	NOT FOUND			
99	S4# 55	PRONAMIDE	173		4	NOT FOUND			
100	S4# 60	PENTACHLORONITROBENZENE	237		4	NOT FOUND			
101	S4# 65	C640 PHENANTHRENE	178	A BB	1092	4	17905.	1.033	7.205
102	S4# 70	C645 ANTHRACENE	178		4	NOT FOUND			
103	S4# 75	2SECBUTYL-4,6-DINITROPHE	211		4	NOT FOUND			
104	S4# 80	C647 CARBAZOLE	167	A BB	1149	4	650.	0.764	0.354
105	S4# 85	C650 DI-N-BUTYLPHTHALATE	149		4	NOT FOUND			
106	S4#100	4-NITROQUINOLINE-1-OXIDE	190		4	NOT FOUND			
107	S4#105	METHAPYRILENE	58		4	NOT FOUND			
108	S4#106	ISODRIN	193		4	NOT FOUND			
109	S4#110	C655 FLUORANTHENE	202	A BB	1300	4	550.	0.900	0.254
110	S4#120	CHLOROBENZILATE	139		4	NOT FOUND			

111 S5# 10 C710 BENZIDINE	184	5	NOT FOUND		
112 S5# 15 C715 PYRENE	202 A BB 1336	5	2225.	1.323	0.729
113 S5# 20 ARAMITE (#1)	185	5	NOT FOUND		
114 S5# 25 ARAMITE (#2)	185	5	NOT FOUND		
115 S5# 30 P-DIMETHYLAMINOAZOBENZEN	120	5	NOT FOUND		
116 S5# 35 3,3'-DIMETHYLBENZIDINE	212	5	NOT FOUND		
117 S5# 37 KEPONE	272	5	NOT FOUND		
118 S5# 40 C720 BUTYLBENZYLPHTHALAT	149	5	NOT FOUND		
119 S5# 45 2-ACETYLAMINOFLUORENE	181	5	NOT FOUND		
120 S5# 50 C730 BENZO(A)ANTHRACENE	228 A BB 1557	5	1896.	1.125	0.731
121 S5# 55 C725 3,3'-DICHLOROBENZID	252	5	NOT FOUND		
122 S5# 60 C740 CHRYSENE	228 A BB 1557	5	1896.	1.014	0.811
123 S5# 65 C745 BIS(2-ETHYLNEXYL)PH	149 A BB 1604	5	3863.	0.980	1.710
124 S5# 85 3-METHYLCOLANTHRENE	268 A*BV 1832	5	454.	0.548	0.359
125 S6# 10 C760 DI-N-OCTYL PHTHALAT	149	6	NOT FOUND		
126 S6# 15 C765 BENZO(B)FLUORANTHEN	252	6	NOT FOUND		
127 S6# 20 7,12-DIMETHYLBENZANTHRAC	256 A BB 1742	6	376.	0.610	0.289
128 S6# 25 C770 BENZO(K)FLUORANTHEN	252	6	NOT FOUND		
129 S6# 30 HEXACHLOROPHENE	196	6	NOT FOUND		
130 S6# 35 C775 BENZO(A)PYRENE	252	6	NOT FOUND		
131 S6# 55 C780 INDENO(1,2,3-CD)PYR	276	6	NOT FOUND		
132 S6# 60 C785 DIBENZ(A,H)ANTHRACE	278	6	NOT FOUND		
133 S6# 65 C790 BENZO(G,H,I)PERYLEN	276	6	NOT FOUND		

BNA, S LIBRARIES

ANALYST: _____ DATE: _____

QUANTERRA Internal Standard Check

Standard Filename: ST16980831

Analyzed: 08/31/98 12:31

Standard	IS# 1			IS# 2			IS# 3		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	28649.		3.78	100432.		5.90	54347.		9.21
Upper Limit	57297.==+200		4.28	200864.==+200		6.40	108695.==+200		9.71
Lower Limit	14324.==+ 50		3.28	50216.==+ 50		5.40	27174.==+ 50		8.71
Filename									
1 30068103	31920.= 111		3.79	111468.= 110		5.91	58555.= 107		9.23

IS# 1 = C130 1,4-DICHLOROBENZENE-D4

IS# 2 = C140 NAPHTHALENE-D8

IS# 3 = C150 ACENAPHTHENE-D10

* - indicates an I.S. is outside QC limit(s)

Standard	IS# 4			IS# 5			IS# 6		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	98520.		12.08	62289.		17.26	50750.		19.83
Upper Limit	197039.==+200		12.58	124578.==+200		17.76	101500.==+200		20.33
Lower Limit	49260.==+ 50		11.58	31145.==+ 50		16.76	25375.==+ 50		19.33
Filename									
1 30068103	96270.= 97		12.10	92258.= 148		17.29	85395.= 168		19.84

IS# 4 = C160 PHENANTHRENE-D10

IS# 5 = C170 CHRYSENE-D12

IS# 6 = C175 PERYLENE-D12

* - indicates an I.S. is outside QC limit(s)

TARGET COMPOUND COMPARISON

COMPOUND: C470 2-METHYLNAPHTHALENE 142 S2#145 91-57-6

RAW DATA: 30068103 #653

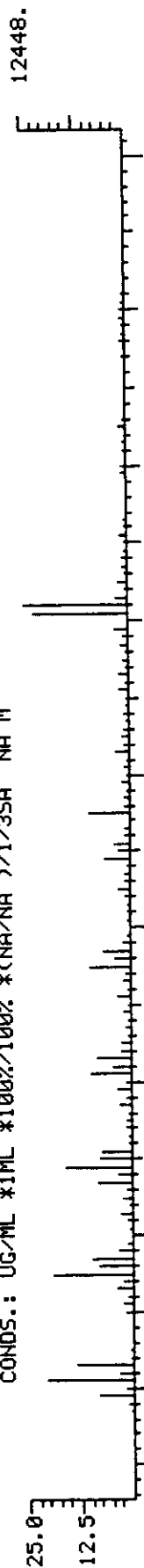
08/31/98 20:01

SAMPLE: T-MMS-2 1/35A/1ML

COND.: UG/ML *1ML *1002/1002 *(NA/NA) >1/35A NA M

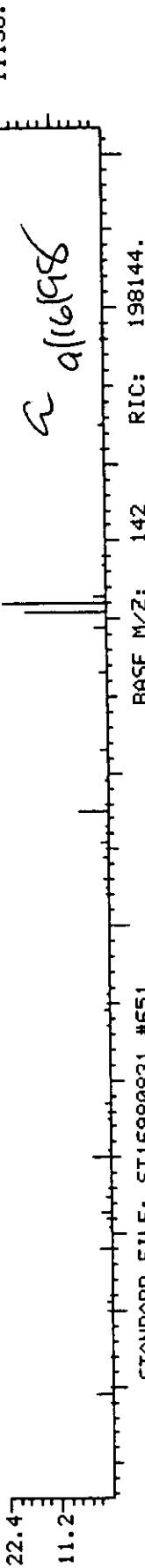
INST. ID: F16

BASE M/Z: 142 RIC: 179200.



ENHANCED DATA: 30068103 #653

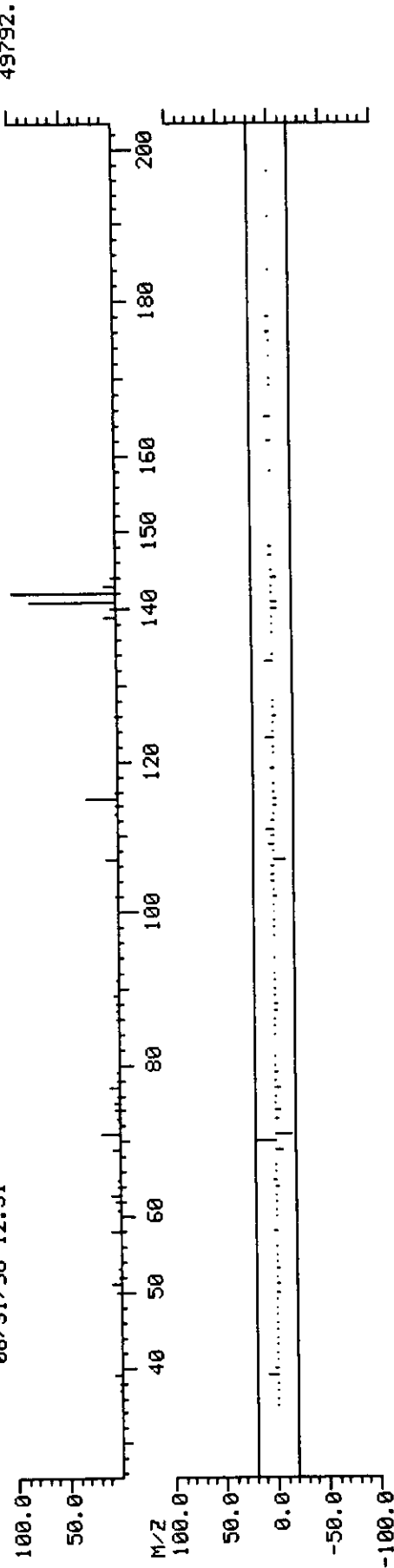
BASE M/Z: 142 RIC: 43904.



STANDARD FILE: ST16980831 #651

08/31/98 12:31

BASE M/Z: 142 RIC: 198144.



Data Reduced by: 2 Date: 9/17/98
Data Reviewed by: 17 Date: 9/17/98

Data File: 30068103

QUANTERRA GC/MS TIC REPORT (Part 1)

Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
Analyst: DAT Date Analyzed: 08/31/98 20:01
Run Factor: 3.00

#	SCAN	Concentration in Sample (UG/A)	CAS #	
1	171	450.	123-42-2	
				2-PENTANONE, 4-HYDROXY-4-METHYL- <i>or isomer</i>
2	222	370.	111-84-2	
				NONANE <i>or isomer</i>
3	444	110.	00-00-0	
				UNKNOWN
4	473	99.	71899-38-2	
				9-EICOSYNE
5	633	180.	54105-67-8	5746400
				HEPTADECANE, 2,6-DIMETHYL-
				<i>unkn. alkane</i>
6	653	240.	2471-83-2	
				1H-INDENE, 1-ETHYLIDENE- <i>or isomer</i>
7	663	180.	1120-21-4	
				UNDECANE <i>or isomer</i>
8	739	220.	18344-37-1	
				HEPTADECANE, 2,6,10,14-TETRAMETHYL- <i>or isomer</i>
9	765	320.	62108-21-8	5746400
				DECANE, 6-ETHYL-2-METHYL-
				<i>unkn. alkane</i>
10	778	200.	00-00-0	5700
				UNKNOWN

841

11 821 260. 18344-37-1
HEPTADECANE, 2,6,10,14-TETRAMETHYL- *or isomer*

12 859 320. 629-62-9
PENTADECANE *or isomer*

13 949 290. 629-92-5
NONADECANE *or isomer*

14 989 190. 54105-67-8
HEPTADECANE, 2,6-DIMETHYL- *or isomer*

15 1037 460. 54105-67-8
HEPTADECANE, 2,6-DIMETHYL- *or isomer*

16 1115 160. ~~29812-79-1~~ 5746400
~~HYDROXYLAMINE, 9-DECYL-~~

unknown alkane

17 1120 200. 629-92-5
NONADECANE *or isomer*

~~18~~ 1192 110. 629-92-5
NONADECANE *or isomer*

~~19~~ 1266 110. 18344-37-1
HEPTADECANE, 2,6,10,14-TETRAMETHYL-

20 1336 130. 54105-67-8
HEPTADECANE, 2,6-DIMETHYL- *or isomer*

~~21~~ 1386 140. 1718-51-0
P-TERPHENYL-D14 *surrogate*

~~22~~ 1404 100. 629-92-5
NONADECANE

23 1531	110.	629-92-5
NONADECANE	or 130mer	

24 1587	160.	7206-19-1
3-OCTADECENE, (E)-	or 130mer	

25 1809	120.	00-00-0	5700
UNKNOWN			

DATA FILE: 30068103

QUANTERRA GC/MS TIC REPORT (Part 2)

CONCENTRATION = AREA(TIC)*CONC(IS)/AREA(IS)

#	FIT	PURITY	INT.			AREA	HEIGHT	AMOUNT			LIB	LIB #
			STD.	RT	RRT			AS ANALYZED		(UG/ML)		
1	954	915	2	1:54	0.206	1275650.	442368.	149.865			NB	3241.
2	988	937	2	2:28	0.267	1058960.	304107.	124.408			NB	5159.
3	905	600	2	4:56	0.535	311232.	78208.	36.564			UK	1.
4	900	659	2	5:15	0.570	280064.	64471.	32.902			NB	39230.
5	940	770	2	7:02	0.763	524416.	118784.	61.609			NB	37462.
6	934	735	2	7:15	0.787	681600.	76288.	80.075			NB	8109.
7	971	876	2	7:22	0.799	517888.	111232.	60.842			NB	11607.
8	928	808	2	8:13	0.890	615424.	133888.	72.301			NB	42196.
9	944	680	2	8:30	0.922	912896.	219904.	107.248			NB	19006.
10	682	318	2	8:39	0.937	556544.	73216.	65.383			UK	1.
11	965	833	2	9:07	0.989	750848.	136960.	88.211			NB	42196.
12	947	756	2	9:33	1.035	919808.	206592.	108.060			NB	25997.
13	965	853	2	10:33	1.143	825600.	197632.	96.992			NB	37465.
14	949	820	3	11:00	0.909	541952.	136192.	63.383			NB	37462.
15	971	854	3	11:32	0.953	1300990.	226560.	152.156			NB	37462.
16	924	767	3	12:24	1.025	443904.	134912.	51.916			NB	15969.
17	966	858	3	12:27	1.029	572928.	170752.	67.006			NB	37465.
18	967	784	3	13:15	1.096	324480.	112256.	37.949			NB	37465.
19	959	808	3	14:04	1.164	306688.	96256.	35.868			NB	42196.
20	888	698	4	14:51	0.860	340129.	83833.	42.894			NB	37462.
21	976	782	4	15:24	0.892	376192.	105344.	47.441			NB	32801.
22	931	695	4	15:36	0.903	272768.	66706.	34.399			NB	37465.
23	945	731	4	17:01	0.985	297984.	50304.	37.579			NB	37465.
24	968	767	4	17:39	1.021	421120.	74240.	53.107			NB	34410.
25	755	342	5	20:07	1.013	234560.	24576.	38.447			UK	1.

Library Search Data: 30068103 # 171 Base m/z: 43
 08/31/98 20:01:00 + 1:54 Cali: 30068103 # 3 RIC: 323072.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 630 matched at least 6 of the 16 largest peaks in the unknown

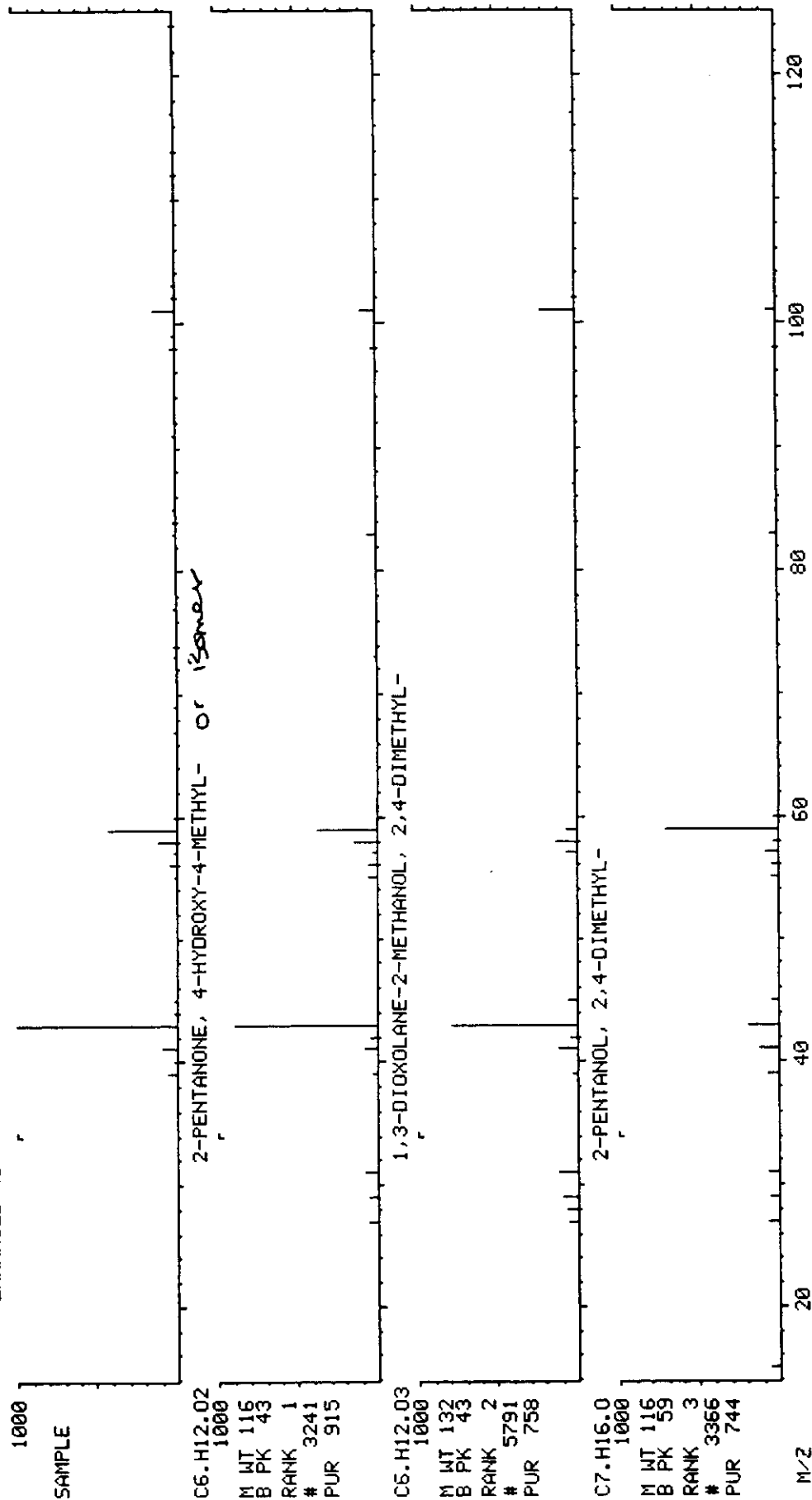
Rank	In.	Name
1	3241	2-PENTANONE, 4-HYDROXY-4-METHYL-
2	5791	1,3-DIOXOLANE-2-METHANOL, 2,4-DIMETHYL-
3	3366	2-PENTANOL, 2,4-DIMETHYL-
4	3288	ACETIC ACID, 1,1-DIMETHYLETHYL ESTER
5	3349	2-HEXANOL, 2-METHYL-
6	919	TERT-BUTYL HYDROPEROXIDE
7	5594	2-PROPANOL, 2-NITROSO-, ACETATE (ESTER)
8	8432	2-HEXANONE, 3-HYDROXY-3,5-DIMETHYL-
9	8425	2-HEPTANONE, 3-HYDROXY-3-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C6.H12.O2	116	43	915	954	928
2	C6.H12.O3	132	43	758	833	785
3	C7.H16.O	116	59	744	767	744
4	C6.H12.O2	116	43	728	780	742
5	C7.H16.O	116	59	709	744	709
6	C4.H10.O2	90	59	709	864	732
7	C5.H9.O3.N	131	43	708	819	799
8	C8.H16.O2	144	59	670	733	697
9	C8.H16.O2	144	59	657	713	678

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	123-42-2
2	---	---	---	---	53951-43-2
3	---	---	---	---	625-06-9
4	---	---	---	---	540-88-5
5	---	---	---	---	625-23-0
6	---	---	---	---	75-91-2
7	---	---	---	---	6931-04-0
8	---	---	---	---	6321-14-8
9	---	---	---	---	13757-91-0

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 20:01:00 + 1:54
SAMPLE: T-MM5-2 1/35A/1ML
CONDOS.: UG/ML *1ML *100Z/100Z *(NA/NA >1/35A NA M
ENHANCED (S 15B 2N 0T)

DATA: 30068103 # 171
CALL: 30068103 # 3
BASE M/Z: 43
RIC: 323072.



Library Search Data: 30068103 # 222 Base m/z: 43
 08/31/98 20:01:00 + 2:28 Cali: 30068103 # 3 RIC: 227328.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 921 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 5159 NONANE
 2 5151 HEPTANE, 3,4-DIMETHYL-
 3 3086 HEXANE, 2,4-DIMETHYL-
 4 5141 HEPTANE, 2,4-DIMETHYL-
 5 19015 DECANE, 2,5,6-TRIMETHYL-
 6 11607 UNDECANE
 7 8085 HEPTANE, 2,4,6-TRIMETHYL-
 8 19028 DECANE, 2,4,6-TRIMETHYL-
 9 5144 HEPTANE, 4-ETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C9.H20	128	43	937	988	937
2	C9.H20	128	43	899	925	899
3	C8.H18	114	43	894	964	904
4	C9.H20	128	43	879	933	888
5	C13.H28	184	57	874	914	884
6	C11.H24	156	43	868	929	905
7	C10.H22	142	43	865	909	881
8	C13.H28	184	43	861	890	906
9	C9.H20	128	43	861	910	868

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	111-84-2
2	---	---	---	---	922-28-1
3	---	---	---	---	589-43-5
4	---	---	---	---	2213-23-2
5	---	---	---	---	62108-23-0
6	---	---	---	---	1120-21-4
7	---	---	---	---	2613-61-8
8	---	---	---	---	62108-27-4
9	---	---	---	---	2216-32-2

DATA: 30058103 # 222
CALI: 30058103 # 3
BASE M/Z: 43
RIC: 227328.

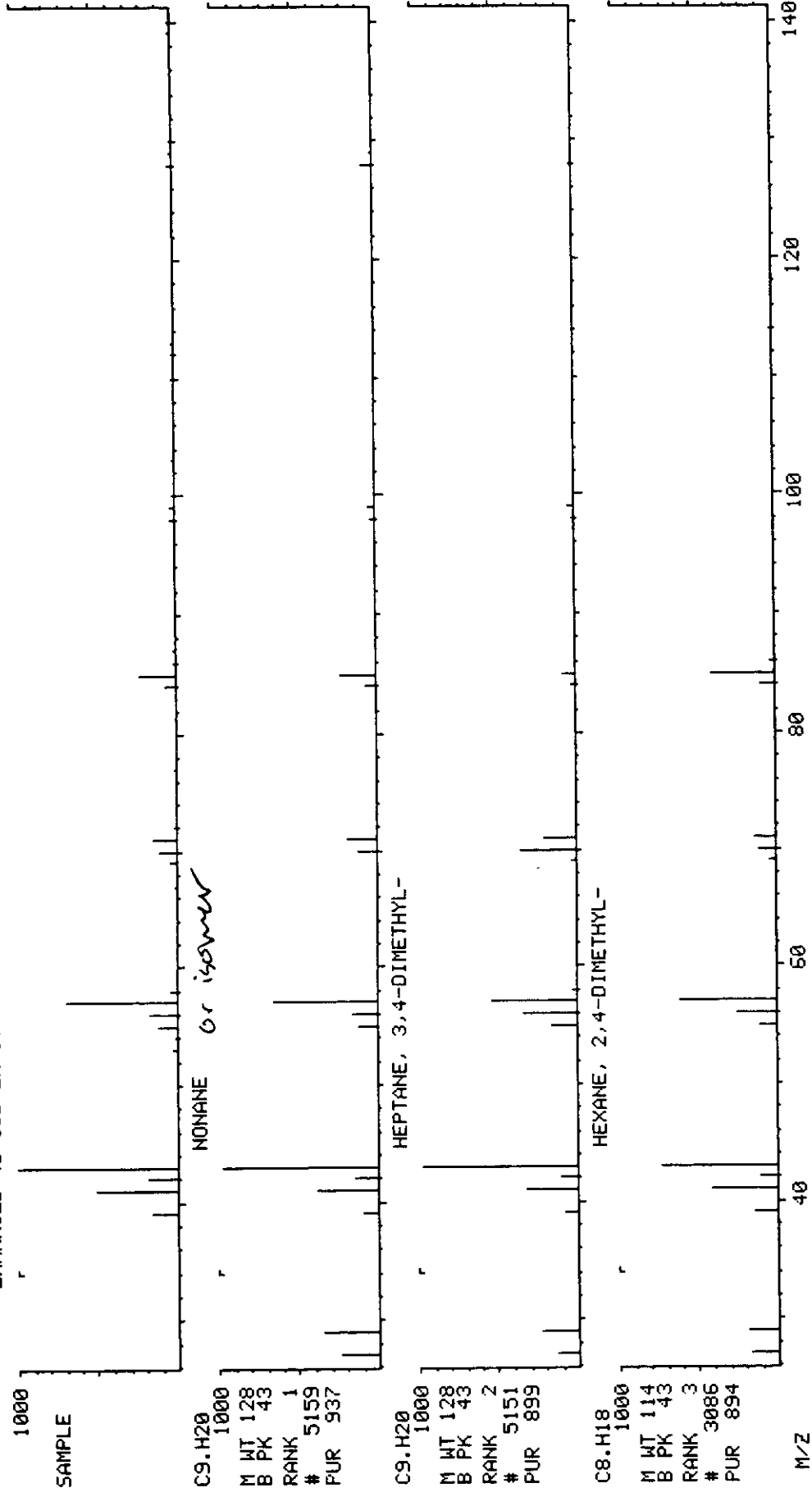
MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 20:01:00 + 2:28

SAMPLE: T-MM5-2 1/35A/1ML INST. ID: F16

CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M

ENHANCED (S 15B 2N 0T)



Library Search Data: 30068103 # 444 Base m/z: 43
 08/31/98 20:01:00 + 4:56 Cali: 30068103 # 3 RIC: 66688.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 463 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 19523 1-DECANOL, 2-ETHYL-
 2 14793 1-UNDECENE, 4-METHYL-
 3 35931 HEXADECANE, 1-CHLORO-
 4 12074 1-HEPTANOL, 2-PROPYL-
 5 15969 HYDROXYLAMINE, O-DECYL-
 6 19015 DECANE, 2,5,6-TRIMETHYL-
 7 40866 OCTADECANE, 1-CHLORO-
 8 11607 UNDECANE
 9 11602 OCTANE, 2,4,6-TRIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H26.O	186	57	600	905	627
2	C12.H24	168	43	597	909	635
3	C16.H33.CL	260	57	595	824	635
4	C10.H22.O	158	43	589	928	607
5	C10.H23.O.N	173	43	588	929	624
6	C13.H28	184	57	583	916	596
7	C18.H37.CL	288	57	583	816	648
8	C11.H24	156	43	579	957	600
9	C11.H24	156	57	576	939	585

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	21078-65-9
2	---	---	---	---	74630-39-0
3	---	---	---	---	4860-03-1
4	---	---	---	---	10042-59-8
5	---	---	---	---	29812-79-1
6	---	---	---	---	62108-23-0
7	---	---	---	---	3386-33-2
8	---	---	---	---	1120-21-4
9	---	---	---	---	62016-37-9

DATA: 30058103 # 444
CALI: 30058103 # 3

BASE M/Z: 43
RIC: 66688.

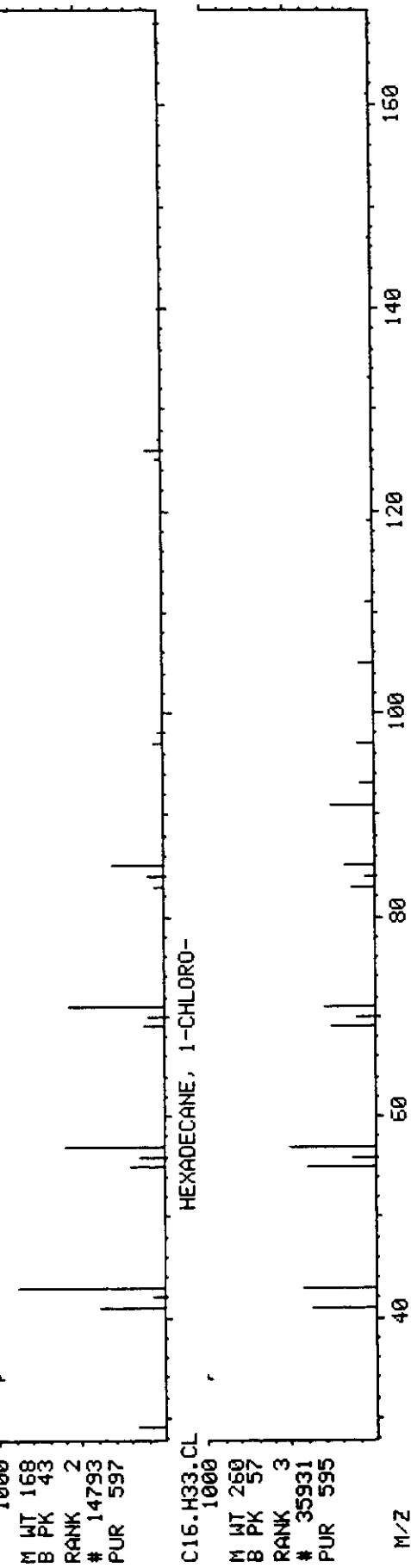
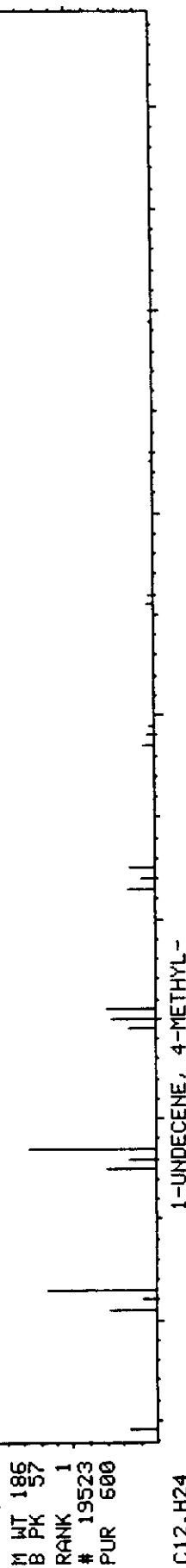
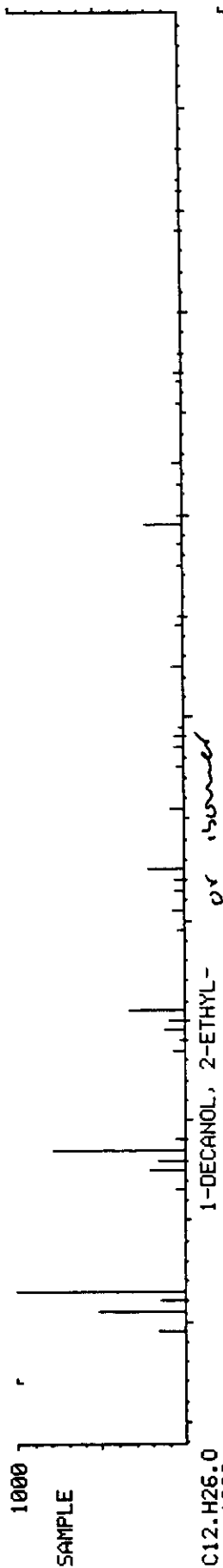
MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 20:01:00 + 4:56

SAMPLE: T-MM5-2 1/35A/1ML INST. ID: F16

CONDS.: UG/ML *1ML *1002/1002 *(NA/NA >1/35A NA M

ENHANCED (S 158 2N 0T)



00
00
00

Library Search Data: 30068103 # 473 Base m/z: 43
 08/31/98 20:01:00 + 5:16 Cali: 30068103 # 3 RIC: 59200.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 459 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 39230 9-EICOSYNE
 2 22482 2-TRIDECEN-1-OL, (E)-
 3 34800 (R)-(-)-(Z)-14-METHYL-8-HEXADECEN-1-OL
 4 25971 OXIRANE, DODECYL-
 5 29222 (Z)6-PENTADECEN-1-OL
 6 37449 OCTADECANAL
 7 36677 7-OCTADECYNE, 2-METHYL-
 8 37444 OXIRANE, HEXADECYL-
 9 17938 1,12-TRIDECADIENE

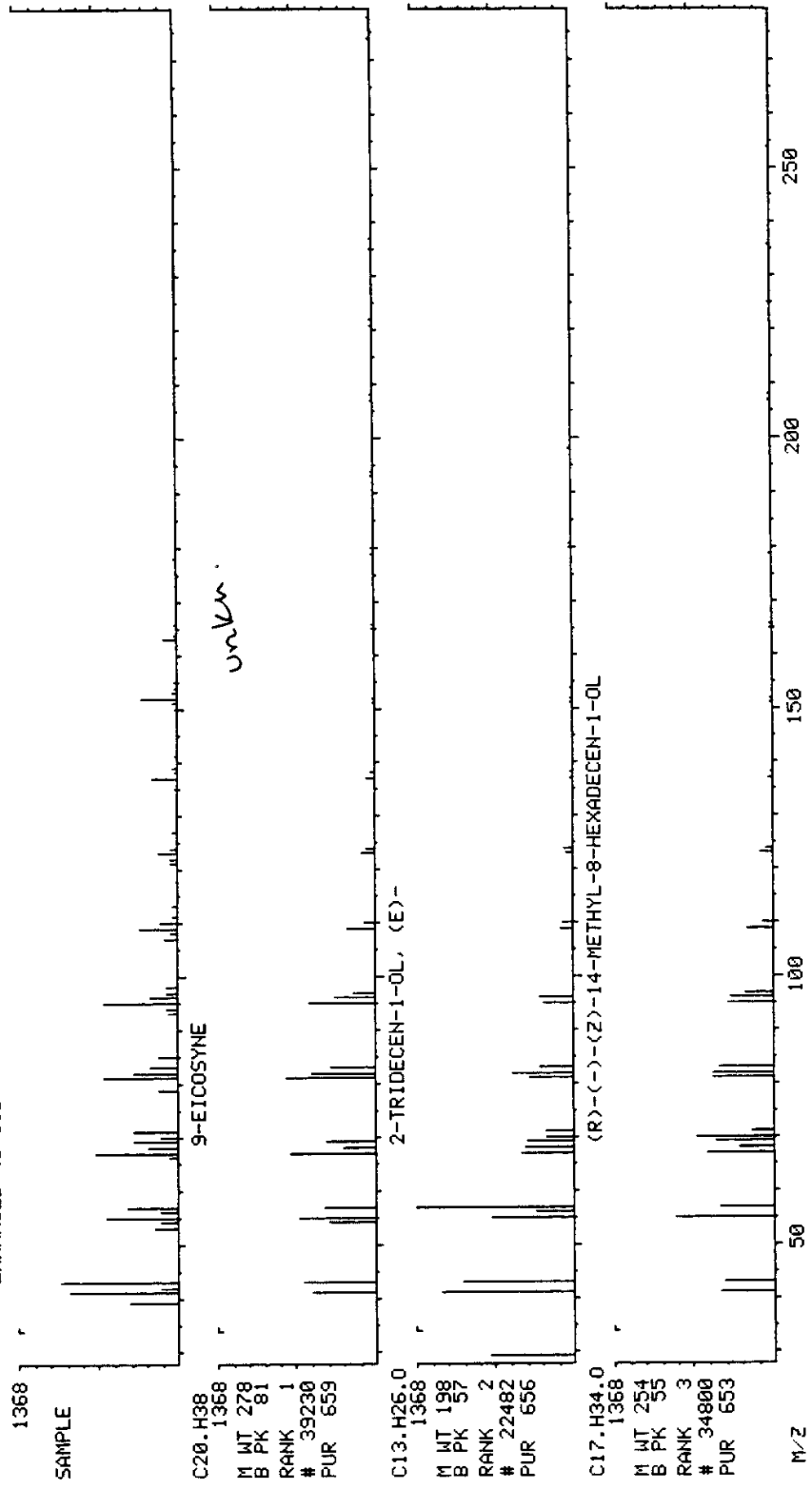
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C20.H38	278	81	659	900	677
2	C13.H26.O	198	57	656	854	662
3	C17.H34.O	254	55	653	850	663
4	C14.H28.O	212	41	650	918	650
5	C15.H30.O	226	67	636	869	654
6	C18.H36.O	268	43	635	828	677
7	C19.H36	264	81	626	889	642
8	C18.H36.O	268	57	624	808	667
9	C13.H24	180	55	624	876	627

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	71899-38-2
2	—	—	—	—	74962-98-4
3	—	—	—	—	30689-78-2
4	—	—	—	—	3234-28-4
5	—	—	—	—	68797-95-5
6	—	—	—	—	638-66-4
7	—	—	—	—	35354-38-2
8	—	—	—	—	7390-81-0
9	—	—	—	—	21964-48-7

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:01:00 + 5:16
 SAMPLE: T-MM5-2 1/35A/1ML
 COND.: UG/ML *1ML *1007/1002 *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068103 # 473
 CALI: 30068103 # 3

BASE M/Z: 43
 RIC: 59200.



Library Search Data: 30068103 # 633 Base m/z: 57
 08/31/98 20:01:00 + 7:02 Cali: 30068103 # 3 RIC: 109056.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (\$ 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 862 matched at least 7 of the 16 largest peaks in the unknown

Rank	In.	Name
1	37462	HEPTADECANE, 2,6-DIMETHYL-
2	22534	TRIDECANE, 7-METHYL-
3	14799	1-DECENE, 3,4-DIMETHYL-
4	15969	HYDROXYLAMINE, O-DECYL-
5	19523	1-DECANOL, 2-ETHYL-
6	42196	HEPTADECANE, 2,6,10,14-TETRAMETHYL-
7	14751	2-UNDECENE, 5-METHYL-
8	32422	OCTANE, 1,1'-OXYBIS-
9	8539	1-PENTANOL, 4-METHYL-2-PROPYL-

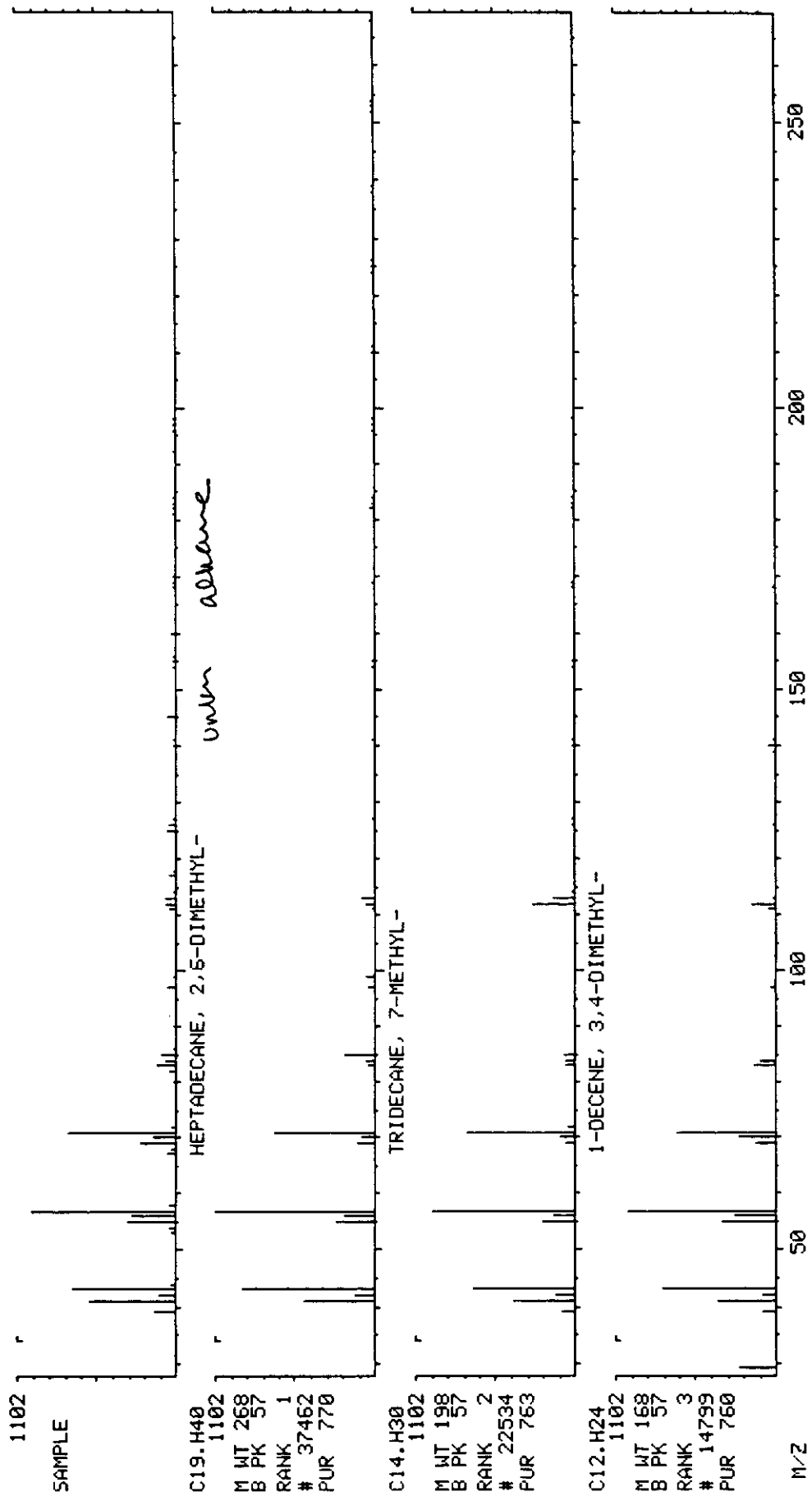
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	770	940	780
2	C14.H30	198	57	763	928	763
3	C12.H24	168	57	760	958	761
4	C10.H23.O.N	173	43	738	943	740
5	C12.H26.O	186	57	736	925	755
6	C21.H44	296	57	735	926	750
7	C12.H24	168	57	732	957	740
8	C16.H34.O	242	57	728	961	736
9	C9.H20.O	144	57	727	971	733

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	54105-67-8
2	---	---	---	---	26730-14-3
3	---	---	---	---	50871-03-9
4	---	---	---	---	29812-79-1
5	---	---	---	---	21078-65-9
6	---	---	---	---	18344-37-1
7	---	---	---	---	56851-34-4
8	---	---	---	---	629-82-3
9	---	---	---	---	54004-41-0

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:01:00 + 7:02
 SAMPLE: T-MM5-2 1/35A/1ML
 COND.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068103 # 633
 CALI: 30068103 # 3

BASE M/2: 57
 RIC: 109056.



Library Search Data: 30068103 # 653 Base m/z: 142
 08/31/98 20:01:00 + 7:16 Cali: 30068103 # 3 RIC: 43008.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

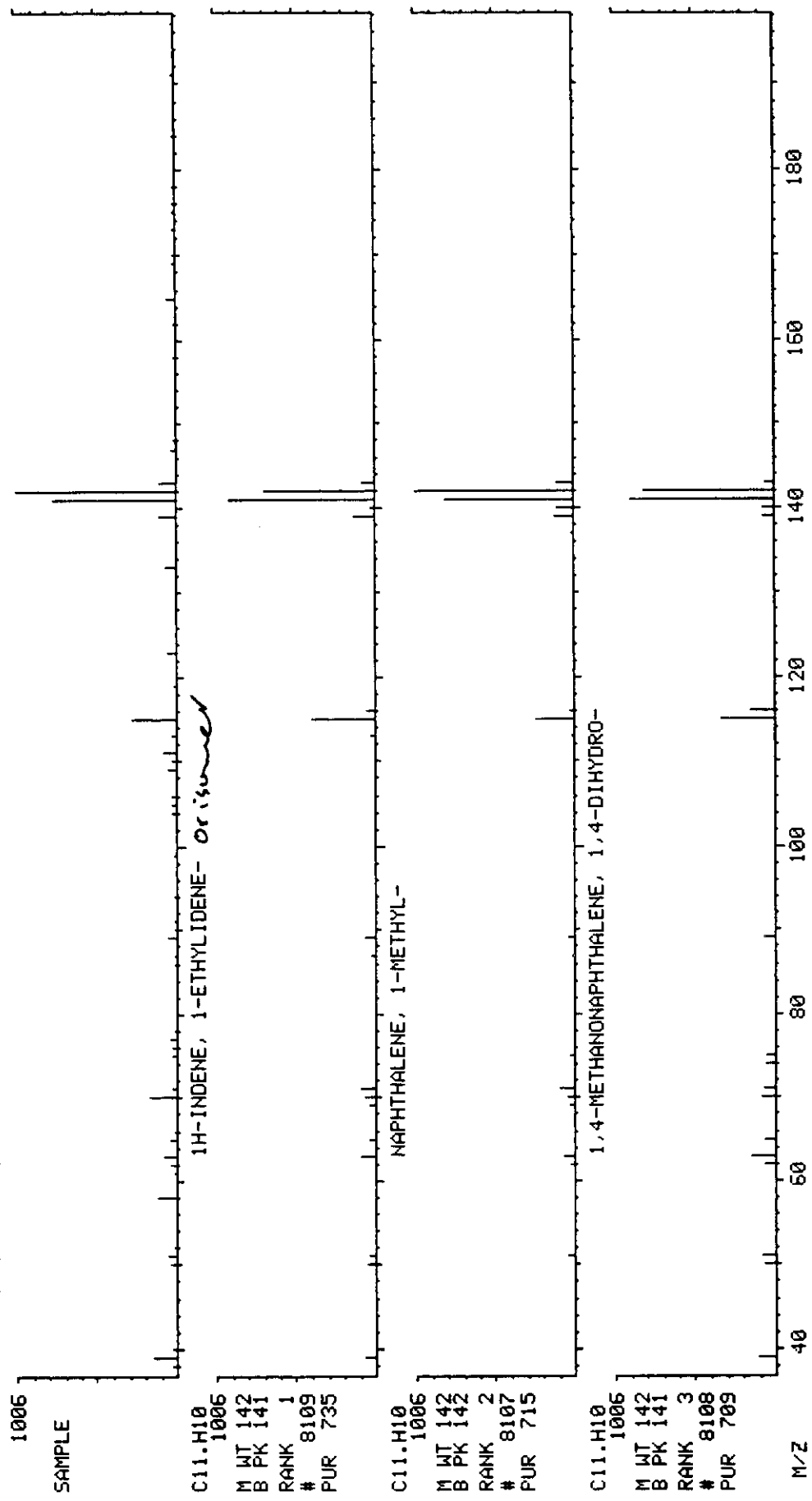
62231 spectra in LIBRARYNB searched for maximum PURITY
 227 matched at least 4 of the 16 largest peaks in the unknown

Rank In. Name
 1 8109 1H-INDENE, 1-ETHYLIDENE-
 2 8107 NAPHTHALENE, 1-METHYL-
 3 8108 1,4-METHANONAPHTHALENE, 1,4-DIHYDRO-
 4 8111 NAPHTHALENE, 2-METHYL-
 5 8110 BENZOCYCLOHEPTATRIENE
 6 15368 BICYCLO[2.2.1]HEPT-2-ENE, 2-PHENYL-
 7 19065 NAPHTHALENE, 2-BUTYL-
 8 7994 BENZENEACETONITRILE, 2-CYANO-
 9 7992 BENZENEACETONITRILE, 4-CYANO-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C11.H10	142	141	735	934	763
2	C11.H10	142	142	715	924	762
3	C11.H10	142	141	709	901	768
4	C11.H10	142	142	703	914	758
5	C11.H10	142	141	672	903	708
6	C13.H14	170	142	564	748	632
7	C14.H16	184	141	563	742	651
8	C9.H6.N2	142	142	560	776	620
9	C9.H6.N2	142	142	551	777	630

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	2471-83-2
2	—	—	—	—	90-12-0
3	—	—	—	—	4453-90-1
4	—	—	—	—	91-57-6
5	—	—	—	—	264-09-5
6	—	—	—	—	4237-08-5
7	—	—	—	—	1134-62-9
8	—	—	—	—	3759-28-2
9	—	—	—	—	876-31-3

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:01:00 + 7:16
 SAMPLE: T-MM5-2 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068103 # 653
 CALI: 30068103 # 3
 BASE M/Z: 142
 RIC: 43008.



Library Search Data: 30068103 # 663 Base m/z: 57
 08/31/98 20:01:00 + 7:22 Cali: 30068103 # 3 RIC: 109824.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 723 matched at least 7 of the 16 largest peaks in the unknown

Rank In.	Name
1	11607 UNDECANE
2	11602 OCTANE, 2,4,6-TRIMETHYL-
3	15356 DECANE, 2,4-DIMETHYL-
4	19016 UNDECANE, 4,7-DIMETHYL-
5	8104 OCTANE, 3,5-DIMETHYL-
6	19015 DECANE, 2,5,6-TRIMETHYL-
7	19026 DECANE, 2,6,8-TRIMETHYL-
8	19028 DECANE, 2,4,6-TRIMETHYL-
9	19013 DECANE, 2,5,9-TRIMETHYL-

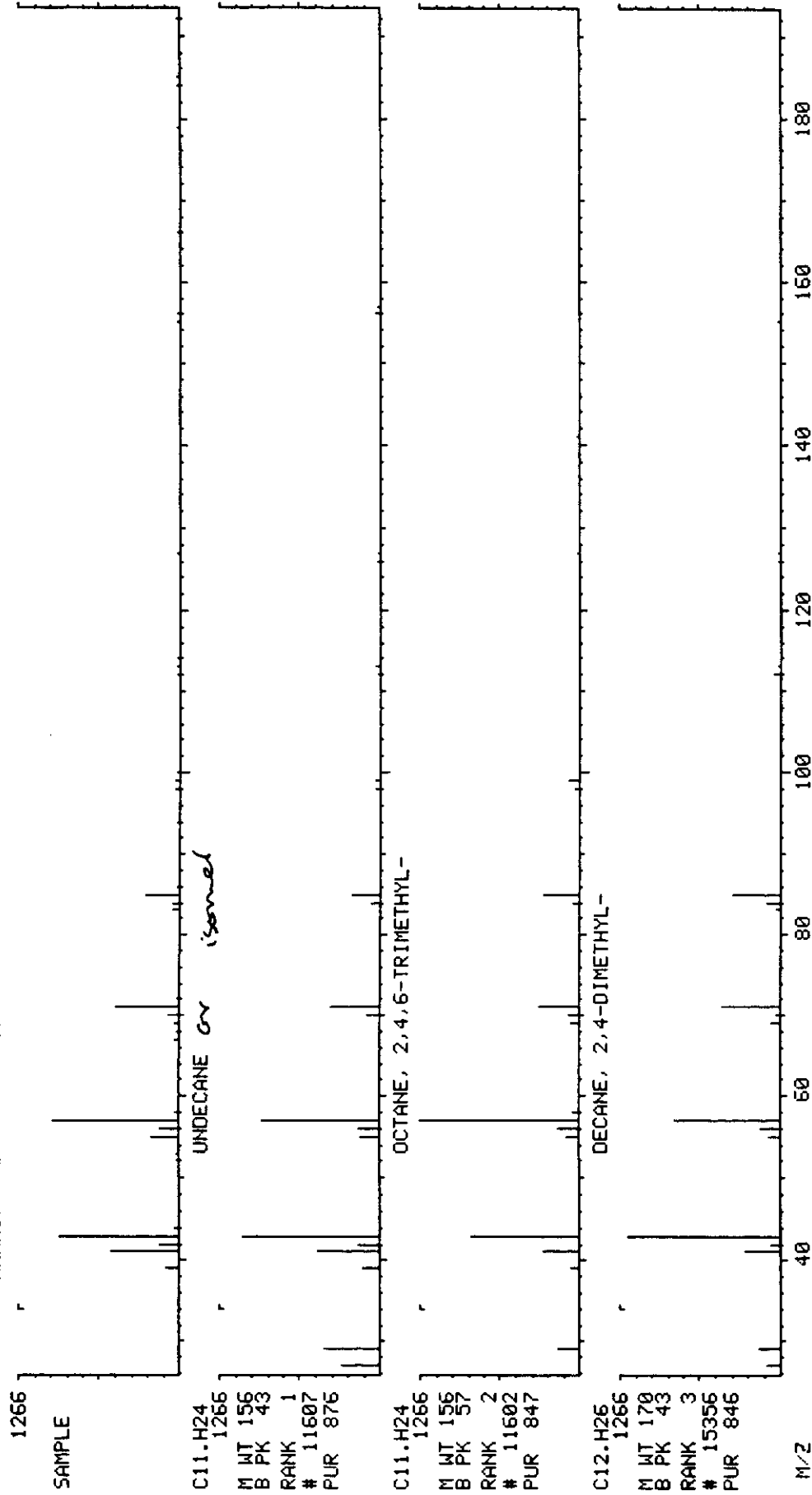
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C11.H24	156	43	876	971	882
2	C11.H24	156	57	847	966	849
3	C12.H26	170	43	846	955	850
4	C13.H28	184	43	846	934	869
5	C10.H22	142	57	845	936	853
6	C13.H28	184	57	841	924	844
7	C13.H28	184	57	840	933	849
8	C13.H28	184	43	839	932	852
9	C13.H28	184	57	838	935	838

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	1120-21-4
2	---	---	---	---	62016-37-9
3	---	---	---	---	2801-84-5
4	---	---	---	---	17301-32-5
5	---	---	---	---	15869-93-9
6	---	---	---	---	62108-23-0
7	---	---	---	---	62108-26-3
8	---	---	---	---	62108-27-4
9	---	---	---	---	62108-22-9

MID LIBRARY SEARCH (LIBRARY.NB)
 08/31/98 20:01:00 + 7:22
 SAMPLE: T-MM5-2 1/35A/1ML
 COND.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068103 # 663
 CALI: 30068103 # 3

BASE M/Z: 57
 RIC: 109824.



Library Search Data: 30068103 # 739 Base m/z: 57
 08/31/98 20:01:00 + 8:13 Cali: 30068103 # 3 RIC: 131328.
 Sample: T-MMS-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

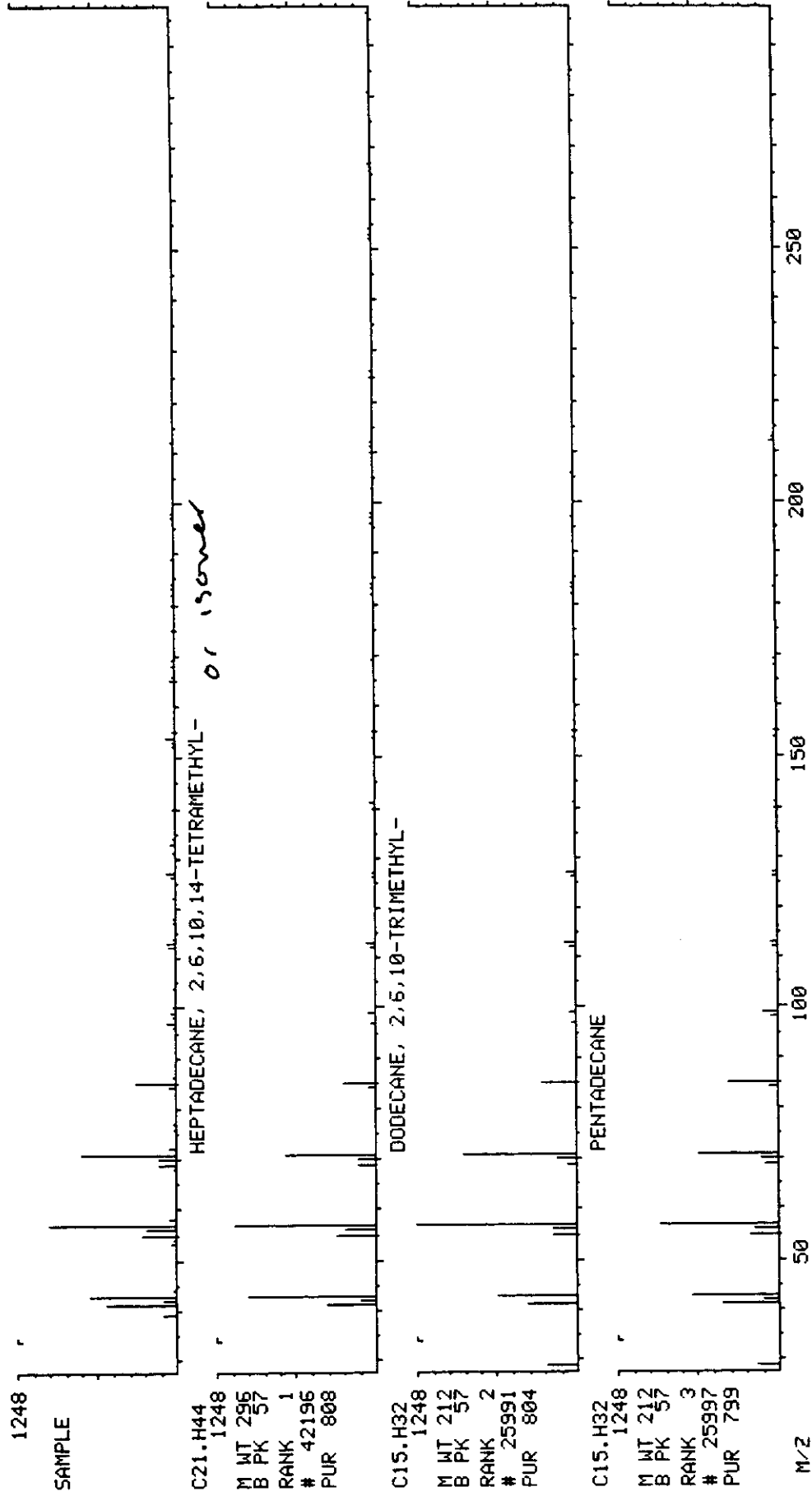
62231 spectra in LIBRARYNB searched for maximum PURITY
 799 matched at least 7 of the 16 largest peaks in the unknown

Rank	In.	Name
1	42196	HEPTADECANE, 2,6,10,14-TETRAMETHYL-
2	25991	DODECANE, 2,6,10-TRIMETHYL-
3	25997	PENTADECANE
4	37462	HEPTADECANE, 2,6-DIMETHYL-
5	22535	DODECANE, 4,6-DIMETHYL-
6	26001	DODECANE, 2,7,10-TRIMETHYL-
7	22530	TETRADECANE
8	18998	UNDECANE, 3,7-DIMETHYL-
9	37465	NONADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C21.H44	296	57	808	928	839
2	C15.H32	212	57	804	960	809
3	C15.H32	212	57	799	941	812
4	C19.H40	268	57	792	924	825
5	C14.H30	198	57	789	941	796
6	C15.H32	212	57	782	972	792
7	C14.H30	198	43	781	916	796
8	C13.H28	184	43	779	929	781
9	C19.H40	268	57	776	904	829

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	18344-37-1
2	—	—	—	—	3891-98-3
3	—	—	—	—	629-62-9
4	—	—	—	—	54105-67-8
5	—	—	—	—	61141-72-8
6	—	—	—	—	74645-98-0
7	—	—	—	—	629-59-4
8	—	—	—	—	17301-29-0
9	—	—	—	—	629-92-5

MID LIBRARY SEARCH <LIBRARYNB>
 08/31/98 20:01:00 + 8:13
 SAMPLE: T-MMS-2 1/35A/1ML
 CONDS.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED <S 15B 2N 0T>
 DATA: 30068103 # 739
 CALI: 30068103 # 3
 BASE M/Z: 57
 RIC: 131328.



Library Search Data: 30068103 # 765 Base m/z: 43
 08/31/98 20:01:00 + 8:31 Cali: 30068103 # 3 RIC: 193536.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 513 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 19006 DECANE, 6-ETHYL-2-METHYL-
 2 22530 TETRADECANE
 3 19007 DODECANE, 3-METHYL-
 4 29249 TETRADECANE, 6,9-DIMETHYL-
 5 39681 1-iodoundecane
 6 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 7 37252 DECANE, 1-iodo-
 8 25996 TRIDECANE, 4,8-DIMETHYL-
 9 11607 UNDECANE

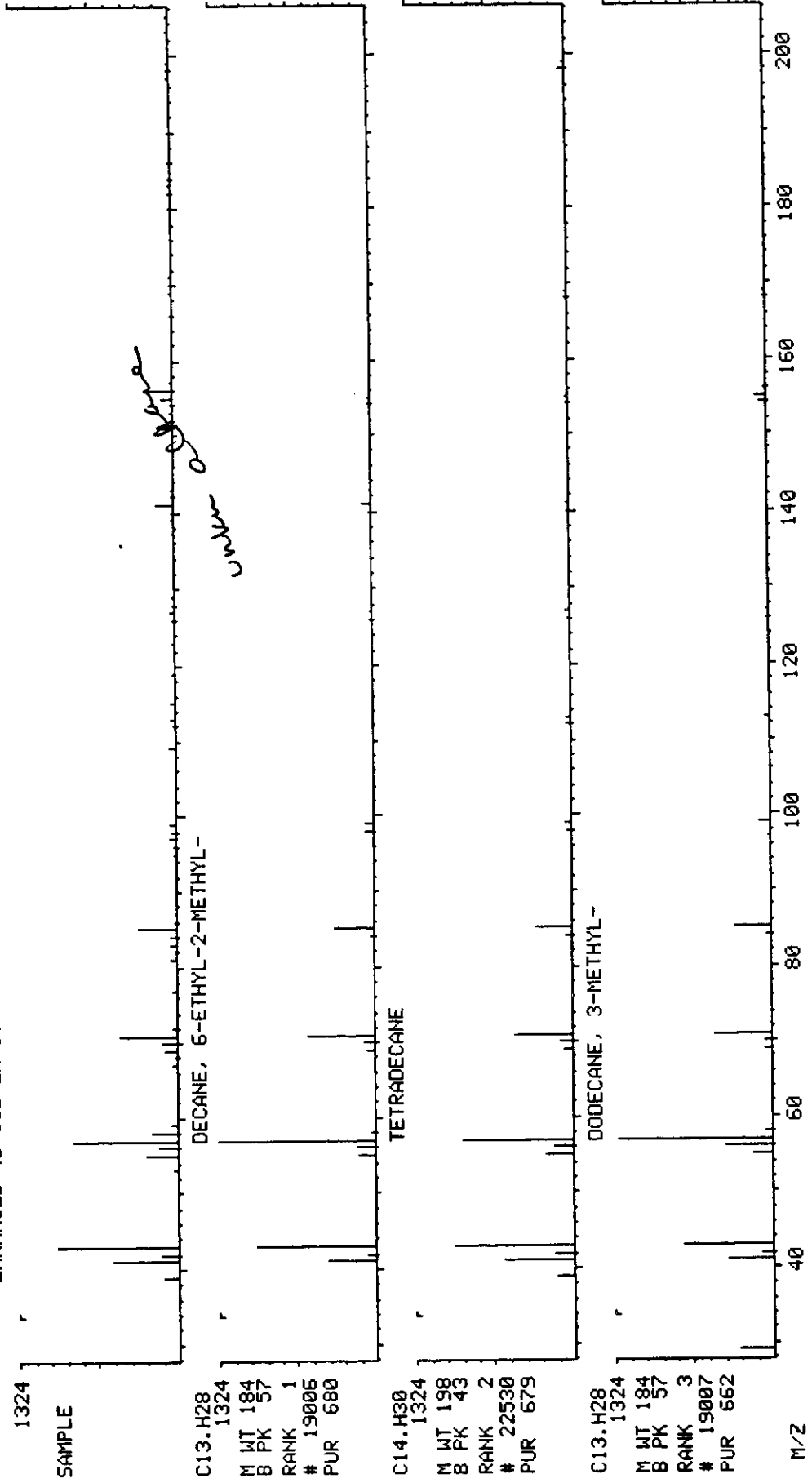
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H28	184	57	680	944	680
2	C14.H30	198	43	679	942	680
3	C13.H28	184	57	662	882	662
4	C16.H34	226	57	659	852	673
5	C11.H23.I	282	57	656	848	660
6	C21.H44	296	57	654	914	667
7	C10.H21.I	268	57	654	877	657
8	C15.H32	212	57	654	918	662
9	C11.H24	156	43	652	957	652

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	62108-21-8
2	—	—	—	—	629-59-4
3	—	—	—	—	17312-57-1
4	—	—	—	—	55045-13-1
5	—	—	—	—	4282-44-4
6	—	—	—	—	18344-37-1
7	—	—	—	—	2050-77-3
8	—	—	—	—	55030-62-1
9	—	—	—	—	1120-21-4

DATA: 30068103 # 765
CALI: 30068103 # 3

MTD LIBRARY SEARCH (LIBRARYNB)
08/31/98 20:01:00 + 8:31
SAMPLE: T-MM5-2 1/35A/1ML
COND5.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 158 2N 0T)

BASE M/Z: 43
RIC: 193536.



Library Search Data: 30068103 # 778 Base m/z: 141
 08/31/98 20:01:00 + 8:39 Cali: 30068103 # 3 RIC: 61952.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 554 matched at least 5 of the 16 largest peaks in the unknown

Rank In. Name
 1 29224 11-DODECEN-1-OL, 2,4,6-TRIMETHYL-, (R,R,R)-
 2 46107 2H-PYRAN-2-CARBOXYLIC ACID, 3,6-DIHYDRO-6-PROPOXY-, 4-(1-METHYLETHY*
 3 56499 9-OCTADECENOIC ACID (Z)-, 2-(ACETYLOXY)-1-[(ACETYLOXY)METHYL]ETHYL *
 4 32188 NONYLAMINE, 1-HEPTYL-
 5 16353 INDAN, 1,1,6,7-TETRAMETHYL-
 6 16347 1H-INDENE, 2,3-DIHYDRO-1,1,4,7-TETRAMETHYL-
 7 21930 CYCLOPROPANOL, 1-(3,7-DIMETHYL-1-OCTENYL)-
 8 56413 9,12-OCTADECADIENOIC ACID (Z,Z)-, 2-(ACETYLOXY)-1-[(ACETYLOXY)METHY*
 9 56303 9,12,15-OCTADECATRIENOIC ACID, 2-(ACETYLOXY)-1-[(ACETYLOXY)METHYL]E*

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C15.H30.O	226	55	318	682	384
2	C19.H32.O4	324	83	318	752	378
3	C25.H44.O6	440	43	305	649	386
4	C16.H35.N	241	128	288	723	305
5	C13.H18	174	159	276	591	280
6	C13.H18	174	159	270	591	275
7	C13.H24.O	196	83	269	768	315
8	C25.H42.O6	438	43	266	741	323
9	C25.H40.O6	436	43	264	759	319

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	27829-54-5
2	—	—	—	—	55124-82-8
3	—	—	—	—	55401-63-3
4	—	—	—	—	24539-83-1
5	—	—	—	—	16204-58-3
6	—	—	—	—	1078-04-2
7	—	—	—	—	65147-72-0
8	—	—	—	—	55320-03-1
9	—	—	—	—	55320-01-9

MID LIBRARY SEARCH (LIBRARYNB)

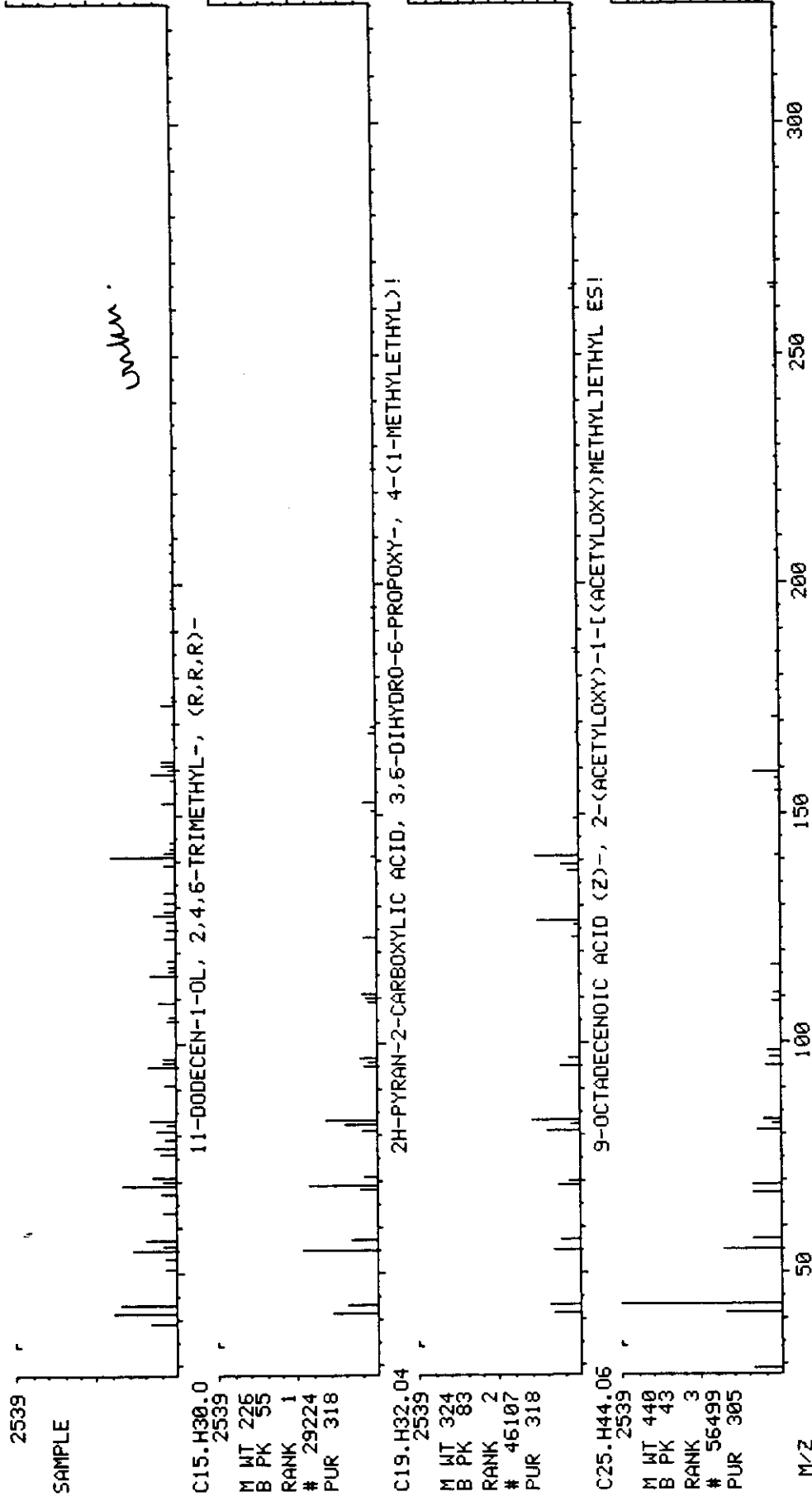
08/31/98 20:01:00 + 8:39

SAMPLE: T-MM5-2 1/35A/1ML

COND.S.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (5 158 2N 0T)

DATA: 30068103 # 778
CALI: 30068103 # 3

BASE M/2: 141
RIC: 61952.



Library Search Data: 30068103 # 821 Base m/z: 57
 08/31/98 20:01:00 + 9:08 Cali: 30068103 # 3 RIC: 128128.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYMB searched for maximum PURITY
 284 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 19016 UNDECANE, 4,7-DIMETHYL-
 4 18987 UNDECANE, 2,8-DIMETHYL-
 5 15353 2,6-DIMETHYLDECANE
 6 19022 UNDECANE, 4,8-DIMETHYL-
 7 37465 NONADECANE
 8 19023 DECANE, 2,6,7-TRIMETHYL-
 9 19004 UNDECANE, 4,6-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C21.H44	296	57	833	965	854
2	C19.H40	268	57	824	946	847
3	C13.H28	184	43	796	948	802
4	C13.H28	184	43	795	966	799
5	C12.H26	170	43	790	924	802
6	C13.H28	184	43	790	960	791
7	C19.H40	268	57	787	917	840
8	C13.H28	184	57	786	950	786
9	C13.H28	184	57	785	964	785

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	18344-37-1
2	---	---	---	---	54105-67-8
3	---	---	---	---	17301-32-5
4	---	---	---	---	17301-25-6
5	---	---	---	---	13150-81-7
6	---	---	---	---	17301-33-6
7	---	---	---	---	629-92-5
8	---	---	---	---	62108-25-2
9	---	---	---	---	17312-82-2

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 20:01:00 + 9:08

SAMPLE: T-MM5-2 1/35A/1ML

CONDS.: UG/ML *1ML *100%/100% *(NA/NA) 1/35A NA M

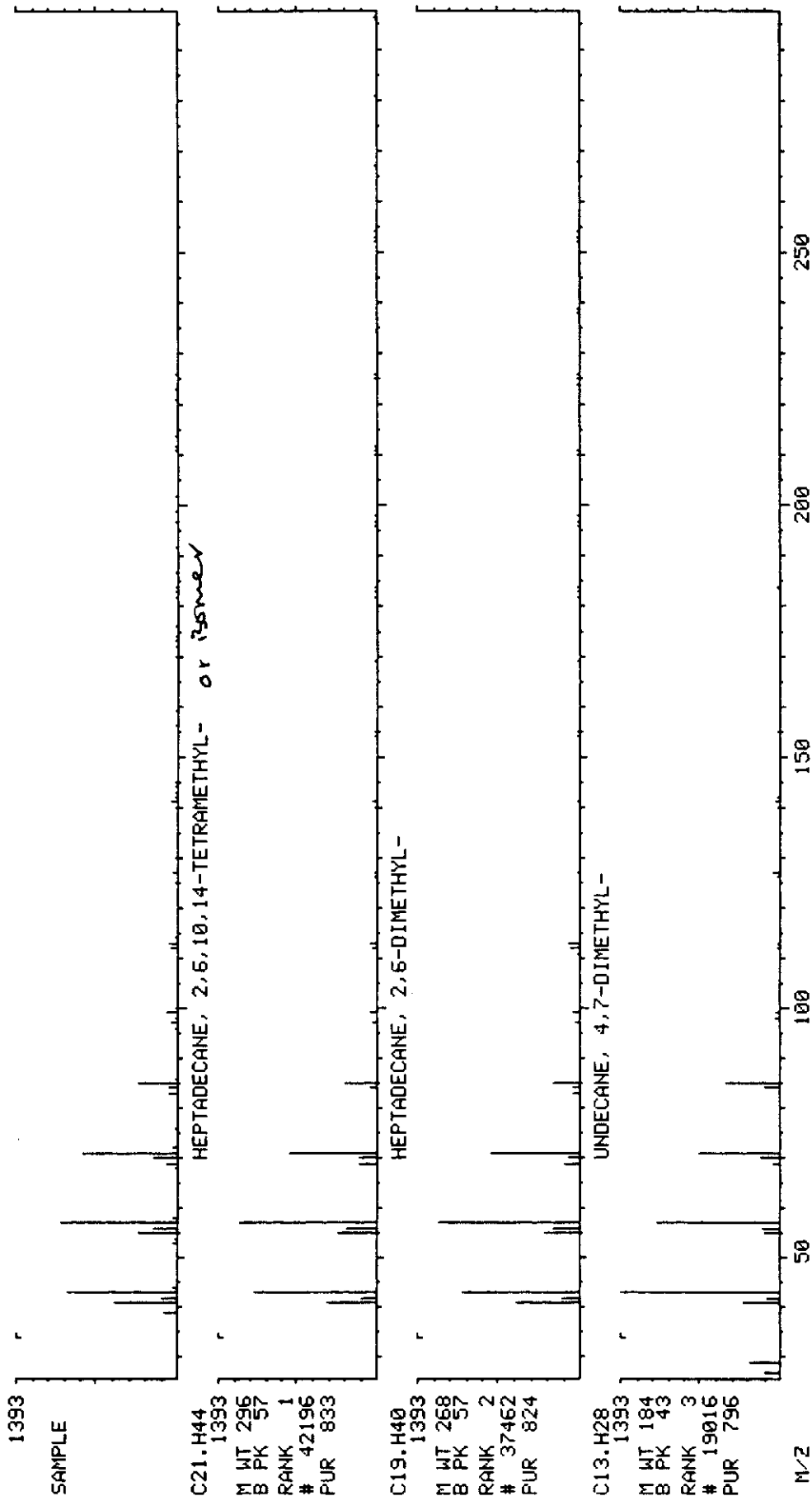
ENHANCED (S 15B 2N 0T)

DATA: 30068103 # 821

CALI: 30068103 # 3

BASE M/Z: 57

RIC: 128128.



Library Search Data: 30068103 # 859 Base m/z: 57
 08/31/98 20:01:00 + 9:33 Cali: 30068103 # 3 RIC: 195328.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 835 matched at least 7 of the 16 largest peaks in the unknown

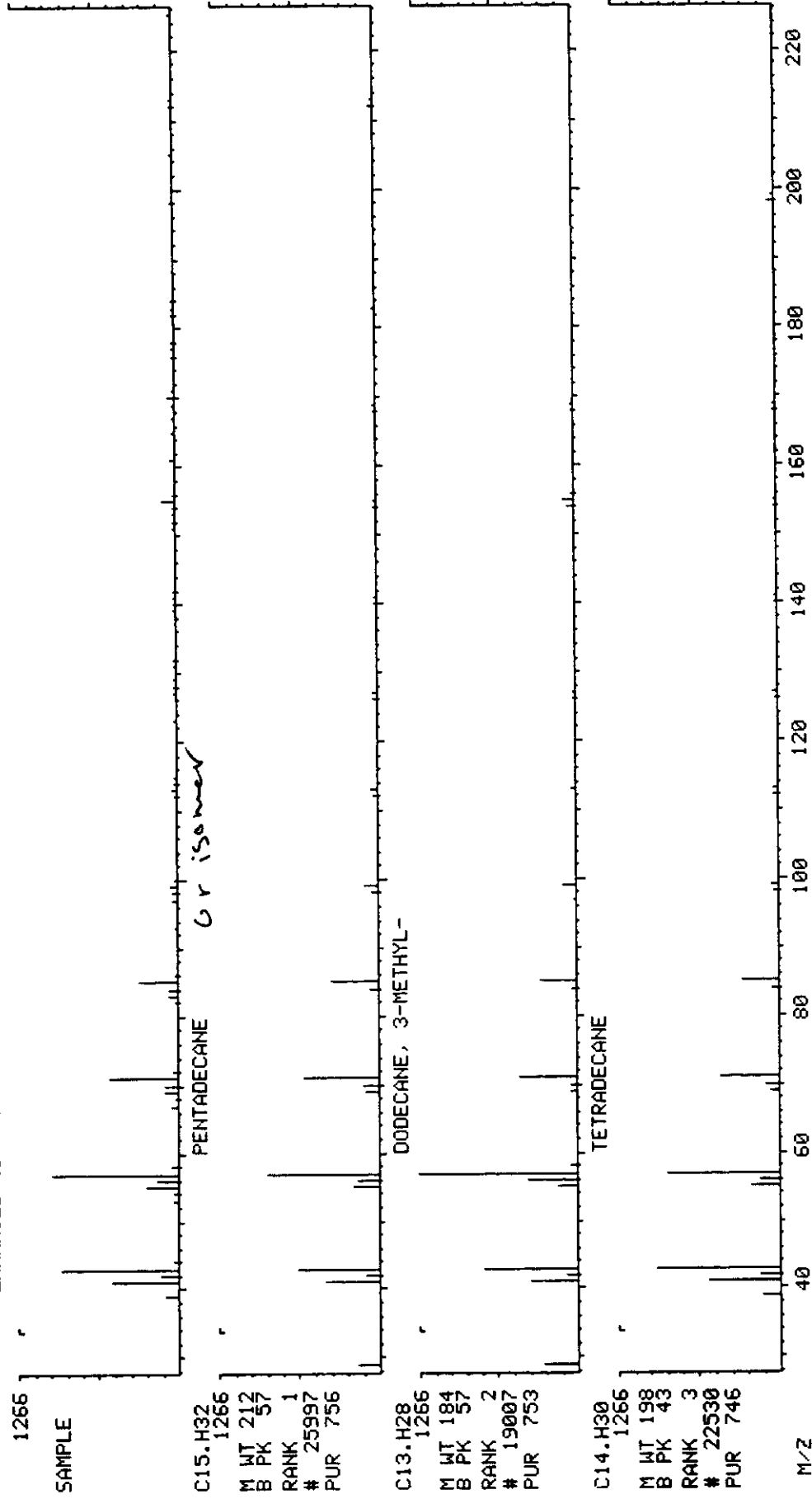
Rank In. Name
 1 25997 PENTADECANE
 2 19007 DODECANE, 3-METHYL-
 3 22530 TETRADECANE
 4 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 5 15343 DODECANE
 6 15353 2,6-DIMETHYLDECANE
 7 39681 1-iodoundecane
 8 37462 HEPTADECANE, 2,6-DIMETHYL-
 9 19006 DECANE, 6-ETHYL-2-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C15.H32	212	57	756	947	769
2	C13.H28	184	57	753	959	757
3	C14.H30	198	43	746	932	759
4	C21.H44	296	57	744	930	762
5	C12.H26	170	57	743	985	748
6	C12.H26	170	43	742	911	747
7	C11.H23.I	282	57	740	965	745
8	C19.H40	268	57	739	918	755
9	C13.H28	184	57	739	952	744

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	629-62-9
2	---	---	---	---	17312-57-1
3	---	---	---	---	629-59-4
4	---	---	---	---	18344-37-1
5	---	---	---	---	112-40-3
6	---	---	---	---	13150-81-7
7	---	---	---	---	4282-44-4
8	---	---	---	---	54105-67-8
9	---	---	---	---	62108-21-8

DATA: 30068103 # 859
CALI: 30068103 # 3

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 20:01:00 + 9:33
SAMPLE: T-MMS-2 1/35A/1ML INST. ID: F16
CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)



Library Search Data: 30068103 # 949 Base m/z: 57
 08/31/98 20:01:00 + 10:33 Cali: 30068103 # 3 RIC: 184064.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARY8 searched for maximum PURITY
 338 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 25997 PENTADECANE
 4 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 5 15969 HYDROXYLAMINE, O-DECYL-
 6 19016 UNDECANE, 4,7-DIMETHYL-
 7 46161 TRICOSANE
 8 25994 DODECANE, 2,6,11-TRIMETHYL-
 9 15353 2,6-DIMETHYLDECANE

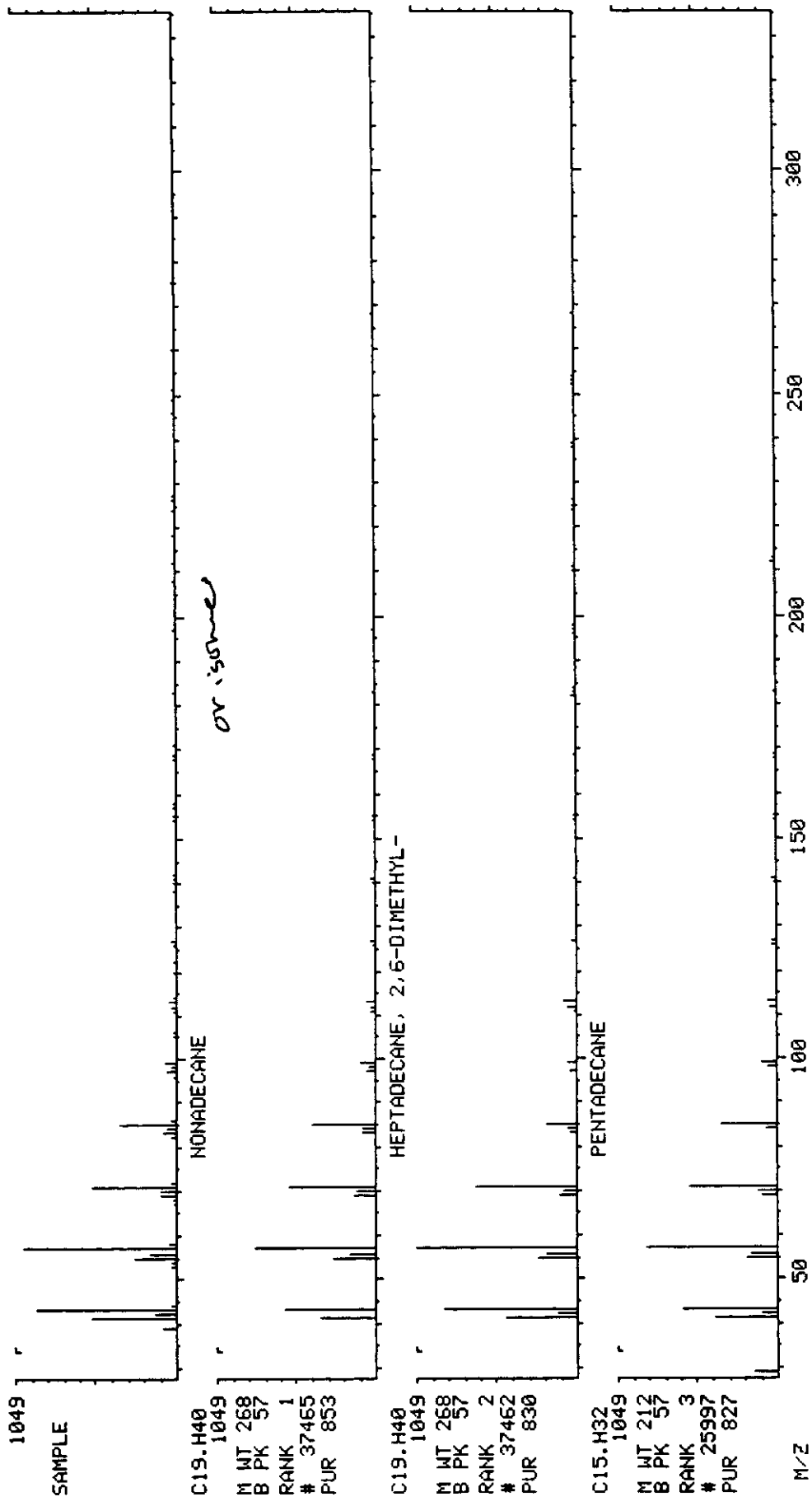
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	853	965	868
2	C19.H40	268	57	830	942	855
3	C15.H32	212	57	827	967	830
4	C21.H44	296	57	823	950	844
5	C10.H23.O.N	173	43	821	967	823
6	C13.H28	184	43	806	963	806
7	C23.H48	324	43	801	904	873
8	C15.H32	212	57	793	962	795
9	C12.H26	170	43	792	927	801

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	629-62-9
4	—	—	—	—	18344-37-1
5	—	—	—	—	29812-79-1
6	—	—	—	—	17301-32-5
7	—	—	—	—	638-67-5
8	—	—	—	—	31295-56-4
9	—	—	—	—	13150-81-7

DATA: 30068103 # 949
CALI: 30068103 # 3
BASE M/Z: 57
RIC: 184054.

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 20:01:00 + 10:33
SAMPLE: T-MM5-2 1/35A/1ML
CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA >)/1/35A NA M
ENHANCED (S 158 2N 0T)

INST. ID: F16



Library Search Data: 30068103 # 989 Base m/z: 57
 08/31/98 20:01:00 + 11:00 Cali: 30068103 # 3 RIC: 133376.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

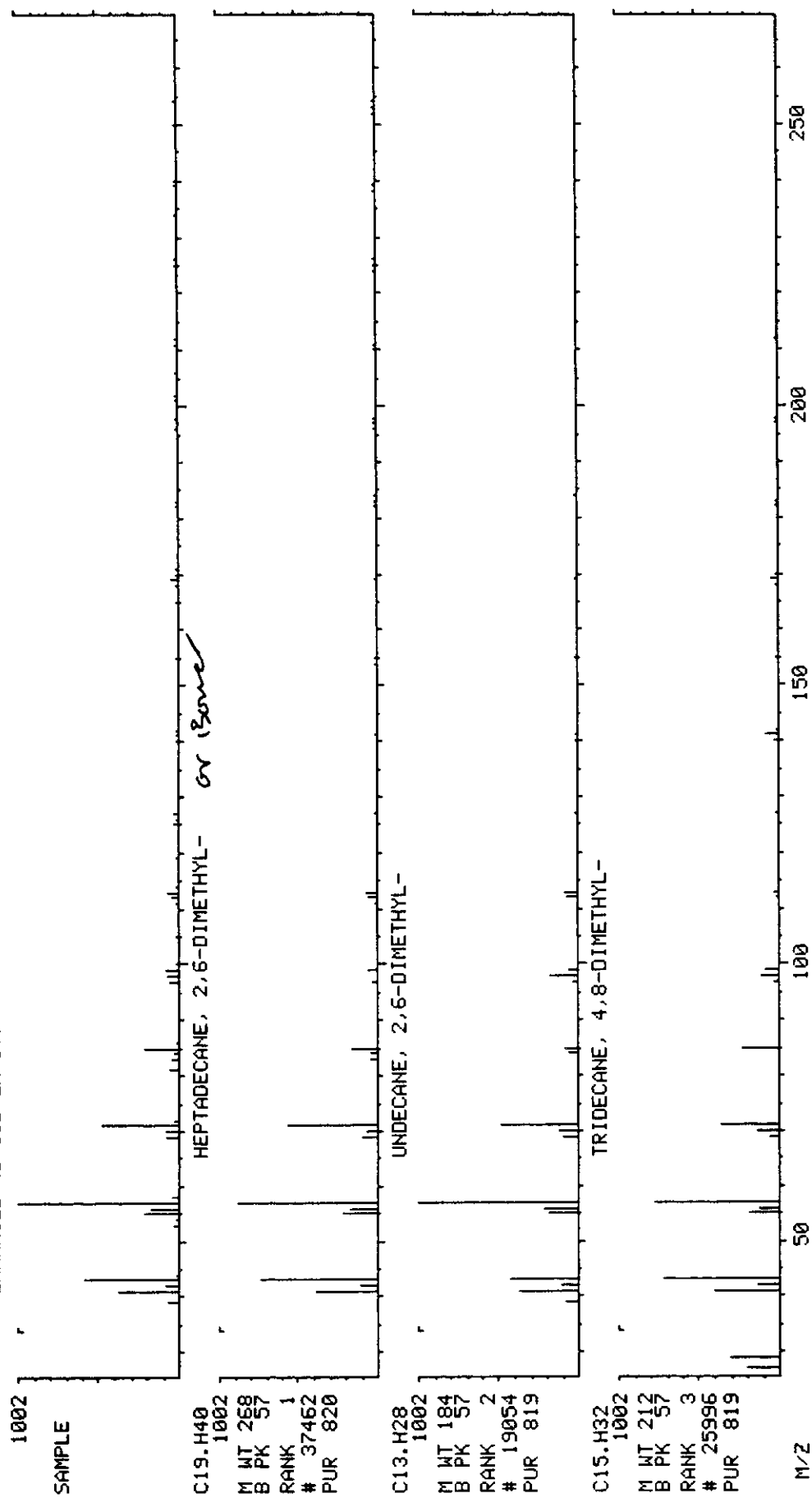
62231 spectra in LIBRARYNB searched for maximum PURITY
 644 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 19054 UNDECANE, 2,6-DIMETHYL-
 3 25996 TRIDECANE, 4,8-DIMETHYL-
 4 37465 NONADECANE
 5 25997 PENTADECANE
 6 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 7 22530 TETRADECANE
 8 22536 TRIDECANE, 3-METHYL-
 9 29263 HEXADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	820	949	832
2	C13.H28	184	57	819	949	826
3	C15.H32	212	57	819	945	826
4	C19.H40	268	57	816	935	841
5	C15.H32	212	57	809	951	817
6	C21.H44	296	57	809	948	820
7	C14.H30	198	43	807	940	816
8	C14.H30	198	57	801	961	801
9	C16.H34	226	57	800	928	815

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	17301-23-4
3	—	—	—	—	55030-62-1
4	—	—	—	—	629-92-5
5	—	—	—	—	629-62-9
6	—	—	—	—	18344-37-1
7	—	—	—	—	629-59-4
8	—	—	—	—	6418-41-3
9	—	—	—	—	544-76-3

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:01:00 + 11:00
 SAMPLE: T-MM5-2 1/35A/1ML
 COND.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30068103 # 989
 CALI: 30068103 # 3
 BASE M/Z: 57
 RIC: 133376.



Library Search Data: 30068103 #1037 Base m/z: 57
 08/31/98 20:01:00 + 11:32 Cali: 30068103 # 3 RIC: 209920.
 Sample: T-MMS-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 840 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 3 37465 NONADECANE
 4 26001 DODECANE, 2,7,10-TRIMETHYL-
 5 25991 DODECANE, 2,6,10-TRIMETHYL-
 6 22535 DODECANE, 4,6-DIMETHYL-
 7 37466 PENTADECANE, 2,6,10,14-TETRAMETHYL-
 8 25994 DODECANE, 2,6,11-TRIMETHYL-
 9 42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-

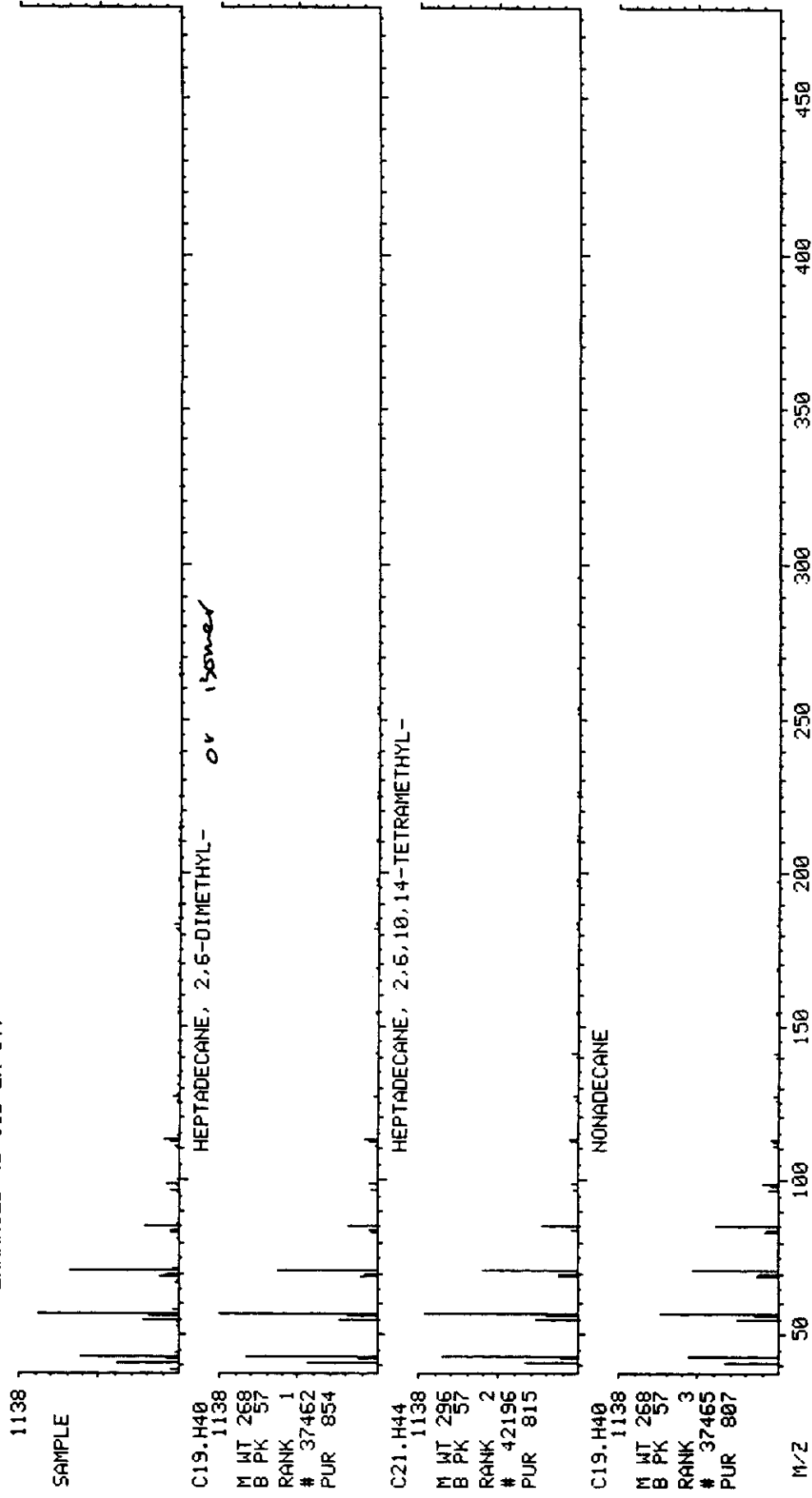
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	854	971	863
2	C21.H44	296	57	815	950	831
3	C19.H40	268	57	807	937	824
4	C15.H32	212	57	805	968	809
5	C15.H32	212	57	803	970	806
6	C14.H30	198	57	800	961	801
7	C19.H40	268	71	796	943	814
8	C15.H32	212	57	786	945	795
9	C21.H44	296	57	782	934	810

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	18344-37-1
3	—	—	—	—	629-92-5
4	—	—	—	—	74645-98-0
5	—	—	—	—	3891-98-3
6	—	—	—	—	61141-72-8
7	—	—	—	—	1921-70-6
8	—	—	—	—	31295-56-4
9	—	—	—	—	54833-48-6

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:01:00 + 11:32
 SAMPLE: T-MM5-2 1/35A/1ML INST. ID: F16
 COND5.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068103 #1037
 CALI: 30068103 # 3

BASE M/Z: 57
 RIC: 209920.



Library Search Data: 30068103 #1115 Base m/z: 57
 08/31/98 20:01:00 + 12:24 Cali: 30068103 # 3 RIC: 109696.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 766 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 15969 HYDROXYLAMINE, O-DECYL-
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 22536 TRIDECANE, 3-METHYL-
 4 11607 UNDECANE
 5 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 6 19035 DODECANE, 2-METHYL-
 7 37465 NONADECANE
 8 19022 UNDECANE, 4,8-DIMETHYL-
 9 18987 UNDECANE, 2,8-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C10.H23.O.N	173	43	767	924	803
2	C19.H40	268	57	757	888	824
3	C14.H30	198	57	749	910	779
4	C11.H24	156	43	743	945	764
5	C21.H44	296	57	743	902	804
6	C13.H28	184	43	737	946	756
7	C19.H40	268	57	736	884	819
8	C13.H28	184	43	736	925	740
9	C13.H28	184	43	729	931	742

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	29812-79-1
2	—	—	—	—	54105-67-8
3	—	—	—	—	6418-41-3
4	—	—	—	—	1120-21-4
5	—	—	—	—	18344-37-1
6	—	—	—	—	1560-97-0
7	—	—	—	—	629-92-5
8	—	—	—	—	17301-33-6
9	—	—	—	—	17301-25-6

BASE M/Z: 57
RIC: 109696.

DATA: 30058103 #1115
CALI: 30058103 # 3

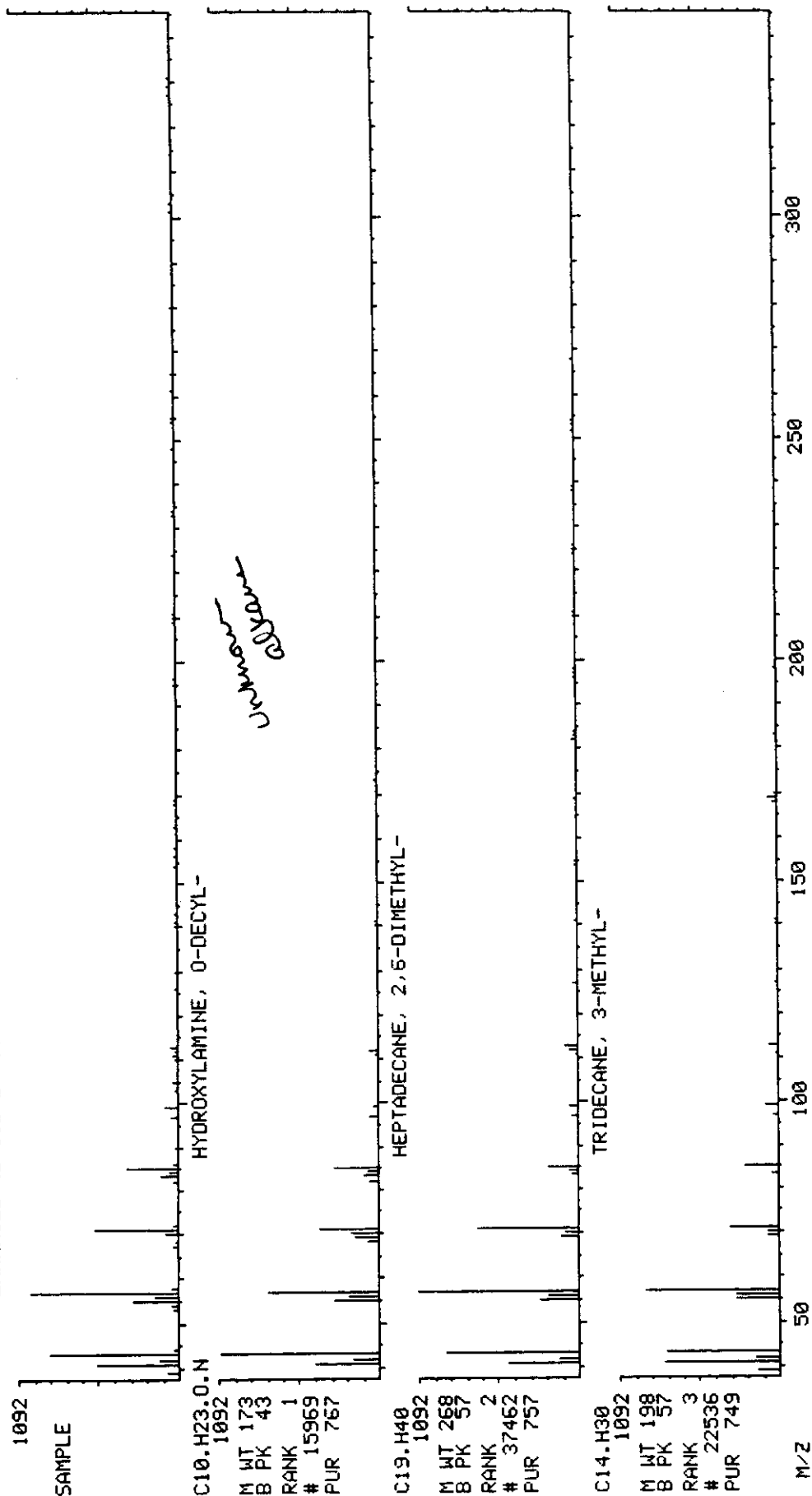
MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 20:01:00 + 12:24

SAMPLE: T-MM5-2 1/35A/1ML INST. ID: F16

CONDS.: UG/ML *1ML *100%/100% *(NA/NA >)/1/35A NA M

ENHANCED (S 15B 2N 0T)



8
CF
6)

Library Search Data: 30068103 #1120 Base m/z: 57
 08/31/98 20:01:00 + 12:27 Cali: 30068103 # 3 RIC: 144384.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

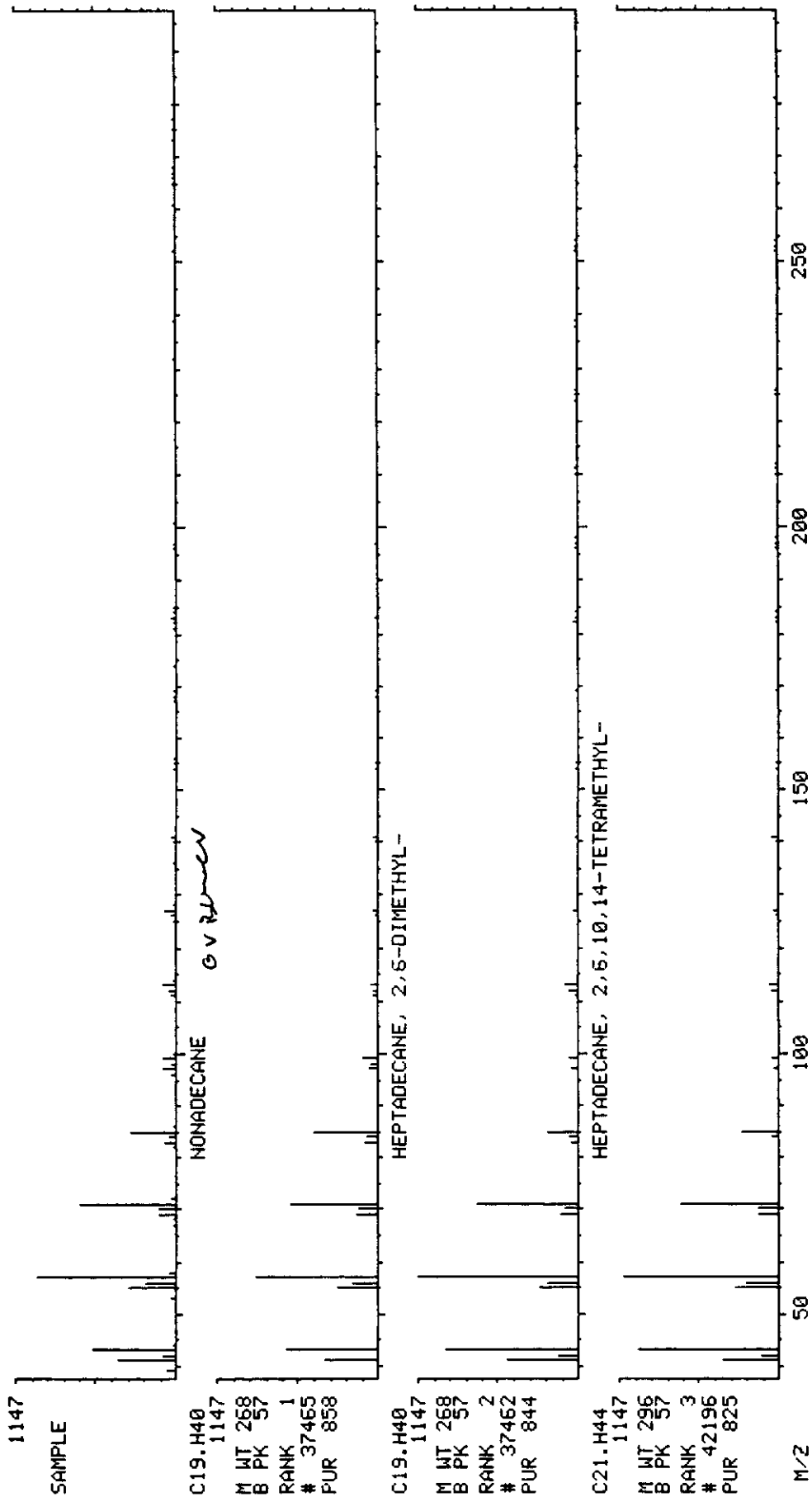
62231 spectra in LIBRARYNB searched for maximum PURITY
 774 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 46161 TRICOSANE
 5 42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-
 6 25991 DODECANE, 2,6,10-TRIMETHYL-
 7 26001 DODECANE, 2,7,10-TRIMETHYL-
 8 32059 HEPTADECANE
 9 37466 PENTADECANE, 2,6,10,14-TETRAMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	858	966	858
2	C19.H40	268	57	844	946	853
3	C21.H44	296	57	825	939	840
4	C23.H48	324	43	824	917	872
5	C21.H44	296	57	812	951	839
6	C15.H32	212	57	811	968	812
7	C15.H32	212	57	811	966	811
8	C17.H36	240	57	811	931	837
9	C19.H40	268	71	808	935	818

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	18344-37-1
4	—	—	—	—	638-67-5
5	—	—	—	—	54833-48-6
6	—	—	—	—	3891-98-3
7	—	—	—	—	74645-98-0
8	—	—	—	—	629-78-7
9	—	—	—	—	1921-70-6

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:01:00 + 12:27
 SAMPLE: T-MM5-2 1/35A/1ML
 COND.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30058103 #1120
 CALI: 30058103 # 3
 BASE M/2: 57
 RIC: 144384.



Library Search Data: 30068103 #1192 Base m/z: 57
 08/31/98 20:01:00 + 13:16 Cali: 30068103 # 3 RIC: 108288.
 Sample: T-MMS-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

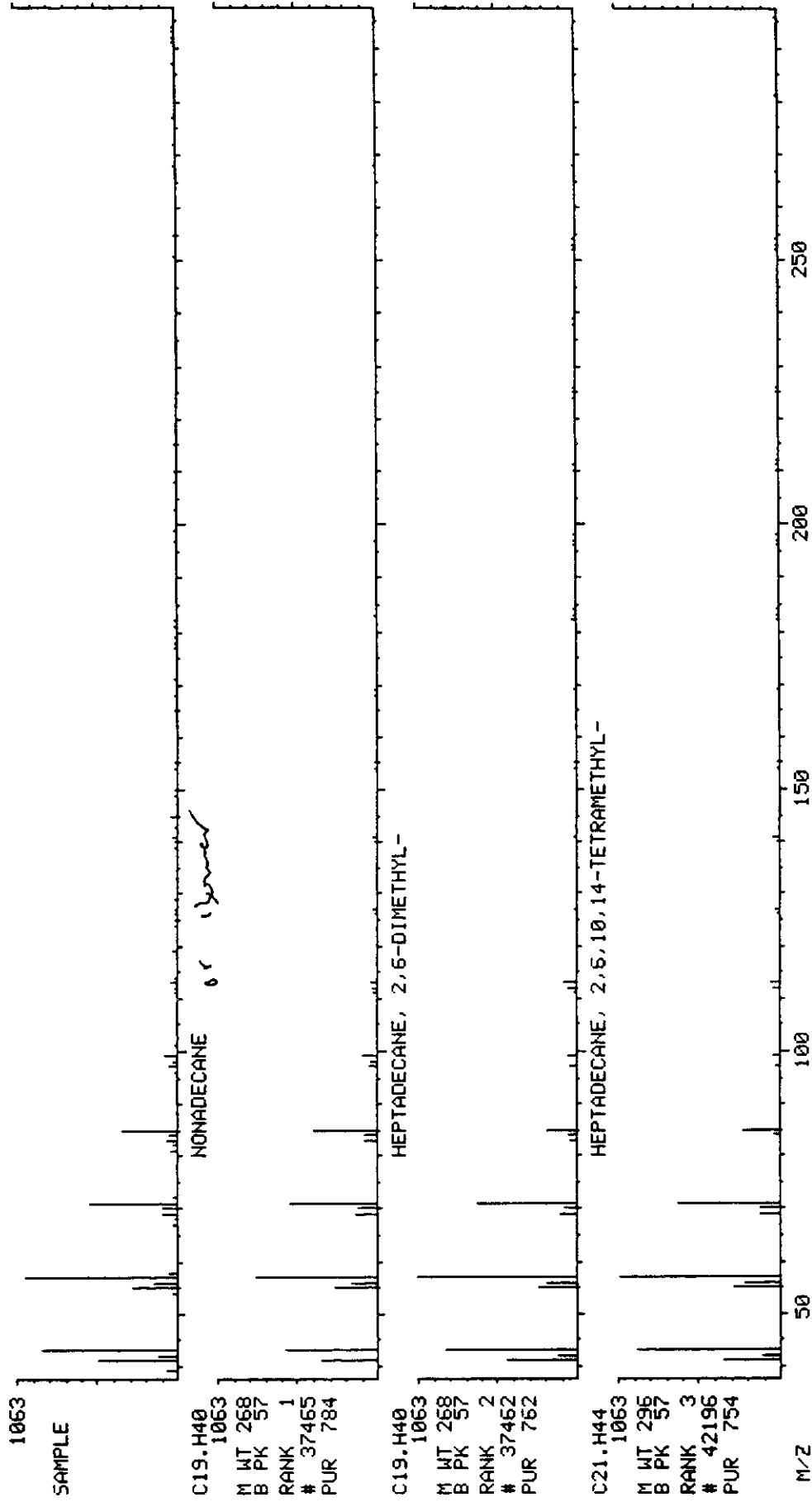
62231 spectra in LIBRARYNB searched for maximum PURITY
 294 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 15969 HYDROXYLAMINE, O-DECYL-
 5 25997 PENTADECANE
 6 19016 UNDECANE, 4,7-DIMETHYL-
 7 46161 TRICOSANE
 8 19523 1-DECANOL, 2-ETHYL-
 9 15353 2,6-DIMETHYLDECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	784	967	795
2	C19.H40	268	57	762	935	784
3	C21.H44	296	57	754	944	780
4	C10.H23.O.N	173	43	748	961	750
5	C15.H32	212	57	741	950	769
6	C13.H28	184	43	735	956	736
7	C23.H48	324	43	735	892	813
8	C12.H26.O	186	57	726	934	726
9	C12.H26	170	43	724	922	733

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	18344-37-1
4	—	—	—	—	29812-79-1
5	—	—	—	—	629-62-9
6	—	—	—	—	17301-32-5
7	—	—	—	—	638-67-5
8	—	—	—	—	21078-65-9
9	—	—	—	—	13150-81-7

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:01:00 + 13:16
 SAMPLE: T-MM5-2 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA) 1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30068103 #1192
 CALI: 30068103 # 3
 BASE M/Z: 57
 RIC: 108288.



Library Search Data: 30068103 #1266 Base m/z: 57
 08/31/98 20:01:00 + 14:05 Cali: 30068103 # 3 RIC: 91520.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 883 matched at least 7 of the 16 largest peaks in the unknown

Rank In.	Name
1 42196	HEPTADECANE, 2,6,10,14-TETRAMETHYL-
2 37465	NONADECANE
3 25997	PENTADECANE
4 37462	HEPTADECANE, 2,6-DIMETHYL-
5 32058	HEXADECANE, 3-METHYL-
6 39858	EICOSANE
7 46161	TRICOSANE
8 42192	HEPTADECANE, 2,6,10,15-TETRAMETHYL-
9 49555	PENTACOSANE

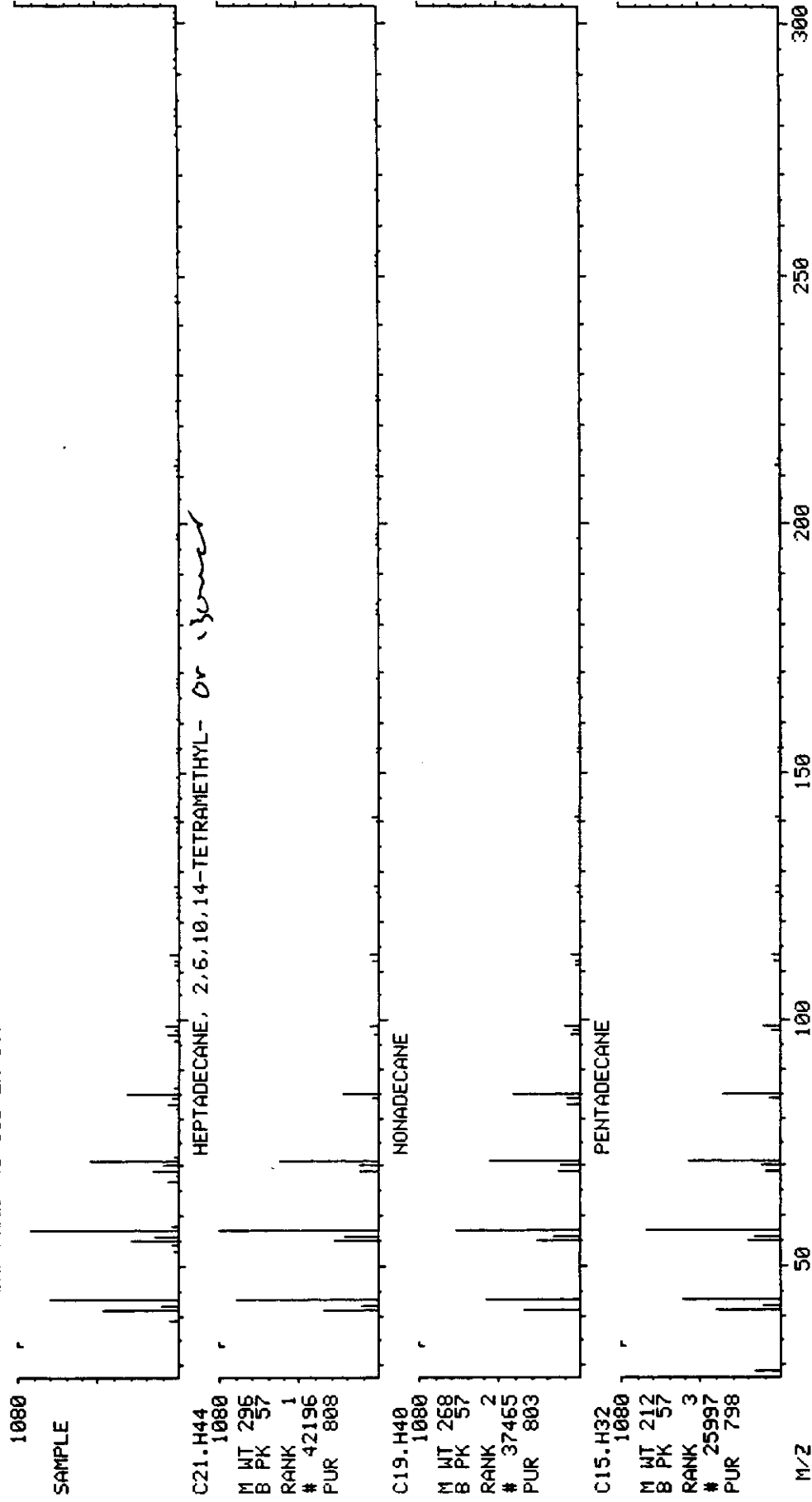
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C21.H44	296	57	808	959	815
2	C19.H40	268	57	803	971	816
3	C15.H32	212	57	798	988	798
4	C19.H40	268	57	797	950	807
5	C17.H36	240	57	785	953	789
6	C20.H42	282	57	785	921	809
7	C23.H48	324	43	784	929	837
8	C21.H44	296	57	768	939	790
9	C25.H52	352	43	765	903	837

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	18344-37-1
2	—	—	—	—	629-92-5
3	—	—	—	—	629-62-9
4	—	—	—	—	54105-67-8
5	—	—	—	—	6418-43-5
6	—	—	—	—	112-95-8
7	—	—	—	—	638-67-5
8	—	—	—	—	54833-48-6
9	—	—	—	—	629-99-2

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:01:00 + 14:05
 SAMPLE: T-MMS-2 1/35A/1ML
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068103 #1266
 CALI: 30068103 # 3

BASE M/Z: 57
 RIC: 91520.



Library Search Data: 30068103 #1336 Base m/z: 57
 08/31/98 20:01:00 + 14:52 Cali: 30068103 # 3 RIC: 80896.
 Sample: T-MMS-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 743 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 37465 NONADECANE
 3 37456 2-METHYLOCTADECANE
 4 29263 HEXADECANE
 5 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 6 46161 TRICOSANE
 7 32058 HEXADECANE, 3-METHYL-
 8 19523 1-DECANOL, 2-ETHYL-
 9 15969 HYDROXYLAMINE, O-DECYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	698	888	727
2	C19.H40	268	57	697	954	723
3	C19.H40	268	43	697	918	722
4	C16.H34	226	57	690	942	712
5	C21.H44	296	57	688	907	710
6	C23.H48	324	43	682	907	739
7	C17.H36	240	57	674	929	693
8	C12.H26.O	186	57	672	949	672
9	C10.H23.O.N	173	43	672	962	675

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	54105-67-8
2	---	---	---	---	629-92-5
3	---	---	---	---	- -
4	---	---	---	---	544-76-3
5	---	---	---	---	18344-37-1
6	---	---	---	---	638-67-5
7	---	---	---	---	6418-43-5
8	---	---	---	---	21078-65-9
9	---	---	---	---	29812-79-1

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 20:01:00 + 14:52

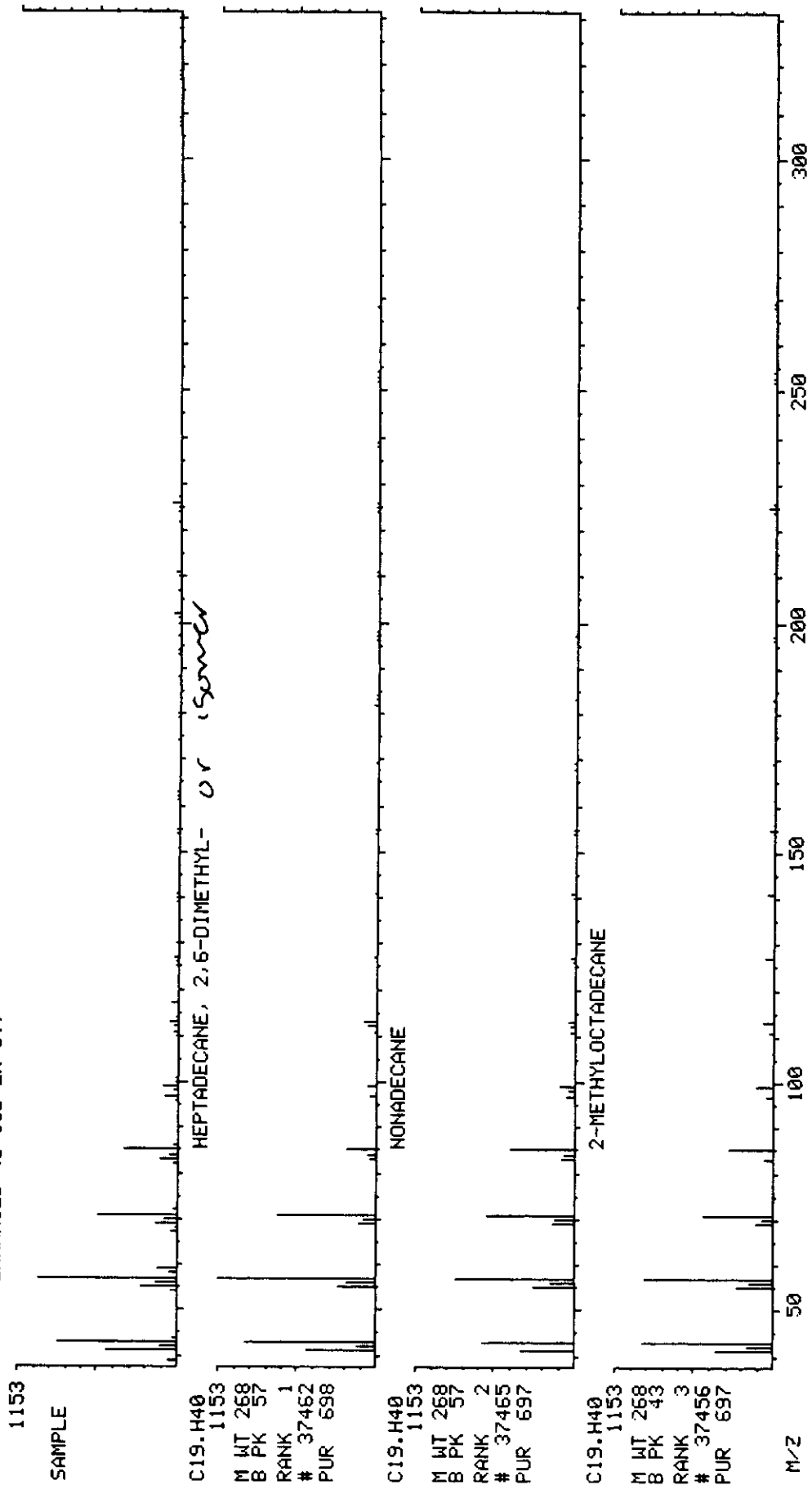
SAMPLE: T-MM5-2 1/35A/1ML

INST. ID: F16

CONDS.: UG/ML *1ML *1002/1002 *(NA/NA)/1/35A NA M

ENHANCED (S 15B 2N 0T)

DATA: 30068103 #1335
CALI: 30068103 # 3
BASE M/Z: 57
RIC: 80896.



Library Search Data: 30068103 #1386 Base m/z: 244
 08/31/98 20:01:00 + 15:25 Cali: 30068103 # 3 RIC: 87808.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 801 matched at least 3 of the 16 largest peaks in the unknown

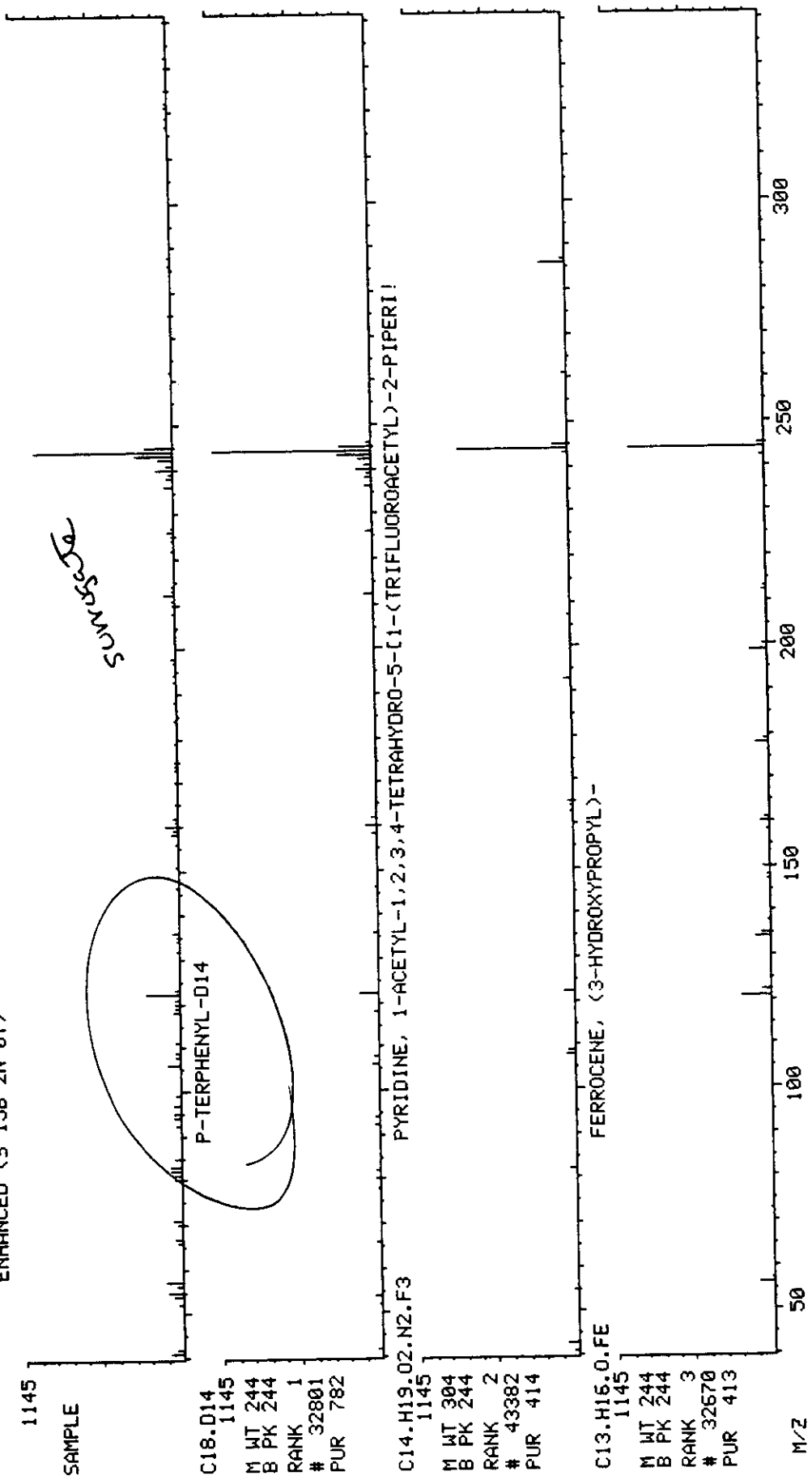
Rank In. Name
 1 32801 P-TERPHENYL-D14
 2 43382 PYRIDINE, 1-ACETYL-1,2,3,4-TETRAHYDRO-5-[1-(TRIFLUOROACETYL)-2-PIPE*
 3 32670 FERROCENE, (3-HYDROXYPROPYL)-
 4 32765 MATRIDIN-15-ONE, 11,12,13,14-TETRADEHYDRO-, (6.BETA.)-
 5 32768 ANAGYRINE
 6 48867 SULFIDE, PENTYL TRITYL
 7 50992 SULFIDE, BENZYL TRITYL
 8 32560 9H-CARBAZOLE, 9-PHENYL-
 9 40723 ETHYLAMINE, N-TRITYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C18.D14	244	244	782	976	791
2	C14.H19.O2.N2.F3	304	244	414	732	442
3	C13.H16.O.FE	244	244	413	742	482
4	C15.H20.O.N2	244	243	401	679	519
5	C15.H20.O.N2	244	98	367	614	542
6	C24.H26.S	346	243	365	579	410
7	C26.H22.S	366	243	363	554	448
8	C18.H13.N	243	243	361	553	394
9	C21.H21.N	287	244	360	630	481

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	1718-51-0
2	—	—	—	—	54966-10-8
3	—	—	—	—	12093-88-8
4	—	—	—	—	6838-34-2
5	—	—	—	—	486-89-5
6	—	—	—	—	20705-42-4
7	—	—	—	—	6622-14-6
8	—	—	—	—	1150-62-5
9	—	—	—	—	7370-34-5

DATA: 30068103 #1386
CALI: 30068103 # 3
BASE M/Z: 244
RIC: 87808.

MTD LIBRARY SEARCH (LIBRARYNB)
08/31/98 20:01:00 + 15:25
SAMPLE: T-MM5-2 1/35A/1ML INST. ID: F16
COND5.: UG/ML *1ML *100%/100% *(NA/NA) >1/35A NA M
ENHANCED (S 15B 2N 0T)



Library Search Data: 30068103 #1404 Base m/z: 57
 08/31/98 20:01:00 + 15:37 Cali: 30068103 # 3 RIC: 62528.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 793 matched at least 7 of the 16 largest peaks in the unknown

Rank In.	Name
1	37465 NONADECANE
2	37462 HEPTADECANE, 2,6-DIMETHYL-
3	18985 TRIDECANE
4	25997 PENTADECANE
5	42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-
6	19016 UNDECANE, 4,7-DIMETHYL-
7	15969 HYDROXYLAMINE, O-DECYL-
8	25994 DODECANE, 2,6,11-TRIMETHYL-
9	22530 TETRADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	695	931	741
2	C19.H40	268	57	688	912	729
3	C13.H28	184	57	687	970	689
4	C15.H32	212	57	686	953	714
5	C21.H44	296	57	684	924	723
6	C13.H28	184	43	684	951	684
7	C10.H23.O.N	173	43	681	950	697
8	C15.H32	212	57	680	956	685
9	C14.H30	198	43	680	938	714

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	629-50-5
4	—	—	—	—	629-62-9
5	—	—	—	—	54833-48-6
6	—	—	—	—	17301-32-5
7	—	—	—	—	29812-79-1
8	—	—	—	—	31295-56-4
9	—	—	—	—	629-59-4

DATA: 30068103 #1404
CALI: 30068103 # 3
BASE M/Z: 57
RIC: 62528.

MID LIBRARY SEARCH (LIBRARYNB)

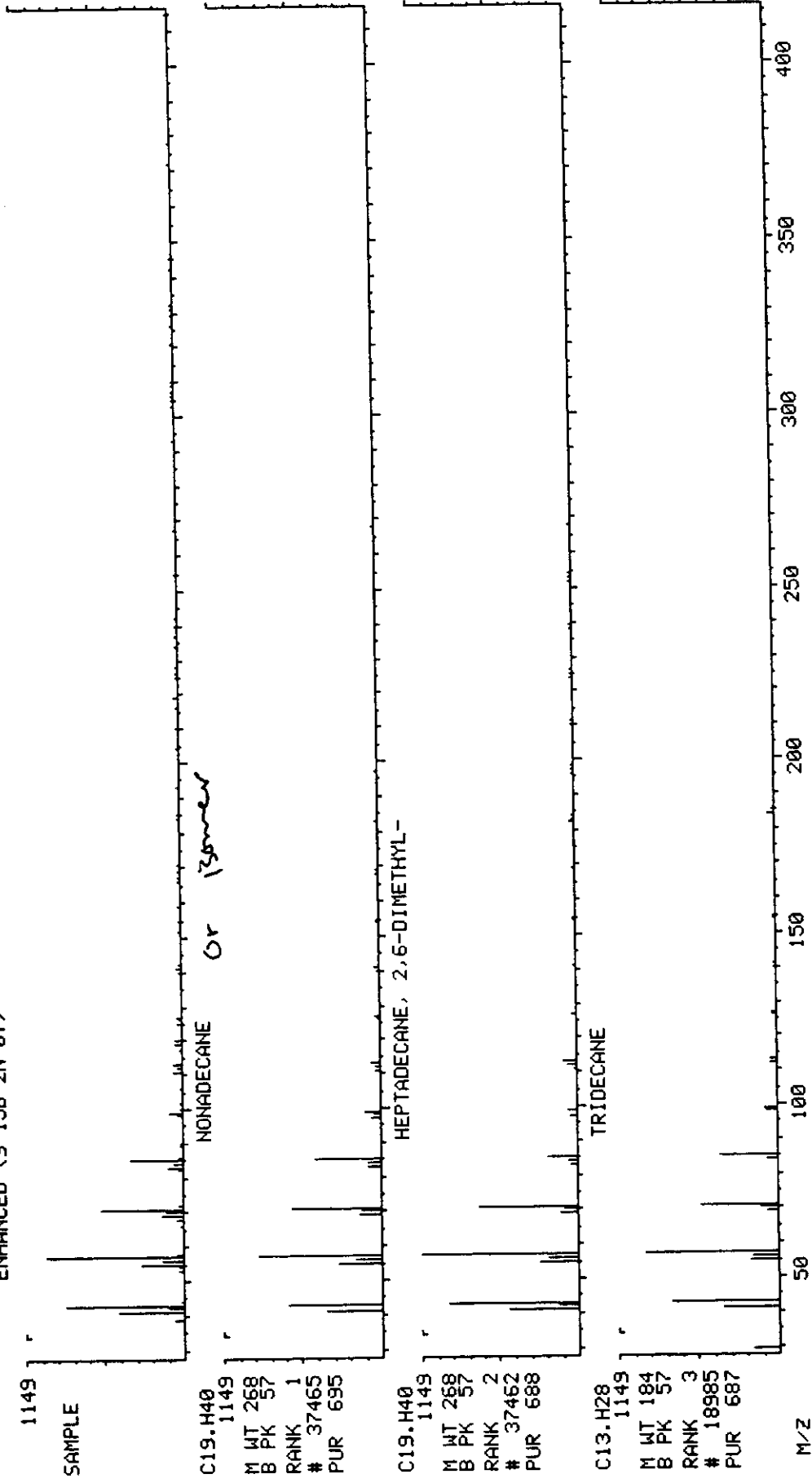
08/31/98 20:01:00 + 15:37

SAMPLE: T-MM5-2 1/35A/1ML

COND.: UG/ML *1ML *100Z/100Z *(NA/NA) /1/35A NA M

ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068103 #1531 Base m/z: 57
 08/31/98 20:01:00 + 17:02 Cali: 30068103 # 3 RIC: 41152.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 783 matched at least 7 of the 16 largest peaks in the unknown

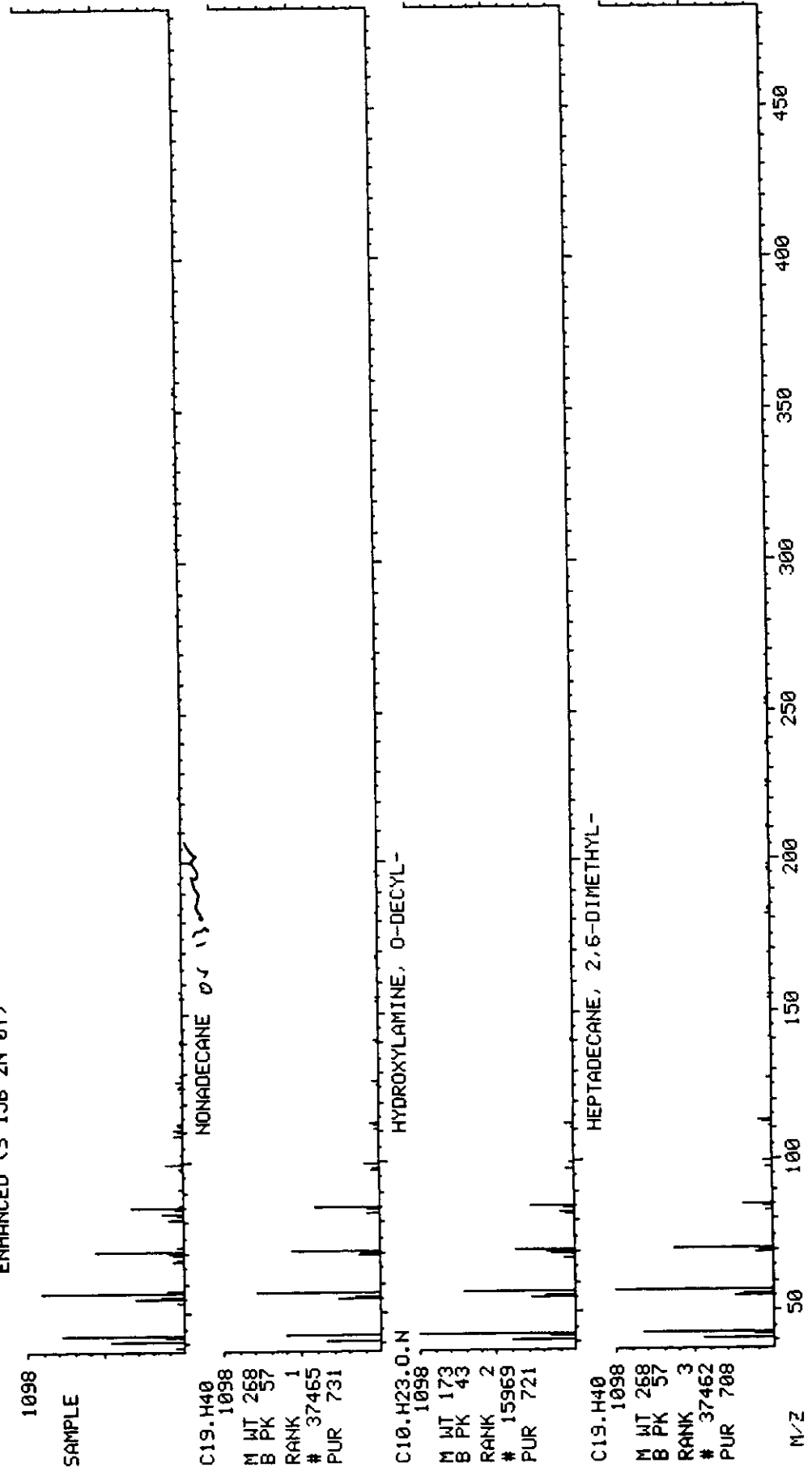
Rank In. Name
 1 37465 NONADECANE
 2 15969 HYDROXYLAMINE, O-DECYL-
 3 37462 HEPTADECANE, 2,6-DIMETHYL-
 4 25997 PENTADECANE
 5 19523 1-DECANOL, 2-ETHYL-
 6 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 7 19016 UNDECANE, 4,7-DIMETHYL-
 8 19006 DECANE, 6-ETHYL-2-METHYL-
 9 22530 TETRADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	731	945	759
2	C10.H23.O.N	173	43	721	957	723
3	C19.H40	268	57	708	924	735
4	C15.H32	212	57	698	949	726
5	C12.H26.O	186	57	698	929	702
6	C21.H44	296	57	694	948	715
7	C13.H28	184	43	692	953	693
8	C13.H28	184	57	691	948	697
9	C14.H30	198	43	688	935	725

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	29812-79-1
3	—	—	—	—	54105-67-8
4	—	—	—	—	629-62-9
5	—	—	—	—	21078-65-9
6	—	—	—	—	18344-37-1
7	—	—	—	—	17301-32-5
8	—	—	—	—	62108-21-8
9	—	—	—	—	629-59-4

DATA: 30068103 #1531
CALI: 30068103 # 3
BASE M/Z: 57
RIC: 41152.

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 20:01:00 + 17:02
SAMPLE: T-MM5-2 1/35A/1ML INST. ID: F16
COND5.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)



Library Search Data: 30068103 #1587 Base m/z: 43
 08/31/98 20:01:00 + 17:39 Cali: 30068103 # 3 RIC: 68864.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 950 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 34410 3-OCTADECENE, (E)-
 2 39515 9-EICOSENE, (E)-
 3 39516 5-EICOSENE, (E)-
 4 34418 9-OCTADECENE, (E)-
 5 34411 5-OCTADECENE, (E)-
 6 31653 1-HEPTADECENE
 7 39517 3-EICOSENE, (E)-
 8 28772 7-HEXADECENE, (Z)-
 9 32420 1-HEXADECANOL

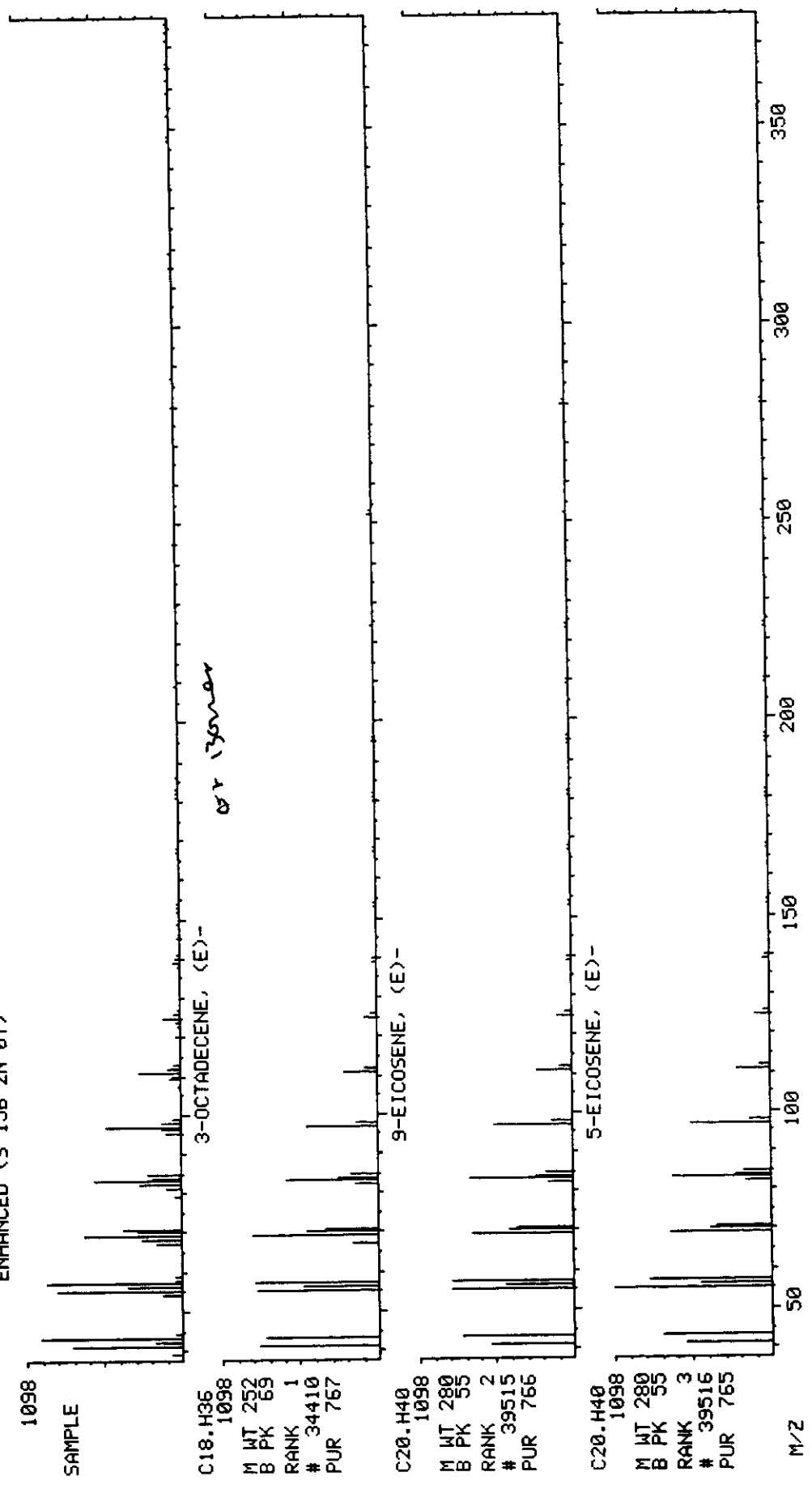
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C18.H36	252	69	767	968	782
2	C20.H40	280	55	766	979	770
3	C20.H40	280	55	765	978	771
4	C18.H36	252	55	763	963	777
5	C18.H36	252	55	762	962	777
6	C17.H34	238	55	756	970	770
7	C20.H40	280	57	755	982	760
8	C16.H32	224	55	749	945	766
9	C16.H34.O	242	55	748	973	758

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	7206-19-1
2	—	—	—	—	74685-29-3
3	—	—	—	—	74685-30-6
4	—	—	—	—	7206-25-9
5	—	—	—	—	7206-21-5
6	—	—	—	—	6765-39-5
7	—	—	—	—	74685-33-9
8	—	—	—	—	35507-09-6
9	—	—	—	—	36653-82-4

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:01:00 + 17:39
 SAMPLE: T-NMS-2 1/35A/1ML
 COND.: UG/ML *1ML *100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068103 #1587
 CALI: 30068103 # 3

BASE M/Z: 43
 RIC: 58854.



08
 11
 21

Library Search Data: 30068103 #1809 Base m/z: 43
 08/31/98 20:01:00 + 20:07 Cali: 30068103 # 3 RIC: 20672.
 Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 761 matched at least 5 of the 16 largest peaks in the unknown

Rank In. Name
 1 52672 SILANE, TRICHLOROCTADECYL-
 2 51637 CHOLESTANE
 3 31486 ACETAMIDE, N-METHYL-N-[4-[4-METHOXY-1-HEXAHYDROPYRIDYL]-2-BUTYNYL]-
 4 40940 PREGNANE
 5 54646 SQUALENE
 6 50599 PENTALENE, OCTAHYDRO-1-(2-OCTYLDECYL)-
 7 46251 DODECANE, 1,2-DIBROMO-
 8 40165 HEXADECANOIC ACID, 2-OXO-, METHYL ESTER
 9 48865 5.ALPHA.-ANDROSTAN-3.BETA.-OL, 4,4-DIMETHYL-, ACETATE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C18.H37.CL3.S1	386	57	342	755	375
2	C27.H48	372	43	316	625	402
3	C13.H22.O2.N2	238	43	305	813	348
4	C21.H36	288	55	300	677	376
5	C30.H50	410	69	294	687	347
6	C26.H50	362	43	292	863	320
7	C12.H24.BR2	326	41	290	770	345
8	C17.H32.O3	284	57	286	817	337
9	C23.H38.O2	346	43	283	672	393

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	112-04-9
2	—	—	—	—	481-21-0
3	—	—	—	—	- -
4	—	—	—	—	481-26-5
5	—	—	—	—	7683-64-9
6	—	—	—	—	55401-65-5
7	—	—	—	—	55334-42-4
8	—	—	—	—	55836-30-1
9	—	—	—	—	7761-08-2

DATA: 30068103 #1809
CALI: 30068103 # 3

BASE M/Z: 43
RIC: 20672.

MID LIBRARY SEARCH <LIBRARY>

08/31/98 20:01:00 + 20:07

SAMPLE: T-MM5-2 1/35A/1ML

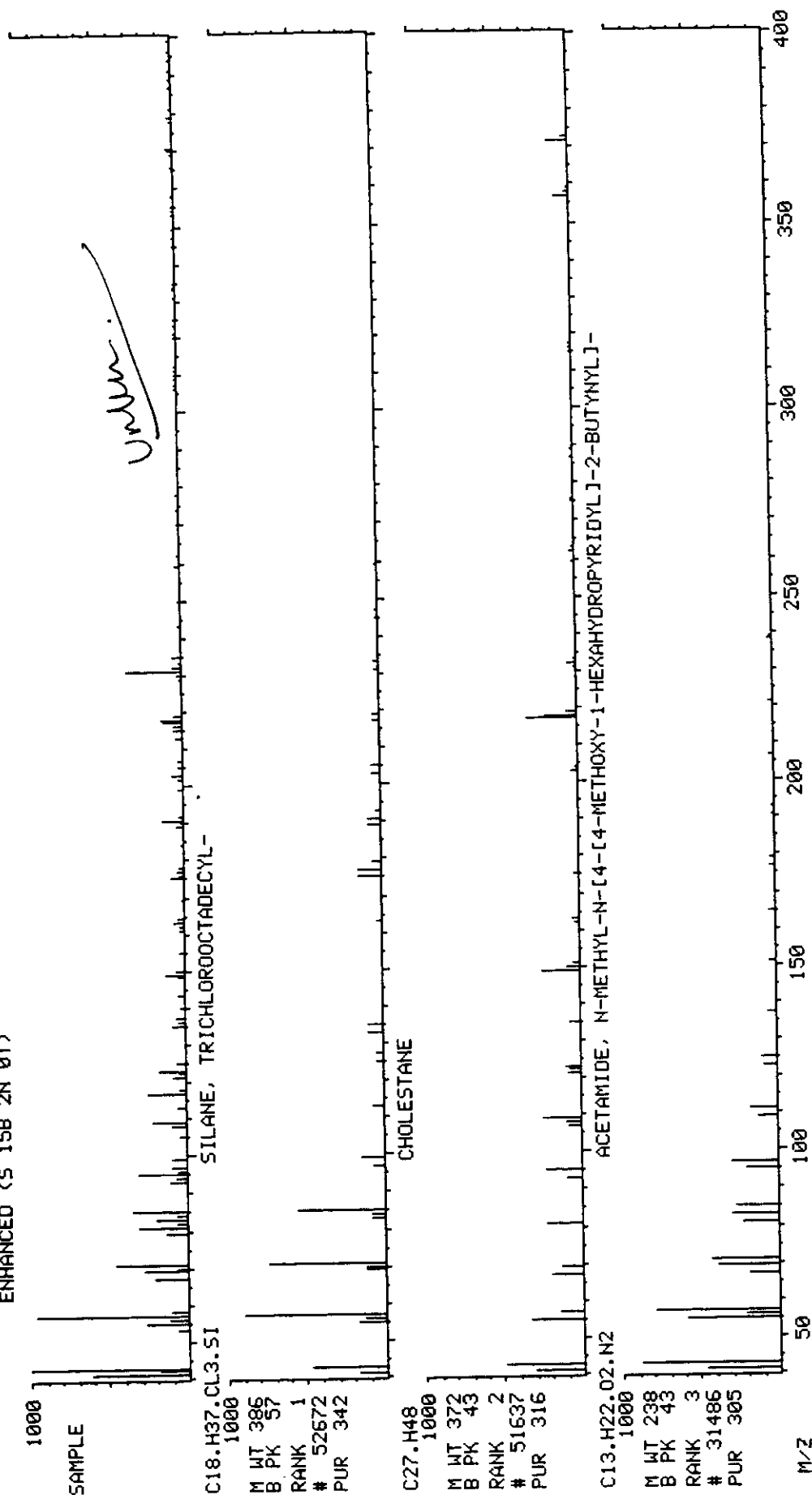
COND.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M

ENHANCED <S 158 2N 0T>

INST. ID: F16

NA M

ENHANCED <S 158 2N 0T>



C
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TIC SELECTION REPORT

DATA FILE: 30068103

THE FOLLOWING PEAKS WERE REJECTED BECAUSE
AT LEAST 40 % OF THEIR SIZE WAS ACCOUNTED FOR BY
TARGET COMPOUNDS ELUTING WITHIN 4 SCANS OF THE
PEAK TOP.

SCAN	SIZE	AMOUNT
199	205984.	24.199
310	455754.	53.543
341	148864.	17.489
398	55273.	6.494
402	58805.	6.908
422	191616.	22.511
531	377600.	44.361
724	383252.	45.025
830	340480.	40.000
1088	342016.	40.000
1554	317184.	40.000
1785	244032.	40.000

TOTAL NUMBER OF UNIDENTIFIED PEAKS WITH SIZE
GREATER THAN 10 % OF THE CLOSEST INTERNAL STANDARD
THAT DOES NOT HAVE INTERFERENCES = 121

INTERNAL STANDARDS THAT HAVE RIC SIZE LESS THAN
50 % OR GREATER THAN 200 % OF THE ESTIMATED RIC SIZE
ARE CONSIDERED TO HAVE INTERFERENCES AND WILL NOT BE USED
FOR QUANTITATION.

#	INTERNAL STANDARD	RIC SIZE	PERCENT OF ESTIMATED RIC SIZE	SCAN
1	C140 NAPHTHALENE-D8	377600.	212 *	531
2	C150 ACENAPHTHENE-D10	340480.	162	830
3	C160 PHENANTHRENE-D10	342016.	156	1088
4	C170 CHRYSENE-D12	317184.	140	1554
5	C175 PERYLENE-D12	244032.	106	1785

* INDICATES INTERFERENCE

SIZE = AREA

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 300681RI03
Std Id: ST16980902

Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 09/02/98 19:59
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Surrogate Spike Recoveries
8270-G, LIMS 10/10/96

Surrogate	Surrogate Amount (ug)		% Recovery	
	Spiked	Measured	Measured	QC limits
CS20 NITROBENZENE-D5	50.00	40.80	81.6	45 107
CS25 2-FLUOROBIPHENY	50.00	53.80	108.	62 110
CS30 TERPHENYL-D14	50.00	39.50	79.0	58 135
CS45 PHENOL-D5	100.0	64.50	64.5	43 130
CS50 2-FLUOROPHENOL	100.0	58.30	58.3	36 111
CS55 2,4,6-TRIBROMOP	100.0	45.00	45.0	* 58 131

Target Compounds: SAP9

Confirms
Reporting
Limit

Parameter	Scan	Result	Units	Limit
CS70 2-CHLOROPHENOL-D4		ND	UG/A	30.0
CS75 1,2-DICHLOROBNZN-D4	342	130.	UG/A	30.0
HEXACHLOROBNZN-D13		ND	UG/A	30.0
C310 N-NITROSODIMETHYLAMINE		ND	UG/A	30.0
PYRIDINE		ND	UG/A	60.0
2-PICOLINE		ND	UG/A	30.0
N-NITROSOMETHYLETHYLAMINE		ND	UG/A	30.0
METHYLMETHANESULFONATE		ND	UG/A	30.0
N-NITROSODIETHYLAMINE		ND	UG/A	30.0
ETHYLMETHANE SULFONATE		ND	UG/A	30.0
PENTACHLOROETHANE		ND	UG/A	30.0
C320 ANILINE		ND	UG/A	30.0
C315 PHENOL		ND	UG/A	30.0
C325 BIS(2-CHLOROETHYL) ETHER		ND	UG/A	30.0
C330 2-CHLOROPHENOL		ND	UG/A	30.0
C335 1,3-DICHLOROBENZENE		ND	UG/A	30.0
C340 1,4-DICHLOROBENZENE		ND	UG/A	30.0
C345 BENZYL ALCOHOL		ND	UG/A	30.0
C350 1,2-DICHLOROBENZENE		ND	UG/A	30.0
C355 2-METHYLPHENOL		ND	UG/A	30.0
C360 2,2'-OXYBIS(1-CLPROPAN)		ND	UG/A	30.0
C361 ACETOPHENONE		ND	UG/A	30.0
N-NITROSOPYRROLIDINE		ND	UG/A	30.0
N-NITROSOMORPHOLINE		ND	UG/A	30.0
3-METHYL PHENOL		ND	UG/A	30.0
C365 4-METHYLPHENOL		ND	UG/A	30.0
C370 N-NITROSO-DI-N-PROPYLAM		ND	UG/A	30.0
O-TOLUIDINE		ND	UG/A	30.0
C375 HEXACHLOROETHANE		ND	UG/A	30.0

876

Reviewed by: _____

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 300681RI03
Std Id: ST16980902

Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 09/02/98 19:59
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
=====				
C410 NITROBENZENE		ND	UG/A	30.0
N-NITROSOPIPERIDINE		ND	UG/A	30.0
C415 ISOPHORONE		ND	UG/A	30.0
C420 2-NITROPHENOL		ND	UG/A	30.0
C425 2,4-DIMETHYLPHENOL		ND	UG/A	30.0
C435 BIS(2-CHLOROETHOXY)METH		ND	UG/A	30.0
C440 2,4-DICHLOROPHENOL		ND	UG/A	30.0
C430 BENZOIC ACID		ND	UG/A	150.0
AA-DIMETHYLPHENETHYLAMINE		ND	UG/A	30.0
C445 1,2,4-TRICHLOROBENZENE		ND	UG/A	30.0
C450 NAPHTHALENE		ND	UG/A	30.0
C455 4-CHLOROANALINE		ND	UG/A	30.0
2,6-DICHLOROPHENOL		ND	UG/A	30.0
HEXACHLOROPROPENE		ND	UG/A	30.0
C460 HEXACHLOROBUTADIENE		ND	UG/A	30.0
P-PHENYLENE DIAMINE		ND	UG/A	30.0
N-NITROSODI-N-BUTYLAMINE		ND	UG/A	30.0
C465 4-CHLORO-3-METHYLPHENO		ND	UG/A	30.0
SAFROLE		ND	UG/A	30.0
C470 2-METHYLNAPHTHALENE	654	60.	UG/A	30.0
1,2,4,5-TETRACHLOROBENZENE		ND	UG/A	30.0
ISOSAFROLE (#1)		ND	UG/A	60.0
C510 HEXACHLOROCYCLOPENTADI		ND	UG/A	30.0
C515 2,4,6-TRICHLOROPHENOL		ND	UG/A	30.0
C520 2,4,5-TRICHLOROPHENOL		ND	UG/A	30.0
ISOSAFROLE (#2)		ND	UG/A	60.0
C525 2-CHLORONAPHTHALENE		ND	UG/A	30.0
1-CHLORONAPHTHALENE		ND	UG/A	15.0
C530 2-NITROANALINE		ND	UG/A	150.0
1,4-NAPHTHOQUINONE		ND	UG/A	30.0
C535 DIMETHYLPHTHALATE		ND	UG/A	30.0
1,3-DINITROBENZENE		ND	UG/A	30.0
C540 ACENAPHTHYLENE		ND	UG/A	30.0
C543 2,6-DINITROTOLUENE		ND	UG/A	30.0
C545 3-NITROANILINE		ND	UG/A	150.0
C550 ACENAPHTHENE		ND	UG/A	30.0
C555 2,4-DINITROPHENOL		ND	UG/A	150.0
C565 DIBENZOFURAN		ND	UG/A	30.0
C560 4-NITROPHENOL		ND	UG/A	150.0
PENTACHLOROBENZENE		ND	UG/A	30.0
C570 2,4-DINITROTOLUENE		ND	UG/A	30.0
1-NAPHTHYLAMINE		ND	UG/A	30.0
2-NAPHTHYLAMINE		ND	UG/A	30.0
2,3,4,6-TETRACHLOROPHENOL		ND	UG/A	60.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 300681RI03
Std Id: ST16980902

Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 09/02/98 19:59
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
=====				
C580 DIETHYLPHTHALATE		ND	UG/A	30.0
C590 FLUORENE		ND	UG/A	30.0
C585 4-CHLOROPHENYL-PHENYLE		ND	UG/A	30.0
5-NITRO-O-TOLUIDINE		ND	UG/A	30.0
C595 4-NITROANALINE		ND	UG/A	150.0
C610 4,6-DINITRO-2-METHYLPH		ND	UG/A	150.0
C615 N-NITROSODIPHENYLAMINE		ND	UG/A	30.0
C620 AZOBENZENE		ND	UG/A	30.0
SYM-TRINITROBENZENE		ND	UG/A	30.0
C625 4-BROMOPHENYL-PHENYLET		ND	UG/A	30.0
PHENACETIN		ND	UG/A	30.0
DIALATE	234	ND	UG/A	30.0
C630 HEXACHLOROBENZENE		ND	UG/A	30.0
4-AMINOBIIPHENYL		ND	UG/A	30.0
C635 PENTACHLOROPHENOL		ND	UG/A	150.0
PRONAMIDE		ND	UG/A	30.0
PENTACHLORONITROBENZENE		ND	UG/A	150.0
C640 PHENANTHRENE		ND	UG/A	30.0
C645 ANTHRACENE		ND	UG/A	30.0
2SECBUTYL-4,6-DINITROPHENOL		ND	UG/A	30.0
C647 CARBAZOLE		ND	UG/A	30.0
C650 DI-N-BUTYLPHTHALATE		ND	UG/A	30.0
4-NITROQUINOLINE-1-OXIDE		ND	UG/A	30.0
METHAPYRILENE		ND	UG/A	30.0
ISODRIN		ND	UG/A	30.0
C655 FLUORANTHENE		ND	UG/A	30.0
CHLOROBENZILATE		ND	UG/A	30.0
C710 BENZIDINE		ND	UG/A	300.0
C715 PYRENE		ND	UG/A	30.0
ARAMITE (#1)		ND	UG/A	30.0
ARAMITE (#2)		ND	UG/A	30.0
P-DIMETHYLAMINOAZOBENZENE		ND	UG/A	30.0
3,3'-DIMETHYLBENZIDINE		ND	UG/A	30.0
KEPONE		ND	UG/A	150.0
C720 BUTYLBENZYLPHTHALATE		ND	UG/A	30.0
2-ACETYLAMINOFLUORENE		ND	UG/A	30.0
C730 BENZO (A) ANTHRACENE		ND	UG/A	30.0
C725 3,3'-DICHLOROBENZIDINE		ND	UG/A	60.0
C740 CHRYSENE		ND	UG/A	30.0
C745 BIS (2-ETHYLHEXYL) PHTHA		ND	UG/A	30.0
3-METHYLCHOLANTHRENE		ND	UG/A	30.0
C760 DI-N-OCTYL PHTHALATE		ND	UG/A	30.0
C765 BENZO (B) FLUORANTHENE		ND	UG/A	30.0
7,12-DIMETHYLBENZANTHRACENE		ND	UG/A	30.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 300681RI03
Std Id: ST16980902

Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 09/02/98 19:59
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C770 BENZO (K) FLUORANTHENE		ND	UG/A	30.0
HEXACHLOROPHENE		ND	UG/A	30.0
C775 BENZO (A) PYRENE		ND	UG/A	30.0
C780 INDENO (1,2,3-CD) PYRENE		ND	UG/A	30.0
C785 DIBENZ (A,H) ANTHRACENE		ND	UG/A	30.0
C790 BENZO (G,H,I) PERYLENE		ND	UG/A	30.0

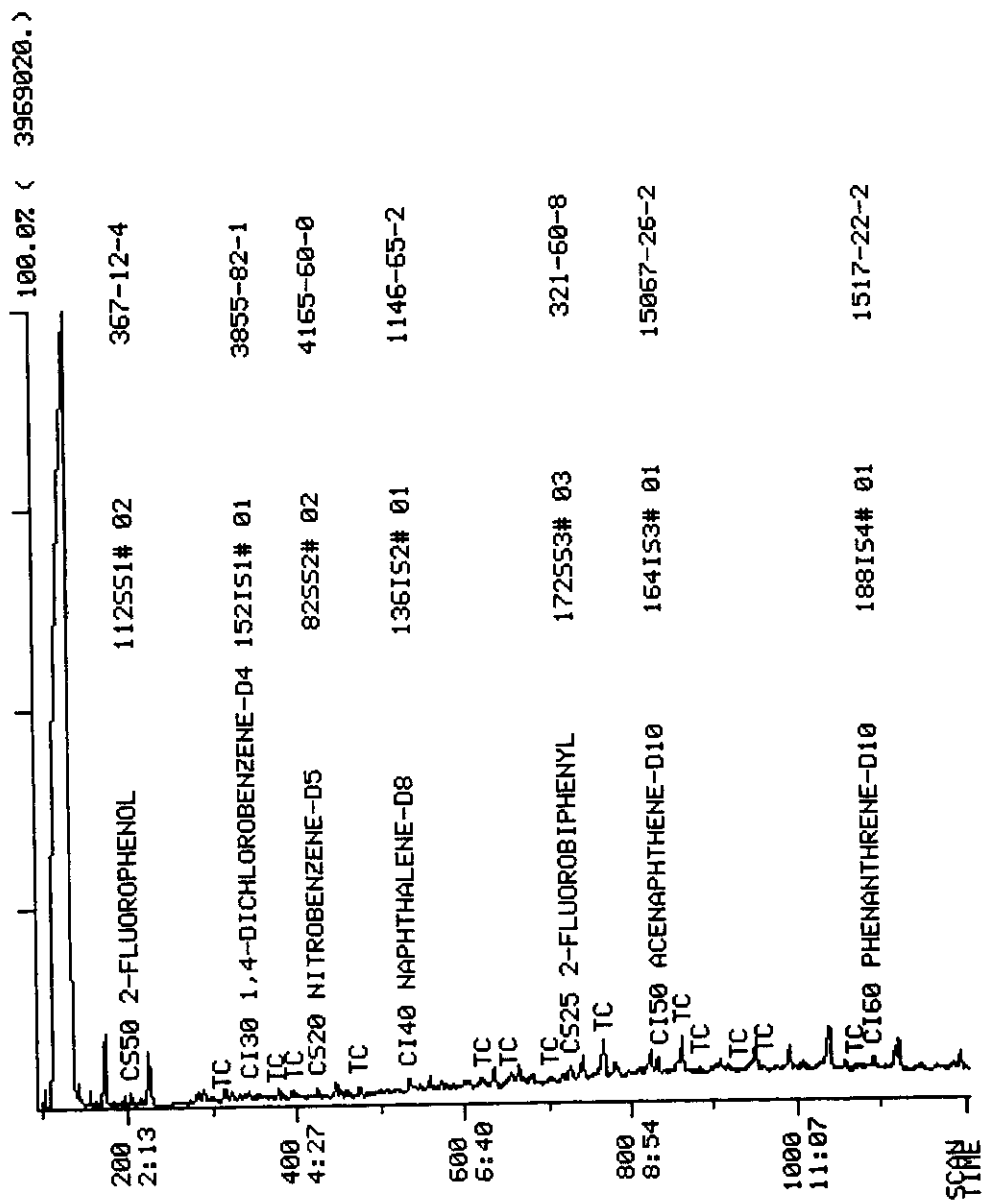
DATA FROM FILE: 300581RI03 SCANS 94 TO 1204 ACQUIRED: 09/02/98 19:59:00

CALI: 300581RI03 #3

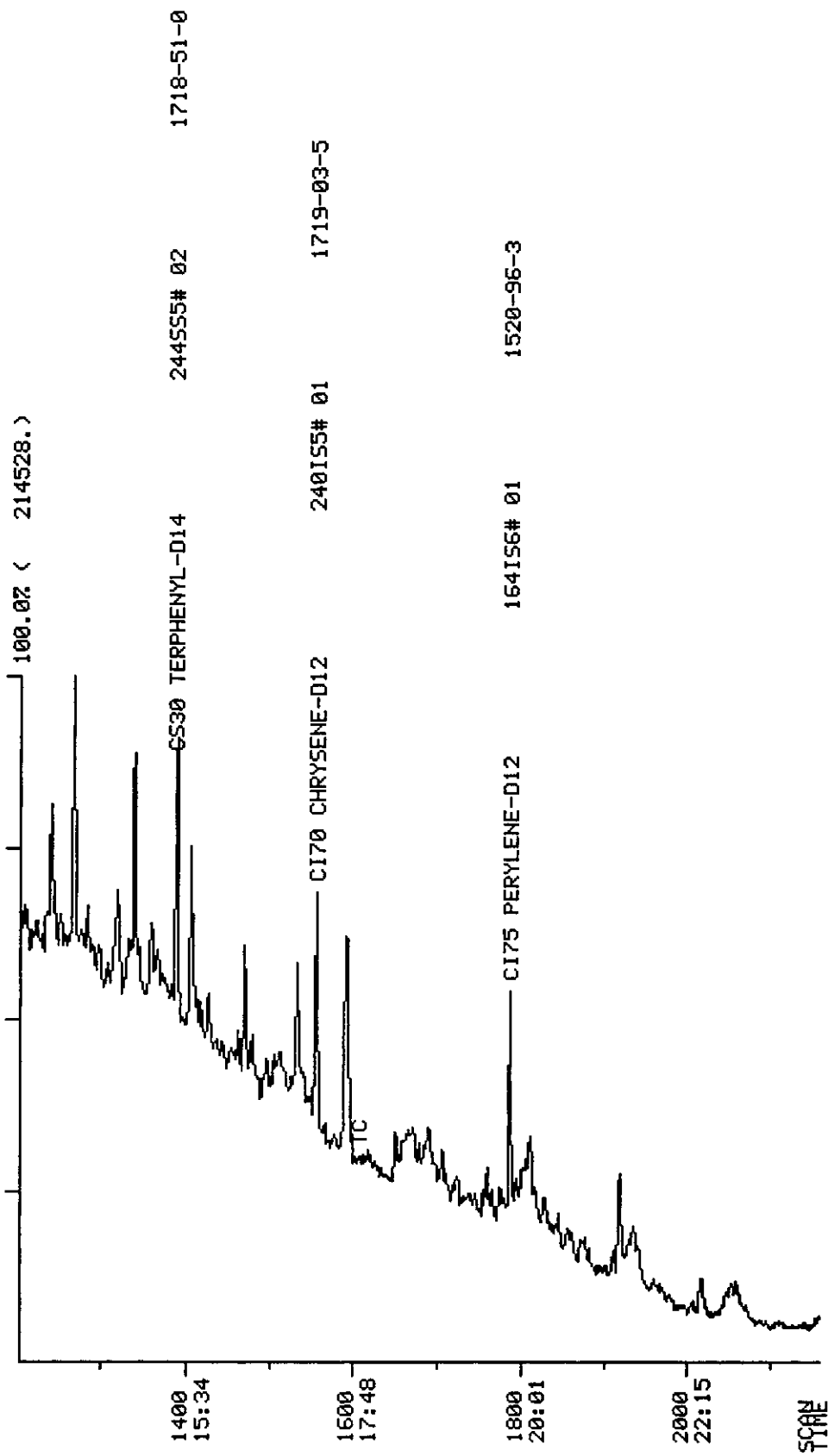
INST. ID: F16

SAMPLE: T-MM5-2 1/35A/1ML

CONDS.: UG/ML *1002/100% *(NA/NA) /1/35A NA M



DATA FROM FILE: 300681RI03 SCANS 1204 TO 2158 ACQUIRED: 09/02/98 19:59:00
 SAMPLE: T-MMS-2 1/35A/1ML CALI: 300681RI03 #3
 COND.: UG/ML *1ML *100%/100% *(NA/NA) /1/35A NA M INST. ID: F16



STANDARD USED FOR THIS REPORT IS ST16980902

REPORT PRINTED: 9/03/98 12:47:25

ADDITIONAL STANDARD USED FOR THIS REPORT IS: ST16980902A

QUANTERRA QUANTITATION SUMMARY

File: 300681RI03

Sample: T-MM5-2 1/3SA/1ML INST. ID: F16
 Analyst: DAT Instrument Id: F16 Analyzed: 09/02/98 19:59

Compounds with amounts less than 0.20 reported as NOT FOUND

Library				Mass	Meth	Scan	Ref	Area	RRF (L)	Amount
No	Entry	Name								
1	S1# 1	CI30 1,4-DICHLOROBENZENE	152	A	BV	342	1	27571.	1.000	40.000
2	S2# 1	CI40 NAPHTHALENE-D8	136	A	BB	531	2	95690.	1.000	40.000
3	S3# 1	CI50 ACENAPHTHENE-D10	164	A	BB	831	3	49039.	1.000	40.000
4	S4# 1	CI60 PHENANTHRENE-D10	188	A	BB	1089	4	81651.	1.000	40.000
5	S5# 1	CI70 CHRYSENE-D12	240	A	BB	1555	5	77811.	1.000	40.000
6	S6# 1	CI75 PERYLENE-D12	264	A	BB	1786	6	72812.	1.000	40.000
7	S2# 2	CS20 NITROBENZENE-D5	82	A	BV	423	2	45877.	0.470	40.841
8	S3# 3	CS25 2-FLUOROBIPHENYL	172	A	VB	725	3	81379.	1.234	53.778
9	S5# 2	CS30 TERPHENYL-D14	244	A	BB	1386	5	76441.	0.995	39.485
10	S1# 3	CS45 PHENOL-D5	99	A	BV	312	1	90022.	2.025	64.480
11	S1# 2	CS50 2-FLUOROPHENOL	112	A	BB	200	1	55021.	1.368	58.341
12	S3# 2	CS55 2,4,6-TRIBROMOPHENO	330	A	BB	973	3	13000.	0.236	44.998
13	S1# 5	CS70 2-CHLOROPHENOL-D4	132			1	NOT FOUND			
14	S1# 6	CS75 1,2-DICHLOROBNZN-D4	152	A	BV	342	1	27571.	0.920	43.476
15	S4# 4	HEXACHLOROBENZENE-C13	294			4	NOT FOUND			
16	S1# 15	C310 N-NITROSODIMETHYLAM	74			1	NOT FOUND			
17	S1# 20	PYRIDINE	79			1	NOT FOUND			
18	S1# 30	2-PICOLINE	93			1	NOT FOUND			
19	S1# 40	N-NITROSOMETHYLETHYLAMIN	42			1	NOT FOUND			
20	S1# 60	METHYLMETHANESULFONATE	80			1	NOT FOUND			
21	S1# 70	N-NITROSODIETHYLAMINE	102			1	NOT FOUND			
22	S1# 85	ETHYLMETHANE SULFONATE	79			1	NOT FOUND			
23	S1# 95	PENTACHLOROETHANE	117	A	VB	307	1	448.	0.602	1.081
24	S1#100	C320 ANILINE	93			1	NOT FOUND			
25	S1#105	C315 PHENOL	94	A	BB	313	1	10332.	1.968	7.617
26	S1#110	C325 BIS(2-CHLOROETHYL)E	93	A	VB	321	1	808.	1.368	0.857
27	S1#115	C330 2-CHLOROPHENOL	128	A	VB	321	1	310.	1.497	0.300
28	S1#125	C335 1,3-DICHLOROBENZENE	146			1	NOT FOUND			
29	S1#130	C340 1,4-DICHLOROBENZENE	146			1	NOT FOUND			
30	S1#145	C345 BENZYL ALCOHOL	108	A	BB	373	1	4387.	0.854	7.449
31	S1#150	C350 1,2-DICHLOROBENZENE	146			1	NOT FOUND			
32	S1#160	C355 2-METHYLPHENOL	108			1	NOT FOUND			
33	S1#165	C360 2,2'-OXYBIS(1-CLPRO	45	A	BB	395	1	3495.	2.539	1.997
34	S1#170	C361 ACETOPHENONE	105	A	BB	402	1	8684.	1.934	6.515
35	S1#175	N-NITROSOPYRROLIDINE	100	A*VB		403	1	350.	0.553	0.918
36	S1#180	N-NITROSOMORPHOLINE	56	A	VV	411	1	2422.	0.821	4.281
37	S1#182	3-METHYL PHENOL	108			1	NOT FOUND			
38	S1#185	C365 4-METHYLPHENOL	108			1	NOT FOUND			
39	S1#190	C370 N-NITROSO-DI-N-PROP	70	A	VB	404	1	1369.	0.906	2.192
40	S1#195	O-TOLUIDINE	106			1	NOT FOUND			
41	S1#200	C375 HEXACHLOROETHANE	117			1	NOT FOUND			
42	S2# 10	C410 NITROBENZENE	77	A	BB	424	2	1347.	0.445	1.266
43	S2# 15	N-NITROSOPIPERIDINE	42			2	NOT FOUND			
44	S2# 20	C415 ISOPHORONE	82	A	VB	467	2	449.	0.741	0.253
45	S2# 25	C420 2-NITROPHENOL	139	A	BB	473	2	817.	0.213	1.600

882

46	S2#	30	C425	2,4-DIMETHYLPHENOL	107	A	BB	501	2	737.	0.330	0.934
47	S2#	35	C435	BIS(2-CHLOROETHOXY)	93			2	NOT FOUND			
48	S2#	40	C440	2,4-DICHLOROPHENOL	162	A	BB	520	2	340.	0.308	0.462
49	S2#	45	C430	BENZOIC ACID	122	A	BB	543	2	6945.	0.276	10.527

50	S2# 50	AA-DIMETHYLPHENETHYLAMIN	58 A BB	707 2	3330. 1.285	1.083
51	S2# 55	C445 1,2,4-TRICHLOROBENZ	180	2 NOT FOUND		
52	S2# 60	C450 NAPHTHALENE	128 A BV	534 2	23287. 1.022	9.523
53	S2# 80	C455 4-CHLOROANALINE	127 A BB	557 2	540. 0.454	0.498
54	S2# 85	2,6-DICHLOROPHENOL	162 A BB	553 2	453. 0.257	0.736
55	S2# 90	HEXACHLOROPROPENE	213	2 NOT FOUND		
56	S2# 95	C460 HEXACHLOROBUTADIENE	225	2 NOT FOUND		
57	S2#115	P-PHENYLENE DIAMINE	108	2 NOT FOUND		
58	S2#120	N-NITROSODI-N-BUTYLAMINE	84 A VB	620 2	1528. 0.213	2.997
59	S2#130	C465 4-CHLORO-3-METHYLPH	107	2 NOT FOUND		
60	S2#140	SAFROLE	162 A BB	650 2	808. 0.248	1.363
61	S2#145	C470 2-METHYLNAPHTHALENE	142 A BV	654 2	35943. 0.751	19.999
62	S3# 10	1,2,4,5-TETRACHLOROBENZE	214	3 NOT FOUND		
63	S3# 15	ISOSAFROLE (#1)	162 A BB	699 3	481. 0.044	8.988
64	S3# 20	C510 HEXACHLOROCYCLOPENT	237	3 NOT FOUND		
65	S3# 25	C515 2,4,6-TRICHLOROPHEN	196	3 NOT FOUND		
66	S3# 30	C520 2,4,5-TRICHLOROPHEN	196	3 NOT FOUND		
67	S3# 35	ISOSAFROLE (#2)	104	3 NOT FOUND		
68	S3# 40	C525 2-CHLORONAPHTHALENE	162	3 NOT FOUND		
69	S3# 42	1-CHLORONAPHTHALEN	162	3 NOT FOUND		
70	S3# 45	C530 2-NITROANALINE	65 A BB	766 3	848. 0.502	1.378
71	S3# 50	1,4-NAPHTHOQUINONE	158	3 NOT FOUND		
72	S3# 55	C535 DIMETHYLPHTHALATE	163	3 NOT FOUND		
73	S3# 60	1,3-DINITROBENZENE	168	3 NOT FOUND		
74	S3# 65	C540 ACENAPHTHYLENE	152	3 NOT FOUND		
75	S3# 70	C543 2,6-DINITROTOLUENE	165	3 NOT FOUND		
76	S3# 75	C545 3-NITROANILINE	138	3 NOT FOUND		
77	S3# 80	C550 ACENAPHTHENE	153 A BB	836 3	2115. 1.159	1.489
78	S3# 85	C555 2,4-DINITROPHENOL	184 A BV	860 3	386. 0.189	1.663
79	S3# 90	C565 DIBENZOFURAN	168 A VB	868 3	1281. 1.669	0.626
80	S3# 95	C560 4-NITROPHENOL	109 A BB	895 3	788. 0.173	3.723
81	S3#100	PENTACHLOROBENZENE	250	3 NOT FOUND		
82	S3#105	C570 2,4-DINITROTOLUENE	165 A BB	883 3	553. 0.397	1.136
83	S3#110	1-NAPHTHYLAMINE	143	3 NOT FOUND		
84	S3#115	2-NAPHTHYLAMINE	143	3 NOT FOUND		
85	S3#120	2,3,4,6-TETRACHLOROPHENO	232	3 NOT FOUND		
86	S3#130	C580 DIETHYLPHTHALATE	149	3 NOT FOUND		
87	S3#135	C590 FLUORENE	166 A BB	928 3	4501. 1.298	2.829
88	S3#140	C585 4-CHLOROPHENYL-PHEN	204 A BB	941 3	161. 0.613	0.214
89	S3#145	5-NITRO-O-TOLUIDINE	152 A VB	943 3	1742. 0.356	3.998
90	S3#150	C595 4-NITROANALINE	138	3 NOT FOUND		
91	S4# 10	C610 4,6-DINITRO-2-METHY	198 A VV	954 4	178. 0.132	0.662
92	S4# 15	C615 N-NITROSODIPHENYLAM	169 A BV	960 4	4922. 0.551	4.378
93	S4# 20	C620 AZOBENZENE	77	4 NOT FOUND		
94	S4# 25	SYM-TRINITROBENZENE	75	4 NOT FOUND		
95	S4# 30	C625 4-BROMOPHENYL-PHENY	248	4 NOT FOUND		
96	S4# 35	PHENACETIN	108	4 NOT FOUND		
97	S4# 37	DIALATE	234	4 NOT FOUND		
98	S4# 40	C630 HEXACHLOROBENZENE	284	4 NOT FOUND		
99	S4# 45	4-AMINOBIIPHENYL	169 A BB	1078 4	2317. 0.609	1.864
100	S4# 50	C635 PENTACHLOROPHENOL	266 A BB	1068 4	242. 0.155	0.765
101	S4# 55	PRONAMIDE	173	4 NOT FOUND		
102	S4# 60	PENTACHLORONITROBENZENE	237 A BB	1066 4	216. 0.080	1.317
103	S4# 65	C640 PHENANTHRENE	178 A BB	1093 4	14817. 1.033	7.030
104	S4# 70	C645 ANTHRACENE	178	4 NOT FOUND		
105	S4# 75	2SECBUTYL-4,6-DINITROPHE	211	4 NOT FOUND		
106	S4# 80	C647 CARBAZOLE	167 A BB	1150 4	482. 0.764	0.309
107	S4# 85	C650 DI-N-BUTYLPHTHALATE	149	4 NOT FOUND		
108	S4#100	4-NITROQUINOLINE-1-OXIDE	190	4 NOT FOUND		
109	S4#105	METHAPYRILENE	58	4 NOT FOUND		
110	S4#106	ISODRIN	193	4 NOT FOUND		

111	S4#110	C655 FLUORANTHENE	202	A BB	1301	4	404.	0.900	0.220
112	S4#120	CHLOROBENZILATE	139		4	NOT FOUND			
113	S5# 10	C710 BENZIDINE	184		5	NOT FOUND			
114	S5# 15	C715 PYRENE	202	A BB	1337	5	2218.	1.323	0.862
115	S5# 20	ARAMITE (#1)	185		5	NOT FOUND			
116	S5# 25	ARAMITE (#2)	185		5	NOT FOUND			
117	S5# 30	P-DIMETHYLAMINOAZOBENZEN	120		5	NOT FOUND			
118	S5# 35	3,3'-DIMETHYLBENZIDINE	212	A BV	1474	5	510.	0.418	0.628
119	S5# 37	KEPONE	272		5	NOT FOUND			
120	S5# 40	C720 BUTYLBENZYLPHTHALAT	149		5	NOT FOUND			
121	S5# 45	2-ACETYLAMINOFLUORENE	181		5	NOT FOUND			
122	S5# 50	C730 BENZO (A) ANTHRACENE	228	A BB	1558	5	1798.	1.125	0.822
123	S5# 55	C725 3,3'-DICHLOROBENZID	252		5	NOT FOUND			
124	S5# 60	C740 CHRYSENE	228	A BB	1558	5	1799.	1.014	0.912
125	S5# 65	C745 BIS (2-ETHYLHEXYL) PH	149	A BB	1605	5	3055.	0.980	1.603
126	S5# 85	3-METHYLCHOLANTHRENE	268	A*BB	1841	5	287.	0.548	0.270
127	S6# 10	C760 DI-N-OCTYL PHTHALAT	149		6	NOT FOUND			
128	S6# 15	C765 BENZO (B) FLUORANTHEN	252		6	NOT FOUND			
129	S6# 20	7,12-DIMETHYLBENZANTHRAC	256	A BB	1743	6	337.	0.610	0.303
130	S6# 25	C770 BENZO (K) FLUORANTHEN	252		6	NOT FOUND			
131	S6# 30	HEXACHLOROPHENE	196		6	NOT FOUND			
132	S6# 35	C775 BENZO (A) PYRENE	252		6	NOT FOUND			
133	S6# 55	C780 INDENO (1,2,3-CD) PYR	276		6	NOT FOUND			
134	S6# 60	C785 DIBENZ (A,H) ANTHRACE	278		6	NOT FOUND			
135	S6# 65	C790 BENZO (G,H,I) PERYLEN	276		6	NOT FOUND			

QUANTERRA Internal Standard Check

Standard Filename: ST16980902

Analyzed: 09/02/98 14:53

Standard	IS# 1			IS# 2			IS# 3		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980902	23319.		3.78	79260.		5.91	43605.		9.22
Upper Limit	46638.=+200		4.28	158519.=+200		6.41	87211.=+200		9.72
Lower Limit	11659.=+ 50		3.28	39630.=+ 50		5.41	21803.=+ 50		8.72
Filename									
1 300681RI03	27571.= 118		3.80	95690.= 120		5.91	49039.= 112		9.24

IS# 1 = CI30 1,4-DICHLOROBENZENE-D4

IS# 2 = CI40 NAPHTHALENE-D8

IS# 3 = CI50 ACENAPHTHENE-D10

* - indicates an I.S. is outside QC limit(s)

Standard	IS# 4			IS# 5			IS# 6		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980902	76218.		12.09	51424.		17.27	44306.		19.84
Upper Limit	152436.=+200		12.59	102848.=+200		17.77	88612.=+200		20.34
Lower Limit	38109.=+ 50		11.59	25712.=+ 50		16.77	22153.=+ 50		19.34
Filename									
1 300681RI03	81651.= 107		12.11	77811.= 151		17.30	72812.= 164		19.87

IS# 4 = CI60 PHENANTHRENE-D10

IS# 5 = CI70 CHRYSENE-D12

IS# 6 = CI75 PERYLENE-D12

* - indicates an I.S. is outside QC limit(s)

Semivolatile Organics
Method 0010/8270

Client Name: Pacific Environmental Services

Client ID: T-MM5-FB-F, FH, XAD, COND, BH

LAB ID: 300681-0004-SA

Matrix: AIRTRAIN

Authorized: 30 JUL 98

Sampled: 25 JUL 98

Prepared: 31 JUL 98

Received: 30 JUL 98

Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
Phenol	ND	ug/Sample	30	R
bis(2-Chloroethyl)ether	ND	ug/Sample	30	
2-Chlorophenol	ND	ug/Sample	30	
1,3-Dichlorobenzene	ND	ug/Sample	30	
1,4-Dichlorobenzene	ND	ug/Sample	30	
Benzyl alcohol	ND	ug/Sample	30	
1,2-Dichlorobenzene	ND	ug/Sample	30	
2-Methylphenol	ND	ug/Sample	30	
2,2'-Oxybis(1-chloropropane)	ND	ug/Sample	30	
3/4-Methylphenol	ND	ug/Sample	30	
N-Nitroso-di-n-propylamine	ND	ug/Sample	30	
Hexachloroethane	ND	ug/Sample	30	
Nitrobenzene	ND	ug/Sample	30	
Isophorone	ND	ug/Sample	30	
2-Nitrophenol	ND	ug/Sample	30	
2,4-Dimethylphenol	ND	ug/Sample	30	
Benzoic acid	ND	ug/Sample	150	
bis(2-Chloroethoxy)-methane	ND	ug/Sample	30	
2,4-Dichlorophenol	ND	ug/Sample	30	
1,2,4-Trichlorobenzene	ND	ug/Sample	30	
Naphthalene	ND	ug/Sample	30	
4-Chloroaniline	ND	ug/Sample	30	
Hexachlorobutadiene	ND	ug/Sample	30	
4-Chloro-3-methylphenol	ND	ug/Sample	30	
2-Methylnaphthalene	ND	ug/Sample	30	
Hexachlorocyclopentadiene	ND	ug/Sample	30	
2,4,6-Trichlorophenol	ND	ug/Sample	30	
2,4,5-Trichlorophenol	ND	ug/Sample	150	
2-Chloronaphthalene	ND	ug/Sample	30	
2-Nitroaniline	ND	ug/Sample	30	
Dimethyl phthalate	ND	ug/Sample	30	
Acenaphthylene	ND	ug/Sample	30	
3-Nitroaniline	ND	ug/Sample	150	
Acenaphthene	ND	ug/Sample	30	
2,4-Dinitrophenol	ND	ug/Sample	150	
4-Nitrophenol	ND	ug/Sample	150	
Dibenzofuran	ND	ug/Sample	30	

Note R = Reporting limit(s) raised due to sample volume limitations.

ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.

Rev 230787

887

Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: T-MM5-FB-F, FH, XAD, COND, BH
LAB ID: 300681-0004-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 25 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
2,4-Dinitrotoluene	ND	ug/Sample	30	
2,6-Dinitrotoluene	ND	ug/Sample	30	
Diethyl phthalate	ND	ug/Sample	30	
4-Chlorophenyl phenyl ether	ND	ug/Sample	30	
Fluorene	ND	ug/Sample	30	
4-Nitroaniline	ND	ug/Sample	150	
4,6-Dinitro-2-methylphenol	ND	ug/Sample	150	
N-Nitrosodiphenylamine	ND	ug/Sample	30	
4-Bromophenyl phenyl ether	ND	ug/Sample	30	
Hexachlorobenzene	ND	ug/Sample	30	
Pentachlorophenol	ND	ug/Sample	150	
Phenanthrene	ND	ug/Sample	30	
Anthracene	ND	ug/Sample	30	
Di-n-butyl phthalate	ND	ug/Sample	30	
Fluoranthene	ND	ug/Sample	30	
Pyrene	ND	ug/Sample	30	
Butyl benzyl phthalate	ND	ug/Sample	30	
3,3'-Dichlorobenzidine	ND	ug/Sample	60	
Benzo(a)anthracene	ND	ug/Sample	30	
bis(2-Ethylhexyl)-phthalate	ND	ug/Sample	30	
Chrysene	ND	ug/Sample	30	
Di-n-octyl phthalate	ND	ug/Sample	30	
Benzo(b)fluoranthene	ND	ug/Sample	30	
Benzo(k)fluoranthene	ND	ug/Sample	30	
Benzo(a)pyrene	ND	ug/Sample	30	
Indeno(1,2,3-cd)pyrene	ND	ug/Sample	30	
Dibenz(a,h)anthracene	ND	ug/Sample	30	
Benzo(g,h,i)perylene	ND	ug/Sample	30	
Acetophenone	ND	ug/Sample	30	
4-Aminobiphenyl	ND	ug/Sample	150	
Aniline	ND	ug/Sample	30	
Benzidine	ND	ug/Sample	300	
3,3'-Dimethylbenzidine	ND	ug/Sample	60	
N-Nitrosodimethylamine	ND	ug/Sample	30	
N-Nitrosomorpholine	ND	ug/Sample	30	
Pentachloronitrobenzene (PCNB)	ND	ug/Sample	150	
o-Toluidine	ND	ug/Sample	60	

ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.
Rev 230787

888

Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services

Client ID: T-MM5-FB-F, FH, XAD, COND, BH

LAB ID: 300681-0004-SA

Matrix: AIRTRAIN

Authorized: 30 JUL 98

Sampled: 25 JUL 98

Prepared: 31 JUL 98

Received: 30 JUL 98

Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
2-Methoxybenzenamine	ND	ug/Sample	--	K
Biphenyl	ND	ug/Sample	--	K
Chloroacetophenone	ND	ug/Sample	--	K
Cumene	ND	ug/Sample	--	K
DBCP (1,2-Dibromo-3-chloropropane)	ND	ug/Sample	--	K
Benzo(e)pyrene	ND	ug/Sample	--	K
N-N-Diethylaniline	ND	ug/Sample	--	K
Dimethylaniline	ND	ug/Sample	--	K
3,3'-Dimethoxybenzidine	ND	ug/Sample	--	K
Hydroquinone	ND	ug/Sample	--	K
4,4'-Methyl-bis(2-chloroaniline)	ND	ug/Sample	--	K
4-Nitrodiphenyl	ND	ug/Sample	--	K
Trifluralin	ND	ug/Sample	--	K

Surrogate	Recovery	Acceptable Range
Nitrobenzene-d5	60 %	45 - 107
2-Fluorobiphenyl	85 %	62 - 110
Terphenyl-d14	94 %	58 - 135
Phenol-d5	65 %	43 - 130
2-Fluorophenol	56 %	36 - 111
2,4,6-Tribromophenol	70 %	58 - 131

Note K = Identified by mass spectrum only; quantitation based on 1:1 response with internal standard.

ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.

Rev 230787

888

Semivolatiles Library Search (20 Compound TID)
Method 8270

Client Name: Pacific Environmental Services
Client ID: T-MM5-FB-F, FH, XAD, COND, BH
LAB ID: 300681-0004-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 25 JUL 98
Prepared: NA

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Result	Units	Reporting Limit	Qualifier
4-Hydroxy-4-methyl-2-pentanone	260	ug/Sample	--	0
n-Nonane	260	ug/Sample	--	0
Benzaldehyde	22	ug/Sample	--	0
Unknown	23	ug/Sample	--	
Hexadecanoic acid	29	ug/Sample	--	
Unknown	54	ug/Sample	--	
Unknown	48	ug/Sample	--	
Unknown	42	ug/Sample	--	0
Unknown	62	ug/Sample	--	
5-Eiconsene, (E) -	160	ug/Sample	--	0
Unknown	55	ug/Sample	--	
Unknown	40	ug/Sample	--	
Unknown	20	ug/Sample	--	
Unknown	58	ug/Sample	--	
Unknown	73	ug/Sample	--	
Unknown alkene	75	ug/Sample	--	
Unknown	74	ug/Sample	--	
Unknown alkene	46	ug/Sample	--	
Unknown	40	ug/Sample	--	
Unknown	29	ug/Sample	--	

Note 0 = Or structurally similar compound (isomer).
NA = Not Applicable

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.
Rev 230787

890

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068104
Std Id: ST16980831

Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 20:30
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Surrogate Spike Recoveries
8270-G, LIMS 10/10/96

Surrogate	Surrogate Amount (ug)		% Recovery	QC Limits
	Spiked	Measured	Measured	
CS20 NITROBENZENE-D5	50.00	29.90	59.8	45 107
CS25 2-FLUOROBIPHENY	50.00	42.70	85.4	62 110
CS30 TERPHENYL-D14	50.00	47.00	94.0	58 135
CS45 PHENOL-D5	100.0	65.10	65.1	43 130
CS50 2-FLUOROPHENOL	100.0	55.70	55.7	36 111
CS55 2,4,6-TRIBROMOP	100.0	69.90	69.9	58 131

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
HEXACHLOROBENZENE-C13		ND	UG/A	30.0
C310 N-NITROSODIMETHYLAMINE		ND	UG/A	30.0
PYRIDINE		ND	UG/A	60.0
2-PICOLINE		ND	UG/A	30.0
N-NITROSOMETHYLETHYLAMINE		ND	UG/A	30.0
METHYLMETHANESULFONATE		ND	UG/A	30.0
N-NITROSODIETHYLAMINE		ND	UG/A	30.0
ETHYLMETHANE SULFONATE		ND	UG/A	30.0
PENTACHLOROETHANE		ND	UG/A	30.0
C320 ANILINE		ND	UG/A	30.0
C315 PHENOL		ND	UG/A	30.0
C325 BIS(2-CHLOROETHYL)ETHER		ND	UG/A	30.0
C330 2-CHLOROPHENOL		ND	UG/A	30.0
C335 1,3-DICHLOROBENZENE		ND	UG/A	30.0
C340 1,4-DICHLOROBENZENE		ND	UG/A	30.0
C345 BENZYL ALCOHOL		ND	UG/A	30.0
C350 1,2-DICHLOROBENZENE		ND	UG/A	30.0
C355 2-METHYLPHENOL		ND	UG/A	30.0
C360 2,2'-OXYBIS(1-CLPROPAN)		ND	UG/A	30.0
C361 ACETOPHENONE		ND	UG/A	30.0
N-NITROSPYRROLIDINE		ND	UG/A	30.0
N-NITROSOMORPHOLINE		ND	UG/A	30.0
3-METHYL PHENOL		ND	UG/A	30.0
C365 4-METHYLPHENOL		ND	UG/A	30.0
C370 N-NITROSO-DI-N-PROPYLAM		ND	UG/A	30.0
O-TOLUIDINE		ND	UG/A	30.0
C375 HEXACHLOROETHANE		ND	UG/A	30.0
C410 NITROBENZENE		ND	UG/A	30.0
N-NITROSOPIPERIDINE		ND	UG/A	30.0

Reviewed by: K9/17/0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068104
Std Id: ST16980831

Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 20:30
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C415 ISOPHORONE		ND	UG/A	30.0
C420 2-NITROPHENOL		ND	UG/A	30.0
C425 2,4-DIMETHYLPHENOL		ND	UG/A	30.0
C435 BIS(2-CHLOROETHOXY)METH		ND	UG/A	30.0
C440 2,4-DICHLOROPHENOL		ND	UG/A	30.0
C430 BENZOIC ACID		ND	UG/A	150.0
AA-DIMETHYLPHENETHYLAMINE		ND	UG/A	30.0
C445 1,2,4-TRICHLOROBENZENE		ND	UG/A	30.0
C450 NAPHTHALENE		ND	UG/A	30.0
C455 4-CHLOROANILINE		ND	UG/A	30.0
2,6-DICHLOROPHENOL		ND	UG/A	30.0
HEXACHLOROPROPENE		ND	UG/A	30.0
C460 HEXACHLOROBUTADIENE		ND	UG/A	30.0
P-PHENYLENE DIAMINE		ND	UG/A	30.0
N-NITROSODI-N-BUTYLAMINE		ND	UG/A	30.0
C465 4-CHLORO-3-METHYLPHENO		ND	UG/A	30.0
SAFROLE		ND	UG/A	30.0
C470 2-METHYLNAPHTHALENE		ND	UG/A	30.0
1,2,4,5-TETRACHLOROBENZENE		ND	UG/A	30.0
ISOSAFROLE (#1)		ND	UG/A	60.0
C510 HEXACHLOROCYCLOPENTADI		ND	UG/A	30.0
C515 2,4,6-TRICHLOROPHENOL		ND	UG/A	30.0
C520 2,4,5-TRICHLOROPHENOL		ND	UG/A	30.0
ISOSAFROLE (#2)		ND	UG/A	60.0
C525 2-CHLORONAPHTHALENE		ND	UG/A	30.0
1-CHLORONAPHTHALENE		ND	UG/A	15.0
C530 2-NITROANILINE		ND	UG/A	150.0
1,4-NAPHTHOQUINONE		ND	UG/A	30.0
C535 DIMETHYLPHTHALATE		ND	UG/A	30.0
1,3-DINITROBENZENE		ND	UG/A	30.0
C540 ACENAPHTHYLENE		ND	UG/A	30.0
C543 2,6-DINITROTOLUENE		ND	UG/A	30.0
C545 3-NITROANILINE		ND	UG/A	150.0
C550 ACENAPHTHENE		ND	UG/A	30.0
C555 2,4-DINITROPHENOL		ND	UG/A	150.0
C565 DIBENZOFURAN		ND	UG/A	30.0
C560 4-NITROPHENOL		ND	UG/A	150.0
PENTACHLOROBENZENE		ND	UG/A	30.0
C570 2,4-DINITROTOLUENE		ND	UG/A	30.0
1-NAPHTHYLAMINE		ND	UG/A	30.0
2-NAPHTHYLAMINE		ND	UG/A	30.0
2,3,4,6-TETRACHLOROPHENOL		ND	UG/A	60.0
C580 DIETHYLPHTHALATE		ND	UG/A	30.0
C590 FLUORENE		ND	UG/A	30.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068104
Std Id: ST16980831

Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 20:30
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
=====				
C585 4-CHLOROPHENYL-PHENYLE		ND	UG/A	30.0
5-NITRO-O-TOLUIDINE		ND	UG/A	30.0
C595 4-NITROANALINE		ND	UG/A	150.0
C610 4,6-DINITRO-2-METHYLPH		ND	UG/A	150.0
C615 N-NITROSODIPHENYLAMINE		ND	UG/A	30.0
C620 AZOBENZENE		ND	UG/A	30.0
SYM-TRINITROBENZENE		ND	UG/A	30.0
C625 4-BROMOPHENYL-PHENYLET		ND	UG/A	30.0
PHENACETIN		ND	UG/A	30.0
DIALATE	234	ND	UG/A	30.0
C630 HEXACHLOROBENZENE		ND	UG/A	30.0
4-AMINOBIIPHENYL		ND	UG/A	30.0
C635 PENTACHLOROPHENOL		ND	UG/A	150.0
PRONAMIDE		ND	UG/A	30.0
PENTACHLORONITROBENZENE		ND	UG/A	150.0
C640 PHENANTHRENE		ND	UG/A	30.0
C645 ANTHRACENE		ND	UG/A	30.0
2SECBUTYL-4,6-DINITROPHENOL		ND	UG/A	30.0
C647 CARBAZOLE		ND	UG/A	30.0
C650 DI-N-BUTYLPHTHALATE		ND	UG/A	30.0
4-NITROQUINOLINE-1-OXIDE		ND	UG/A	30.0
METHAPYRILENE		ND	UG/A	30.0
ISODRIN		ND	UG/A	30.0
C655 FLUORANTHENE		ND	UG/A	30.0
CHLOROBENZILATE		ND	UG/A	30.0
C710 BENZIDINE		ND	UG/A	300.0
C715 PYRENE		ND	UG/A	30.0
ARAMITE (#1)		ND	UG/A	30.0
ARAMITE (#2)		ND	UG/A	30.0
P-DIMETHYLAMINOAZOBENZENE		ND	UG/A	30.0
3,3'-DIMETHYLBENZIDINE		ND	UG/A	30.0
KEPONE		ND	UG/A	150.0
C720 BUTYLBENZYLPHTHALATE		ND	UG/A	30.0
2-ACETYLAMINOFLUORENE		ND	UG/A	30.0
C730 BENZO(A)ANTHRACENE		ND	UG/A	30.0
C725 3,3'-DICHLOBENZIDINE		ND	UG/A	60.0
C740 CHRYSENE		ND	UG/A	30.0
C745 BIS(2-ETHYLHEXYL)PHTHA		ND	UG/A	30.0
3-METHYLCHOLANTHRENE		ND	UG/A	30.0
C760 DI-N-OCTYL PHTHALATE		ND	UG/A	30.0
C765 BENZO(B)FLUORANTHENE		ND	UG/A	30.0
7,12-DIMETHYLBENZANTHRACENE		ND	UG/A	30.0
C770 BENZO(K)FLUORANTHENE		ND	UG/A	30.0
HEXACHLOROPHENE		ND	UG/A	30.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

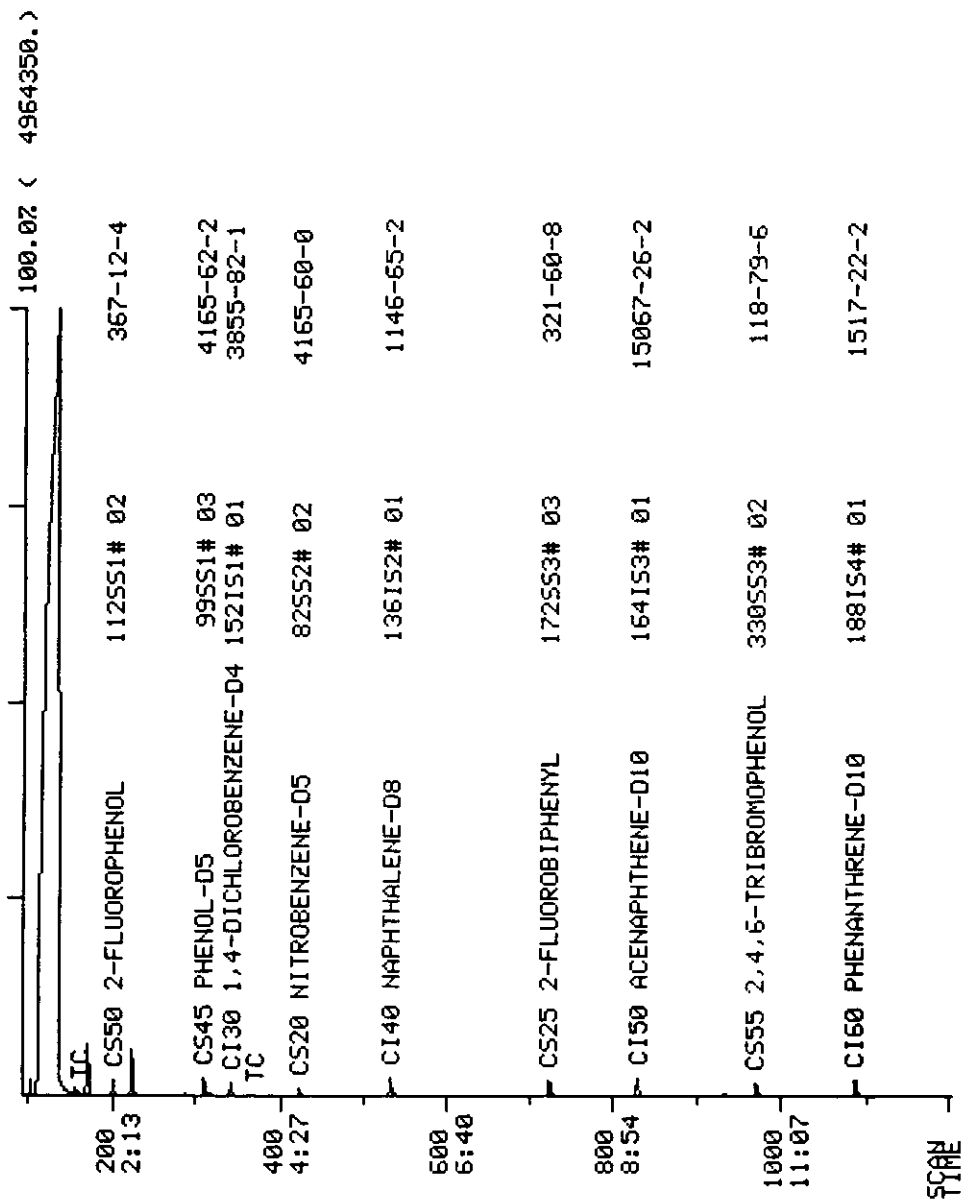
Data File: 30068104
Std Id: ST16980831

Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 20:30
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

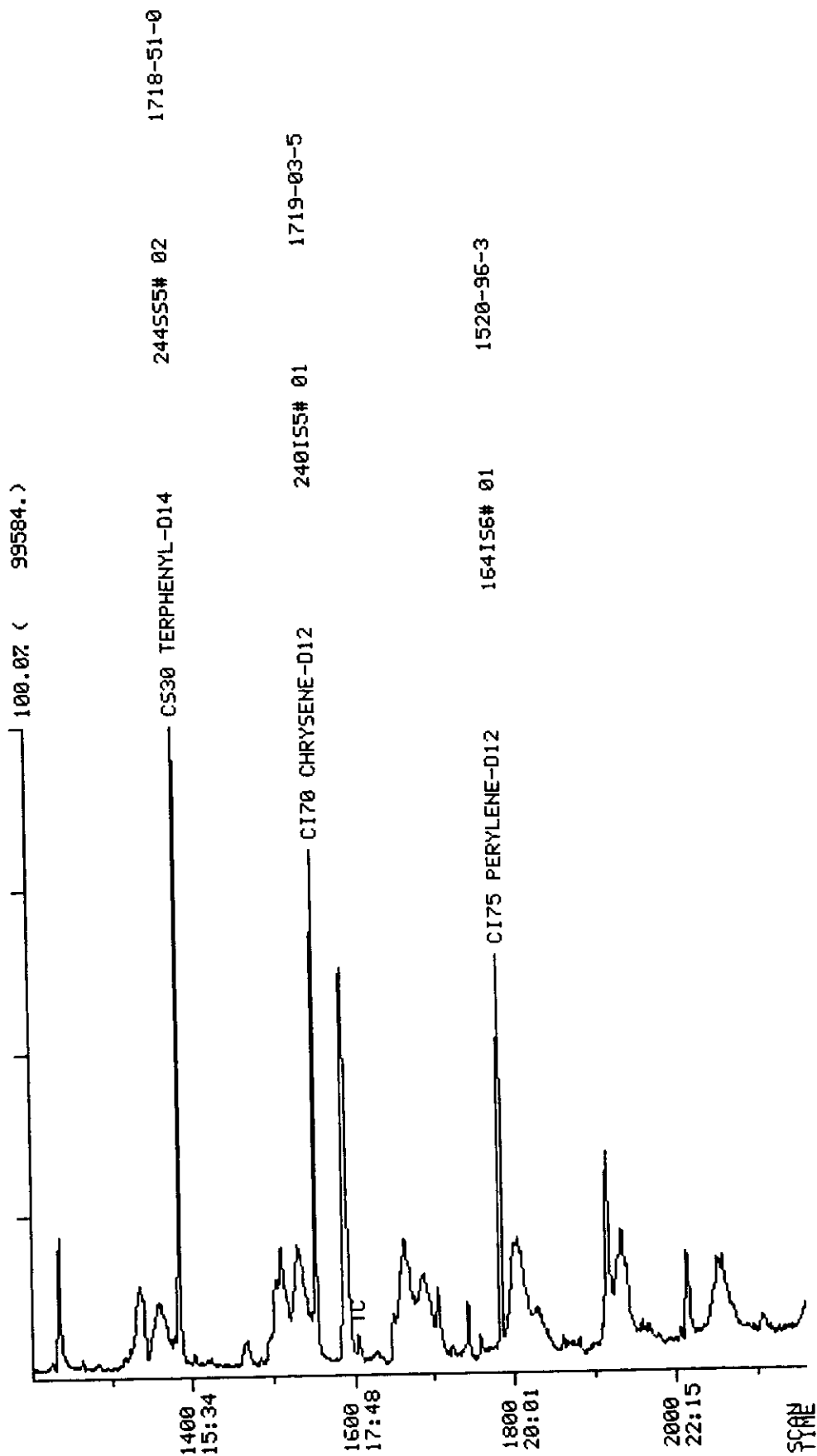
Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C775 BENZO(A)PYRENE		ND	UG/A	30.0
C780 INDENO(1,2,3-CD)PYRENE		ND	UG/A	30.0
C785 DIBENZ(A,H)ANTHRACENE		ND	UG/A	30.0
C790 BENZO(G,H,I)PERYLENE		ND	UG/A	30.0

DATA FROM FILE: 30068104 SCANS 93 TO 1203 ACQUIRED: 08/31/98 20:30:00
 CALL: 30068104 #3
 SAMPLE: T-MMS-FB 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *1002/1002 *KNA/NA >1/35A NA M



DATA FROM FILE: 30068104 SCANS 1203 TO 2158 ACQUIRED: 08/31/98 20:30:00
 CALI: 30068104 #3
 SAMPLE: T-MMS-FB 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA) /1/35A NA M



ADDITIONAL STANDARD USED FOR THIS REPORT IS: ST16980831A

QUANTERRA QUANTITATION SUMMARY

File: 30068104

Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Analyst: DAT Instrument Id: F16 Analyzed: 08/31/98 20:30

Compounds with amounts less than 0.20 reported as NOT FOUND

Library				Units: UG/ML			
No	Entry	Name	Mass Meth	Scan Ref	Area	RRF(L)	Amount
1	S1#	1 C130 1,4-DICHLOROBENZENE	152 A BB	341 1	33946.	1.000	40.000
2	S2#	1 C140 NAPHTHALENE-DB	136 A BB	529 2	129608.	1.000	40.000
3	S3#	1 C150 ACENAPHTHENE-D10	164 A BB	828 3	68092.	1.000	40.000
4	S4#	1 C160 PHENANTHRENE-D10	188 A BB	1085 4	115982.	1.000	40.000
5	S5#	1 C170 CHRYSENE-D12	240 A BB	1551 5	71692.	1.000	40.000
6	S6#	1 C175 PERYLENE-D12	264 A BB	1782 6	58915.	1.000	40.000
7	S2#	2 CS20 NITROBENZENE-D5	82 A BB	421 2	45553.	0.470	29.940
8	S3#	3 CS25 2-FLUOROBIPHENYL	172 A BB	722 3	89717.	1.234	42.699
9	S5#	2 CS30 TERPHENYL-D14	244 A BB	1383 5	83889.	0.995	47.030
10	S1#	3 CS45 PHENOL-D5	99 A BB	308 1	111982.	2.025	65.145
11	S1#	2 CS50 2-FLUOROPHENOL	112 A BB	199 1	64661.	1.368	55.686
12	S3#	2 CS55 2,4,6-TRIBROMOPHENO	330 A BB	968 3	28049.	0.236	69.918
13	S4#	4 HEXACHLOROBENZENE-C13	294	4 NOT FOUND			
14	S1#	15 C310 N-NITROSODIMETHYLAM	74	1 NOT FOUND			
15	S1#	20 PYRIDINE	79	1 NOT FOUND			
16	S1#	30 2-PICOLINE	93	1 NOT FOUND			
17	S1#	40 N-NITROSOMETHYLETHYLAMIN	42 A VB	159 1	3558.	0.811	5.169
18	S1#	60 METHYLMETHANESULFONATE	80	1 NOT FOUND			
19	S1#	70 N-NITROSODIETHYLAMINE	102	1 NOT FOUND			
20	S1#	85 ETHYLMETHANE SULFONATE	79	1 NOT FOUND			
21	S1#	95 PENTACHLOROETHANE	117	1 NOT FOUND			
22	S1#	100 C320 ANILINE	93	1 NOT FOUND			
23	S1#	105 C315 PHENOL	94 A BB	309 1	388.	1.968	0.232
24	S1#	110 C325 BIS(2-CHLOROETHYL)E	93	1 NOT FOUND			
25	S1#	115 C330 2-CHLOROPHENOL	128	1 NOT FOUND			
26	S1#	125 C335 1,3-DICHLOROBENZENE	146	1 NOT FOUND			
27	S1#	130 C340 1,4-DICHLOROBENZENE	146	1 NOT FOUND			
28	S1#	145 C345 BENZYL ALCOHOL	108 A BB	369 1	2780.	0.854	3.834
29	S1#	150 C350 1,2-DICHLOROBENZENE	146	1 NOT FOUND			
30	S1#	160 C355 2-METHYLPHENOL	108	1 NOT FOUND			
31	S1#	165 C360 2,2'-OXYBIS(1-CLPRO	45	1 NOT FOUND			
32	S1#	170 C361 ACETOPHENONE	105 A BB	404 1	412.	1.934	0.251
33	S1#	175 N-NITROSOPYRROLIDINE	100	1 NOT FOUND			
34	S1#	180 N-NITROSOMORPHOLINE	56 A BB	413 1	224.	0.821	0.322
35	S1#	182 3-METHYL PHENOL	108 A BB	422 1	402.	2.252	0.210
36	S1#	185 C365 4-METHYLPHENOL	108	1 NOT FOUND			
37	S1#	190 C370 N-NITROSO-DI-N-PROP	70	1 NOT FOUND			
38	S1#	195 O-TOLUIDINE	106	1 NOT FOUND			
39	S1#	200 C375 HEXACHLOROETHANE	117	1 NOT FOUND			
40	S2#	10 C410 NITROBENZENE	77	2 NOT FOUND			
41	S2#	15 N-NITROSOPIPERIDINE	42	2 NOT FOUND			
42	S2#	20 C415 ISOPHORONE	82	2 NOT FOUND			
43	S2#	25 C420 2-NITROPHENOL	139	2 NOT FOUND			
44	S2#	30 C425 2,4-DIMETHYLPHENOL	107	2 NOT FOUND			
45	S2#	35 C435 BIS(2-CHLOROETHOXY)	93	2 NOT FOUND			

46	S2#	40	C440	2,4-DICHLOROPHENOL	162	2	NOT FOUND
47	S2#	45	C430	BENZOIC ACID	122	2	NOT FOUND
48	S2#	50	AA-DIMETHYLPHENETHYLAMIN	58	2	NOT FOUND	
49	S2#	55	C445	1,2,4-TRICHLOROBENZ	180	2	NOT FOUND

50	S2# 60	C450 NAPHTHALENE	128	2	NOT FOUND
51	S2# 80	C455 4-CHLOROANALINE	127	2	NOT FOUND
52	S2# 85	2,6-DICHLOROPHENOL	162	2	NOT FOUND
53	S2# 90	HEXACHLOROPROPENE	213	2	NOT FOUND
54	S2# 95	C460 HEXACHLOROBUTADIENE	225	2	NOT FOUND
55	S2#115	P-PHENYLENE DIAMINE	108	2	NOT FOUND
56	S2#120	N-NITROSODI-N-BUTYLAMINE	84	2	NOT FOUND
57	S2#130	C465 4-CHLORO-3-METHYLPH	107	2	NOT FOUND
58	S2#140	SAFROLE	162	2	NOT FOUND
59	S2#145	C470 2-METHYLNAPHTHALENE	142	2	NOT FOUND
60	S3# 10	1,2,4,5-TETRACHLOROBENZE	214	3	NOT FOUND
61	S3# 15	ISOSAFROLE (#1)	162	3	NOT FOUND
62	S3# 20	C510 HEXACHLOROCYCLOPENT	237	3	NOT FOUND
63	S3# 25	C515 2,4,6-TRICHLOROPHEN	196	3	NOT FOUND
64	S3# 30	C520 2,4,5-TRICHLOROPHEN	196	3	NOT FOUND
65	S3# 35	ISOSAFROLE (#2)	104	3	NOT FOUND
66	S3# 40	C525 2-CHLORONAPHTHALENE	162	3	NOT FOUND
67	S3# 42	1-CHLORONAPHTHALENE	162	3	NOT FOUND
68	S3# 45	C530 2-NITROANALINE	65	3	NOT FOUND
69	S3# 50	1,4-NAPHTHOQUINONE	158	3	NOT FOUND
70	S3# 55	C535 DIMETHYLPHTHALATE	163	3	NOT FOUND
71	S3# 60	1,3-DINITROBENZENE	168	3	NOT FOUND
72	S3# 65	C540 ACENAPHTHYLENE	152	3	NOT FOUND
73	S3# 70	C543 2,6-DINITROTOLUENE	165	3	NOT FOUND
74	S3# 75	C545 3-NITROANILINE	138	3	NOT FOUND
75	S3# 80	C550 ACENAPHTHENE	153	3	NOT FOUND
76	S3# 85	C555 2,4-DINITROPHENOL	184	3	NOT FOUND
77	S3# 90	C565 DIBENZOFURAN	168	3	NOT FOUND
78	S3# 95	C560 4-NITROPHENOL	109	3	NOT FOUND
79	S3#100	PENTACHLOROBENZENE	250	3	NOT FOUND
80	S3#105	C570 2,4-DINITROTOLUENE	165	3	NOT FOUND
81	S3#110	1-NAPHTHYLAMINE	143	3	NOT FOUND
82	S3#115	2-NAPHTHYLAMINE	143	3	NOT FOUND
83	S3#120	2,3,4,6-TETRACHLOROPHENO	232	3	NOT FOUND
84	S3#130	C580 DIETHYLPHTHALATE	149	3	NOT FOUND
85	S3#135	C590 FLUORENE	166	3	NOT FOUND
86	S3#140	C585 4-CHLOROPHENYL-PHEN	204	3	NOT FOUND
87	S3#145	5-NITRO-O-TOLUIDINE	152	3	NOT FOUND
88	S3#150	C595 4-NITROANALINE	138	3	NOT FOUND
89	S4# 10	C610 4,6-DINITRO-2-METHY	198	4	NOT FOUND
90	S4# 15	C615 N-NITROSODIPHENYLAM	169	4	NOT FOUND
91	S4# 20	C620 AZOBENZENE	77	4	NOT FOUND
92	S4# 25	SYM-TRINITROBENZENE	75	4	NOT FOUND
93	S4# 30	C625 4-BROMOPHENYL-PHENY	248	4	NOT FOUND
94	S4# 35	PHENACETIN	108	4	NOT FOUND
95	S4# 37	DIALATE	234	4	NOT FOUND
96	S4# 40	C630 HEXACHLOROBENZENE	284	4	NOT FOUND
97	S4# 45	4-AMINOBIIPHENYL	169	4	NOT FOUND
98	S4# 50	C635 PENTACHLOROPHENOL	266	4	NOT FOUND
99	S4# 55	PRONAMIDE	173	4	NOT FOUND
100	S4# 60	PENTACHLORONITROBENZENE	237	4	NOT FOUND
101	S4# 65	C640 PHENANTHRENE	178	4	NOT FOUND
102	S4# 70	C645 ANTHRACENE	178	4	NOT FOUND
103	S4# 75	2SEC BUTYL-4,6-DINITROPHE	211	4	NOT FOUND
104	S4# 80	C647 CARBAZOLE	167	4	NOT FOUND
105	S4# 85	C650 DI-N-BUTYLPHTHALATE	149	A BB 1226 4	2068. 1.247 0.572
106	S4#100	4-NITROQUINOLINE-1-OXIDE	190	4	NOT FOUND
107	S4#105	METHAPYRILENE	58	4	NOT FOUND
108	S4#106	ISODRIN	193	4	NOT FOUND
109	S4#110	C655 FLUORANTHENE	202	4	NOT FOUND
110	S4#120	CHLOROBENZILATE	139	4	NOT FOUND

111 S5# 10 C710 BENZIDINE	184	5	NOT FOUND		
112 S5# 15 C715 PYRENE	202	5	NOT FOUND		
113 S5# 20 ARAMITE (#1)	185	5	NOT FOUND		
114 S5# 25 ARAMITE (#2)	185	5	NOT FOUND		
115 S5# 30 P-DIMETHYLAMINOAZOBENZEN	120	5	NOT FOUND		
116 S5# 35 3,3'-DIMETHYLBENZIDINE	212	5	NOT FOUND		
117 S5# 37 KEPONE	272	5	NOT FOUND		
118 S5# 40 C720 BUTYLBENZYLPHTHALAT	149 A BB	1487 5		344. 0.685	0.280
119 S5# 45 2-ACETYLAMINOFLUORENE	181	5	NOT FOUND		
120 S5# 50 C730 BENZO(A)ANTHRACENE	228	5	NOT FOUND		
121 S5# 55 C725 3,3'-DICHLOROBENZID	252	5	NOT FOUND		
122 S5# 60 C740 CHRYSENE	228	5	NOT FOUND		
123 S5# 65 C745 BIS(2-ETHYLHEXYL)PH	149 A BB	1603 5		3064. 0.980	1.745
124 S5# 85 3-METHYLCHOLANTHRENE	268	5	NOT FOUND		
125 S6# 10 C760 DI-N-OCTYL PHTHALAT	149	6	NOT FOUND		
126 S6# 15 C765 BENZO(B)FLUORANTHEN	252	6	NOT FOUND		
127 S6# 20 7,12-DIMETHYLBENZANTHRAC	256	6	NOT FOUND		
128 S6# 25 C770 BENZO(K)FLUORANTHEN	252	6	NOT FOUND		
129 S6# 30 HEXACHLOROPHENE	196	6	NOT FOUND		
130 S6# 35 C775 BENZO(A)PYRENE	252	6	NOT FOUND		
131 S6# 55 C780 INDENO(1,2,3-CD)PYR	276	6	NOT FOUND		
132 S6# 60 C785 DIBENZ(A,H)ANTHRACE	278	6	NOT FOUND		
133 S6# 65 C790 BENZO(G,H,I)PERYLEN	276	6	NOT FOUND		

QUANTERRA Internal Standard Check

Standard Filename: ST16980831

Analyzed: 08/31/98 12:31

Standard	IS# 1			IS# 2			IS# 3		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	28649.		3.78	100432.		5.90	54347.		9.21
Upper Limit	57297. =+200		4.28	200864. =+200		6.40	108695. =+200		9.71
Lower Limit	14324. =+ 50		3.28	50216. =+ 50		5.40	27174. =+ 50		8.71
Filename									
1 30068104	33946. = 118		3.79	129608. = 129		5.88	68092. = 125		9.21

IS# 1 = C130 1,4-DICHLOROBENZENE-D4

IS# 2 = C140 NAPHTHALENE-D8

IS# 3 = C150 ACENAPHTHENE-D10

* - indicates an I.S. is outside QC limit(s)

Standard	IS# 4			IS# 5			IS# 6		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	98520.		12.08	62289.		17.26	50750.		19.83
Upper Limit	197039. =+200		12.58	124578. =+200		17.76	101500. =+200		20.33
Lower Limit	49260. =+ 50		11.58	31145. =+ 50		16.76	25375. =+ 50		19.33
Filename									
1 30068104	115982. = 117		12.07	71692. = 115		17.25	58915. = 116		19.82

IS# 4 = C160 PHENANTHRENE-D10

IS# 5 = C170 CHRYSENE-D12

IS# 6 = C175 PERYLENE-D12

* - indicates an I.S. is outside QC limit(s)

Data Reduced by: 2 Date: 9/17/98
Data Reviewed by: 1 Date: 9/17/98

Data File: 30068104

QUANTERRA GC/MS TIC REPORT (Part 1)

Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
Analyst: DAT Date Analyzed: 08/31/98 20:30
Run Factor: 3.00

Concentration in Sample		
# SCAN	(UG/A)	CAS #
1 170	260.	123-42-2
2-PENTANONE, 4-HYDROXY-4-METHYL-		or isomer
2 222	260.	111-84-2
NONANE or isomer		
3 286	22.	100-52-7
BENZALDEHYDE or isomer		
4 932	23.	74381-40-1
PROPANOIC ACID, 2-METHYL-, 1-(1,1-DIMETHYLETHYL)-2-METHYL-1,3-PROPANE!		
		unknown 5700
5 1235	29.	57-10-3
HEXADECANOIC ACID or isomer		
6 1333	54.	00-00-0
UNKNOWN		5700
		↓
7 1357	48.	00-00-0
UNKNOWN		
8 1466	12.	62108-23-0
DECANE, 2,5,6-TRIMETHYL-		or isomer
9 1502	21.	00-00-0
UNKNOWN		5700
		↓
10 1509	42.	00-00-0
UNKNOWN		

11 1528 62. 00-00-0 5700
UNKNOWN

12 1586 160. 74685-30-6
5-EICOSENE, (E)- or Isomer

13 1661 55. 00-00-0 5700
UNKNOWN

14 1686 40. 00-00-0
UNKNOWN

~~15 1704 21. . .~~
~~HEXYL 1-OCTANOL~~ unknown 5700

16 1741 20. 00-00-0
UNKNOWN

17 1799 58. 00-00-0
UNKNOWN

18 1804 73. 00-00-0
UNKNOWN

19 1914 75. 1653-33-4 5720000
~~4-TETRADECANOL~~ unknown alkene

20 1932 74. 00-00-0 5700
UNKNOWN

21 2012 46. 74685-29-3 5720000
~~9-EICOSENE, (E)-~~ unknown alkene

22 2049 40. 00-00-0 5700
UNKNOWN

Use!
23 2056
UNKNOWN

29.

00-00-0

DATA FILE: 30068104

QUANTERRA GC/MS TIC REPORT (Part 2)

CONCENTRATION = AREA(TIC)*CONC(IS)/AREA(IS)

#	FIT	PURITY	INT.		RT	RRT	AREA	HEIGHT	AMOUNT		LIB	LIB #
			STD.						AS ANALYZED	(UG/ML)		
1	954	918	1		1:53	0.321	708264.	317440.	87.787		NB	3241.
2	987	933	1		2:28	0.420	697660.	287744.	86.472		NB	5159.
3	967	920	1		3:10	0.541	60288.	15312.	7.472		NB	2017.
4	930	779	2		10:21	1.126	60114.	20417.	7.544		NB	40501.
5	957	752	3		13:44	1.137	77738.	20052.	9.738		NB	35182.
6	764	558	4		14:49	0.859	101780.	10208.	17.848		UK	1.
7	816	655	4		15:05	0.875	92128.	6768.	16.156		UK	1.
8	905	707	4		16:18	0.945	23692.	3066.	4.155		NB	19015.
9	711	295	4		16:42	0.968	40720.	7264.	7.141		UK	1.
10	842	583	4		16:47	0.973	79168.	11824.	13.883		UK	1.
11	873	536	4		16:59	0.985	116992.	10160.	20.516		UK	1.
12	983	808	4		17:38	1.023	313592.	57920.	54.992		NB	39516.
13	788	502	4		18:28	1.071	104880.	10496.	18.392		UK	1.
14	621	481	5		18:45	0.946	65416.	5616.	13.475		UK	1.
15	960	726	5		18:57	0.956	34164.	7268.	7.037		NB	26409.
16	866	568	5		19:21	0.977	33084.	8195.	6.815		UK	1.
17	678	326	5		20:00	1.010	93760.	12192.	19.313		UK	1.
18	865	474	5		20:03	1.012	118128.	12992.	24.333		UK	1.
19	982	754	5		21:17	1.074	121984.	25021.	25.127		NB	26416.
20	910	435	5		21:29	1.084	119248.	11408.	24.563		UK	1.
21	945	692	5		22:22	1.129	73644.	13181.	15.170		NB	39515.
22	814	463	5		22:47	1.150	64720.	8112.	13.331		UK	1.
23	717	429	5		22:52	1.154	46608.	7984.	9.601		UK	1.

Library Search Data: 30068104 # 170 Base m/z: 43
 08/31/98 20:30:00 + 1:53 Cali: 30068104 # 3 RIC: 236032.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 495 matched at least 6 of the 16 largest peaks in the unknown

Rank In.	Name
1	3241 2-PENTANONE, 4-HYDROXY-4-METHYL-
2	5791 1,3-DIOXOLANE-2-METHANOL, 2,4-DIMETHYL-
3	3366 2-PENTANOL, 2,4-DIMETHYL-
4	3288 ACETIC ACID, 1,1-DIMETHYLETHYL ESTER
5	919 TERT-BUTYL HYDROPEROXIDE
6	5594 2-PROPANOL, 2-NITROSO-, ACETATE (ESTER)
7	3349 2-HEXANOL, 2-METHYL-
8	11889 2-HEXANONE, 6-(ACETYLOXY)-
9	1576 OXIRANE, TETRAMETHYL-

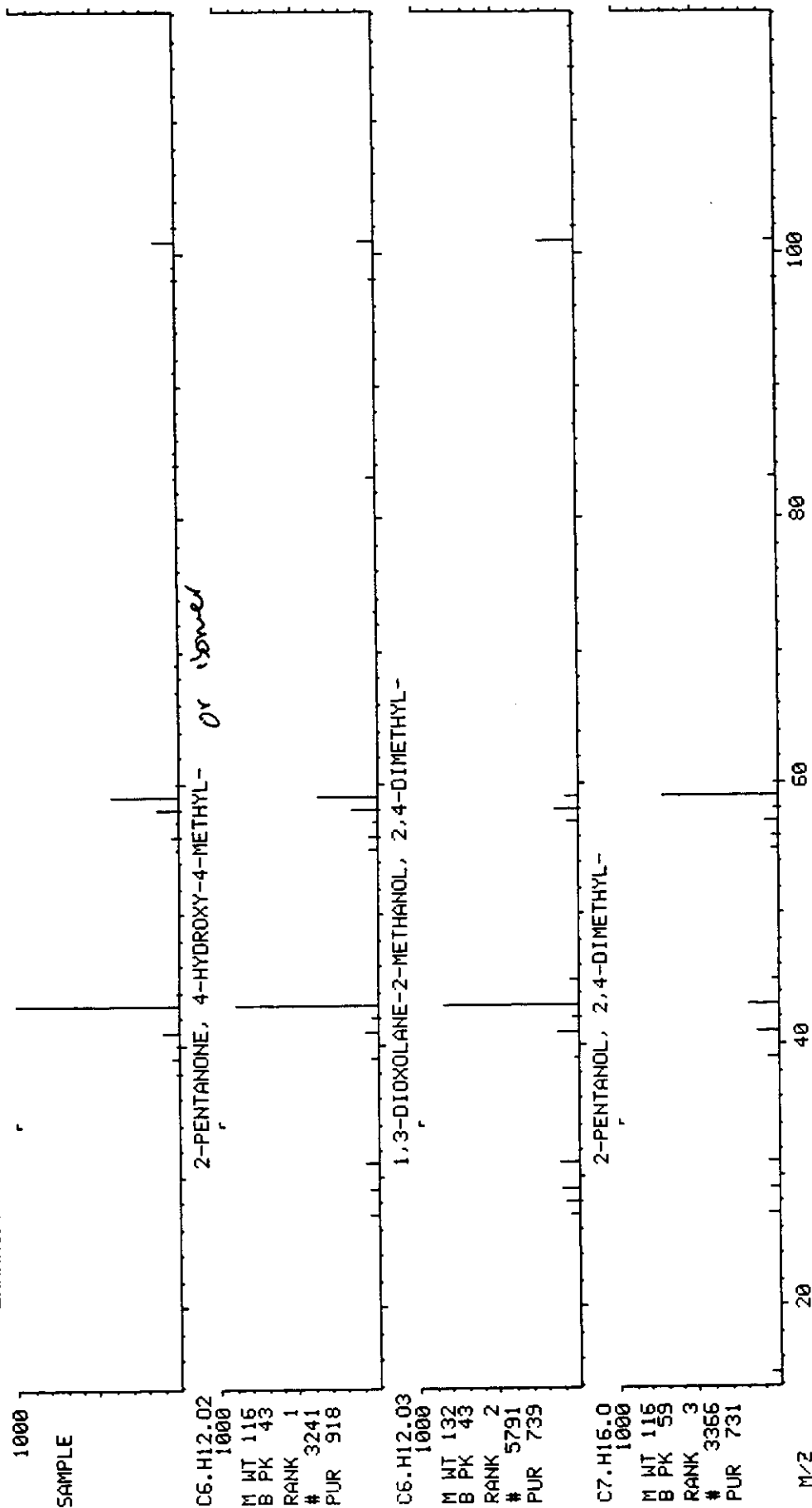
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C6.H12.O2	116	43	918	954	941
2	C6.H12.O3	132	43	739	813	805
3	C7.H16.O	116	59	731	753	749
4	C6.H12.O2	116	43	731	778	745
5	C4.H10.O2	90	59	723	870	742
6	C5.H9.O3.N	131	43	705	817	793
7	C7.H16.O	116	59	686	719	722
8	C8.H14.O3	158	43	651	681	711
9	C6.H12.O	100	59	635	781	676

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	123-42-2
2	---	---	---	---	53951-43-2
3	---	---	---	---	625-06-9
4	---	---	---	---	540-88-5
5	---	---	---	---	75-91-2
6	---	---	---	---	6931-04-0
7	---	---	---	---	625-23-0
8	---	---	---	---	4305-26-4
9	---	---	---	---	5076-20-0

DATA: 30068104 # 170
CALI: 30068104 # 3

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 20:30:00 + 1:53
SAMPLE: T-MMS-FB 1/35A/1ML
CONDS.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068104 # 222 Base m/z: 43
 08/31/98 20:30:00 + 2:28 Cali: 30068104 # 3 RIC: 214784.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 921 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name

1 5159 NONANE
 2 5151 HEPTANE, 3,4-DIMETHYL-
 3 5141 HEPTANE, 2,4-DIMETHYL-
 4 3086 HEXANE, 2,4-DIMETHYL-
 5 19015 DECANE, 2,5,6-TRIMETHYL-
 6 11607 UNDECANE
 7 5154 HEXANE, 4-ETHYL-2-METHYL-
 8 8085 HEPTANE, 2,4,6-TRIMETHYL-
 9 5144 HEPTANE, 4-ETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C9.H20	128	43	933	987	933
2	C9.H20	128	43	901	933	901
3	C9.H20	128	43	883	941	884
4	C8.H18	114	43	880	960	891
5	C13.H28	184	57	873	918	883
6	C11.H24	156	43	866	931	902
7	C9.H20	128	57	860	920	860
8	C10.H22	142	43	857	904	872
9	C9.H20	128	43	857	910	865

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	111-84-2
2	—	—	—	—	922-28-1
3	—	—	—	—	2213-23-2
4	—	—	—	—	589-43-5
5	—	—	—	—	62108-23-0
6	—	—	—	—	1120-21-4
7	—	—	—	—	3074-75-7
8	—	—	—	—	2613-61-8
9	—	—	—	—	2216-32-2

BASE M/Z: 43
RIC: 214784.

DATA: 30058104 # 222
CALI: 30058104 # 3

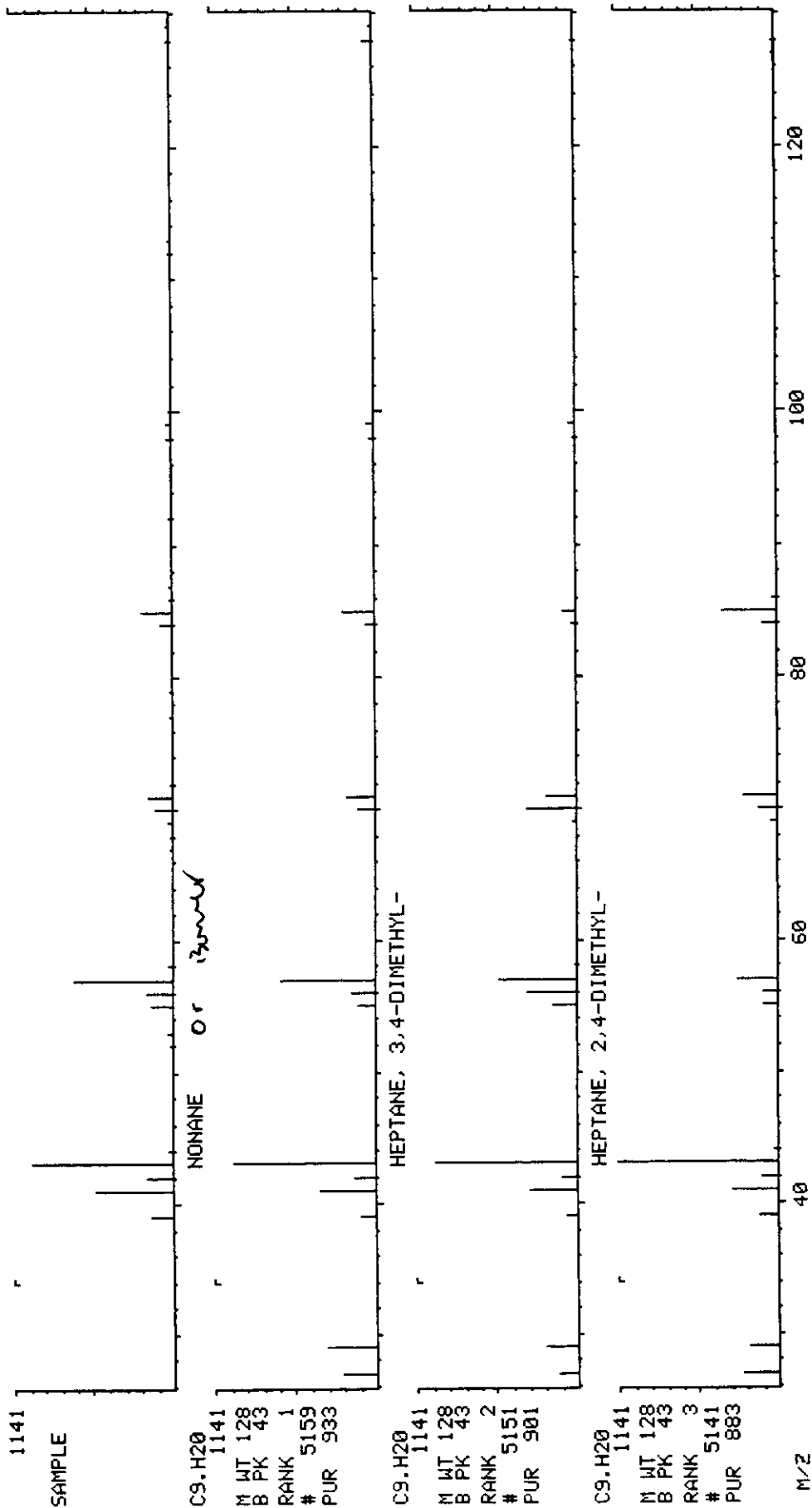
MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 20:30:00 + 2:28

SAMPLE: T-MM5-FB 1/35A/1ML INST. ID: F16

COND5.: UG/ML *1ML *100%/100% *(NA/NA) 1/35A NA M

ENHANCED (S 15B 2N 0T)



Library Search Data: 30068104 # 286 Base m/z: 77
 08/31/98 20:30:00 + 3:11 Cali: 30068104 # 3 RIC: 12928.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 354 matched at least 5 of the 16 largest peaks in the unknown

Rank In. Name
 1 2017 BENZALDEHYDE
 2 7284 BENZOYL CHLORIDE
 3 6466 ETHANONE, 2-HYDROXY-1-PHENYL-
 4 2014 2,4-HEPTADIEN-6-YNAL, (E,E)-
 5 9645 BENZOYLFORMIC ACID
 6 13112 BENZOYL ISOTHIOCYANATE
 7 6857 BENZENECARBOTHIOIC ACID
 8 6111 1,2-BENZENEDICARBOXALDEHYDE
 9 13359 BENZENEACETIC ACID, .ALPHA.-OXO-, METHYL ESTER

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C7.H6.O	106	77	920	967	941
2	C7.H5.O <u>CL</u>	140	105	809	871	826
3	C8.H8.O2	136	105	806	843	822
4	C7.H6.O	106	105	805	844	809
5	C8.H6.O3	150	105	779	849	803
6	C8.H5.O.N.S	163	105	767	818	785
7	C7.H6.O.S	138	77	760	846	807
8	C8.H6.O2	134	105	759	806	915
9	C9.H8.O3	164	105	758	834	786

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	100-52-7
2	—	—	—	—	98-88-4
3	—	—	—	—	582-24-1
4	—	—	—	—	7200-04-6
5	—	—	—	—	611-73-4
6	—	—	—	—	532-55-8
7	—	—	—	—	98-91-9
8	—	—	—	—	643-79-8
9	—	—	—	—	15206-55-0

DATA: 30068104 # 286
CALI: 30068104 # 3
BASE M/Z: 77
RIC: 12928.

MID LIBRARY SEARCH (LIBRARYNB)

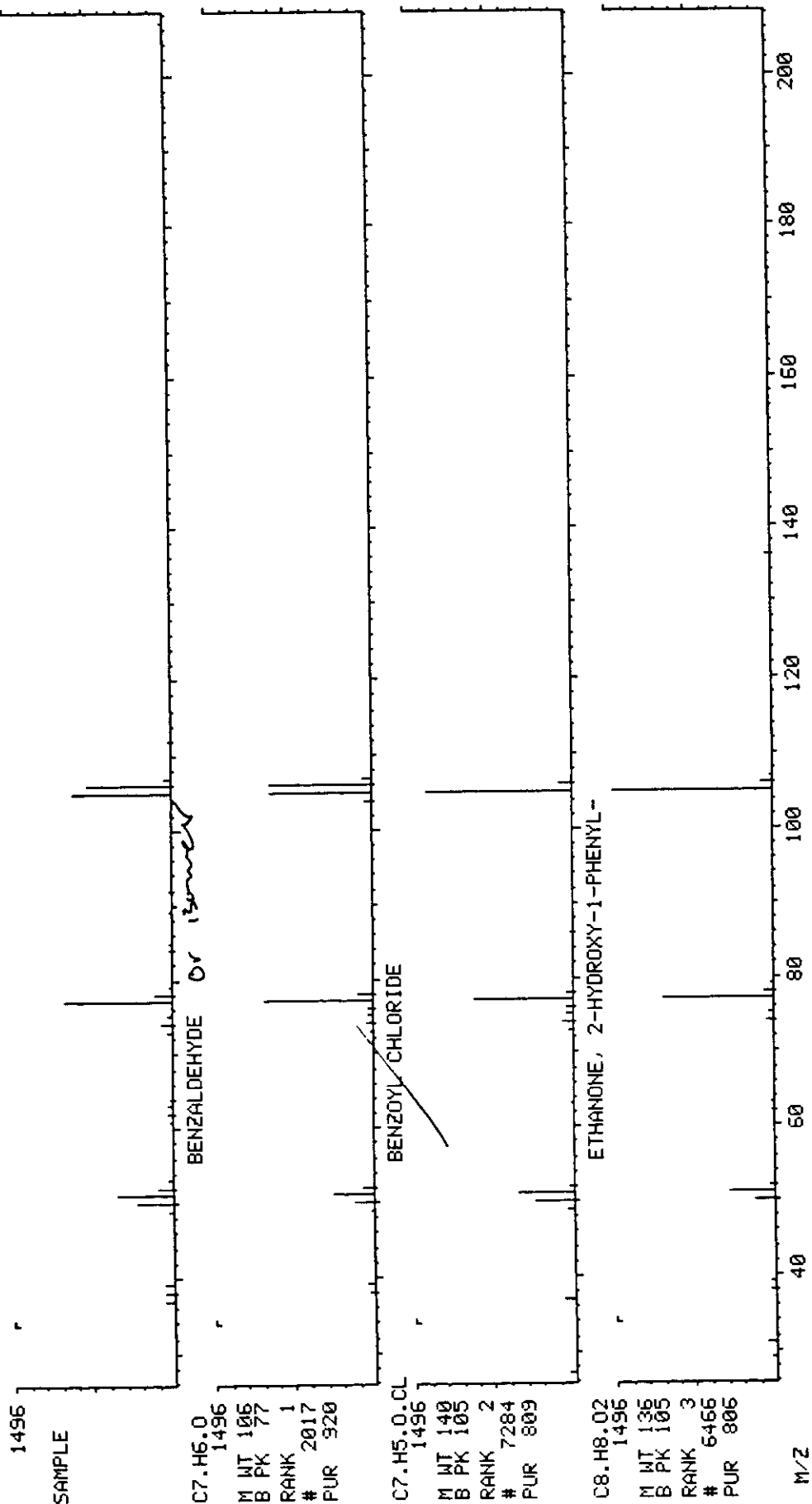
08/31/98 20:30:00 + 3:11

SAMPLE: T-MMS-FB 1/35A/1ML

COND.S.: UG/ML *1ML *100%/100% *(NA/NA) /1/35A NA M

ENHANCED (S 15B 2N 0T)

INST. ID: F16



[Handwritten signature]

Library Search Data: 30068104 # 932 Base m/z: 71
 08/31/98 20:30:00 + 10:22 Cali: 30068104 # 3 RIC: 17504.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 410 matched at least 6 of the 16 largest peaks in the unknown

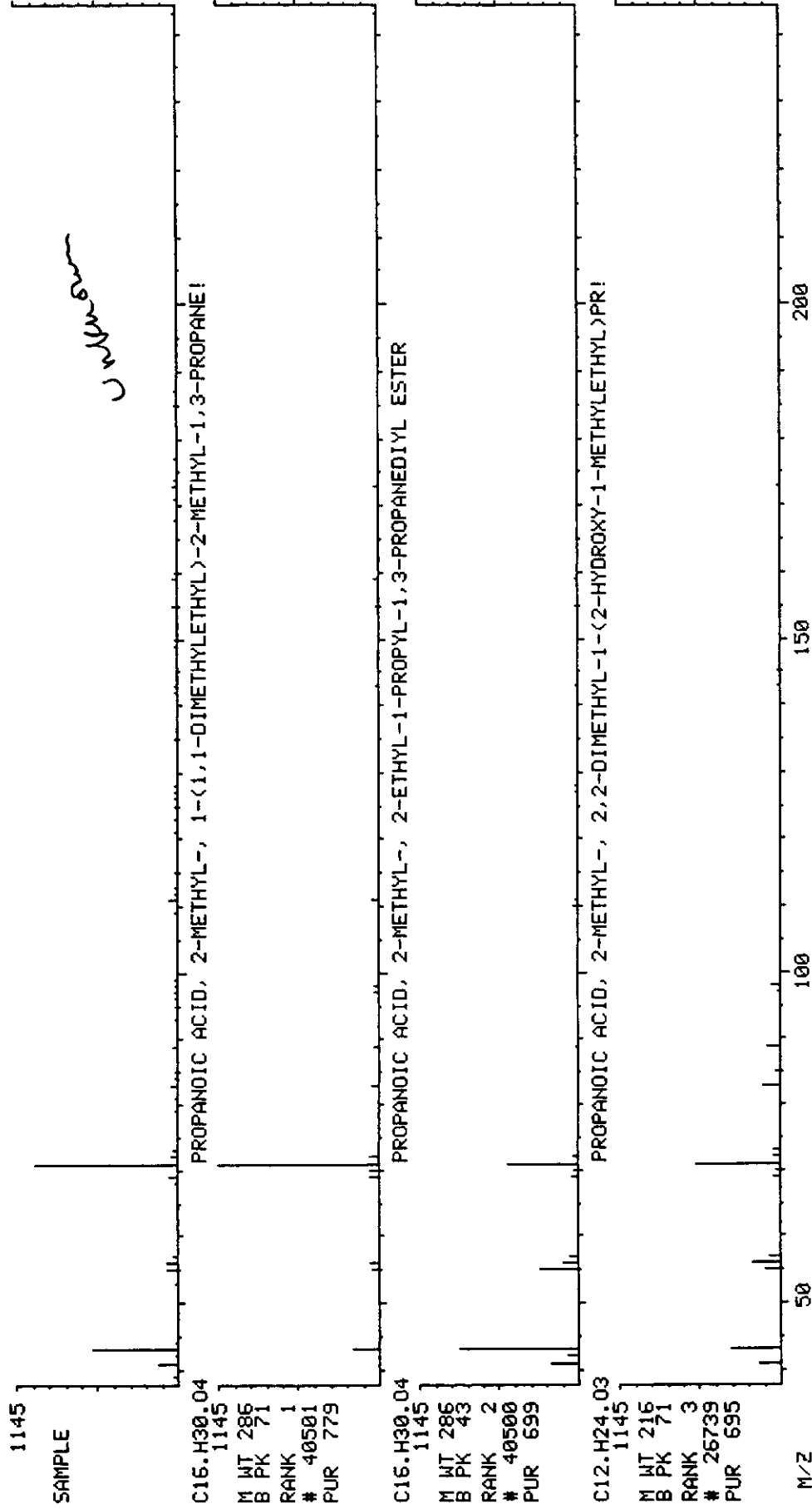
Rank In.	Name
1 40501	PROPANOIC ACID, 2-METHYL-, 1-(1,1-DIMETHYLETHYL)-2-METHYL-1,3-PROPA*
2 40500	PROPANOIC ACID, 2-METHYL-, 2-ETHYL-1-PROPYL-1,3-PROPANEDIYL ESTER
3 26739	PROPANOIC ACID, 2-METHYL-, 2,2-DIMETHYL-1-(2-HYDROXY-1-METHYLETHYL)*
4 26740	PROPANOIC ACID, 2-METHYL-, 3-HYDROXY-2,4,4-TRIMETHYLPENTYL ESTER
5 4625	1-HEXENE, 3,4,5-TRIMETHYL-
6 11945	PROPANOIC ACID, 2-METHYL-, 1-METHYLBUTYL ESTER
7 19470	BUTANOIC ACID, 1-METHYLHEXYL ESTER
8 26737	PROPANOIC ACID, 2-METHYL-, 2-ETHYL-3-HYDROXYHEXYL ESTER
9 14787	1-NONENE, 4,6,8-TRIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C16.H30.O4	286	71	779	930	800
2	C16.H30.O4	286	43	699	866	726
3	C12.H24.O3	216	71	695	883	714
4	C12.H24.O3	216	71	638	794	643
5	C9.H18	126	43	575	800	658
6	C9.H18.O2	158	71	539	799	610
7	C11.H22.O2	186	71	539	755	578
8	C12.H24.O3	216	71	528	691	551
9	C12.H24	168	43	519	715	531

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	74381-40-1
2	—	—	—	—	74367-30-9
3	—	—	—	—	74367-33-2
4	—	—	—	—	74367-34-3
5	—	—	—	—	56728-10-0
6	—	—	—	—	54340-93-1
7	—	—	—	—	39026-94-3
8	—	—	—	—	74367-31-0
9	—	—	—	—	54410-98-9

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:30:00 + 10:22
 SAMPLE: T-MM5-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA) 1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30058104 # 932
 CALI: 30058104 # 3
 BASE M/Z: 71
 RIC: 17504.



Library Search Data: 30068104 #1235 Base m/z: 43
 08/31/98 20:30:00 + 13:44 Cali: 30068104 # 3 RIC: 16144.
 Sample: T-MMS-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

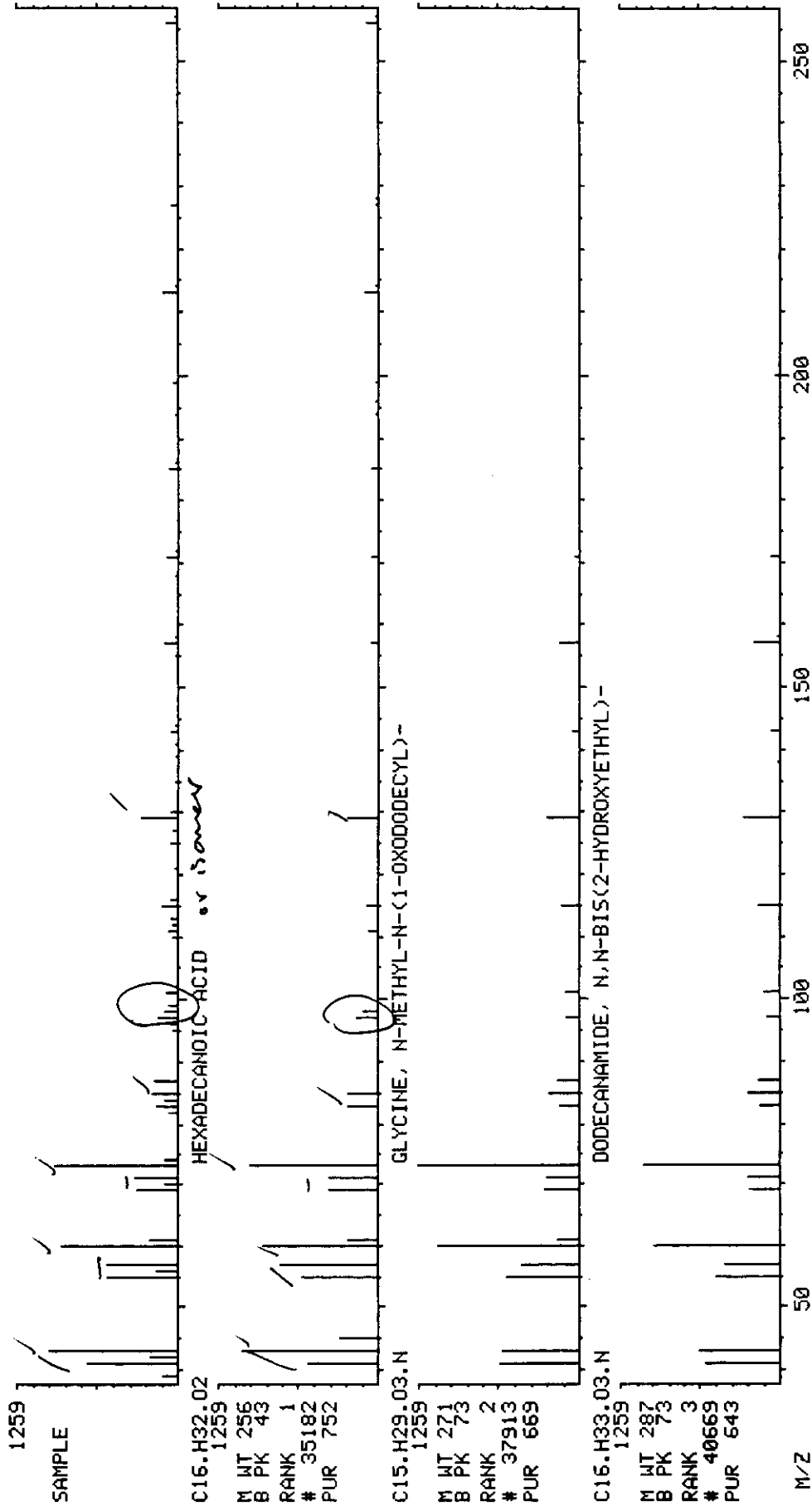
62231 spectra in LIBRARYNB searched for maximum PURITY
 912 matched at least 6 of the 16 largest peaks in the unknown

Rank In.	Name
1	35182 HEXADECANOIC ACID
2	37913 GLYCINE, N-METHYL-N-(1-OXODODECYL)-
3	40669 DODECANAMIDE, N,N-BIS(2-HYDROXYETHYL)-
4	29642 TETRADECANOIC ACID
5	40184 OCTADECANOIC ACID
6	22949 DODECANOIC ACID
7	26351 TRIDECANOIC ACID
8	19469 UNDECANOIC ACID
9	32392 PENTADECANOIC ACID

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C16.H32.O2	256	43	752	957	774
2	C15.H29.O3.N	271	73	669	982	669
3	C16.H33.O3.N	287	73	643	966	655
4	C14.H28.O2	228	73	625	898	673
5	C18.H36.O2	284	43	607	767	770
6	C12.H24.O2	200	60	602	902	660
7	C13.H26.O2	214	73	595	878	616
8	C11.H22.O2	186	60	586	941	598
9	C15.H30.O2	242	43	573	801	688

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	57-10-3
2	—	—	—	—	97-78-9
3	—	—	—	—	120-40-1
4	—	—	—	—	544-63-8
5	—	—	—	—	57-11-4
6	—	—	—	—	143-07-7
7	—	—	—	—	638-53-9
8	—	—	—	—	112-37-8
9	—	—	—	—	1002-84-2

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:30:00 + 13:44
 SAMPLE: T-MM5-FB 1/35A/1ML
 COND5.: UG/ML *1ML *1002/1002 *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068104 #1235
 CALI: 30068104 # 3
 BASE M/Z: 43
 RIC: 16144.



Library Search Data: 30068104 #1333 Base m/z: 59
 08/31/98 20:30:00 + 14:50 Cal: 30068104 # 3 RIC: 7024.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 636 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 24247 2-PROPANOL, 1-[2-(2-METHOXY-1-METHYLETHOXY)-1-METHYLETHOXY]-
 2 16192 2-PROPANOL, 1-[1-METHYL-2-(2-PROPENYLOXY)ETHOXY]-
 3 3502 BUTANOIC ACID, 3-HYDROXY-3-METHYL-
 4 5522 2-HEXANOL, 2,3-DIMETHYL-
 5 8460 SILANE, OCTYL-
 6 20626 2-PROPANOL, 1,1'-[(1-METHYL-1,2-ETHANEDIYL)BIS(OXY)]BIS-
 7 3522 2,3-BUTANEDIOL, 2,3-DIMETHYL-
 8 3000 OXIRANE, 2,2-DIMETHYL-3-PROPYL-
 9 8506 4,5-DIMETHYL-3-HEPTANOL

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C10.H22.O4	206	59	558	764	624
2	C9.H18.O3	174	59	534	728	574
3	C5.H10.O3	118	59	532	771	618
4	C8.H18.O	130	59	526	821	547
5	C8.H20.SI	144	59	508	866	521
6	C9.H20.O4	192	59	503	720	563
7	C6.H14.O2	118	59	500	885	516
8	C7.H14.O	114	43	493	813	524
9	C9.H20.O	144	59	482	815	550

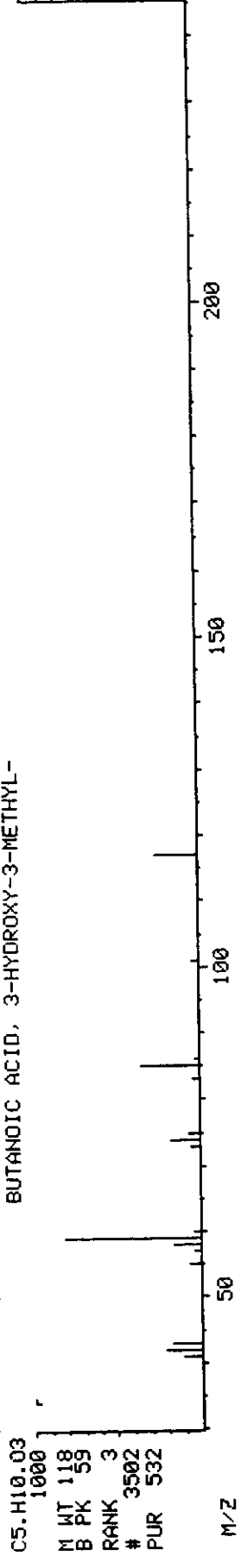
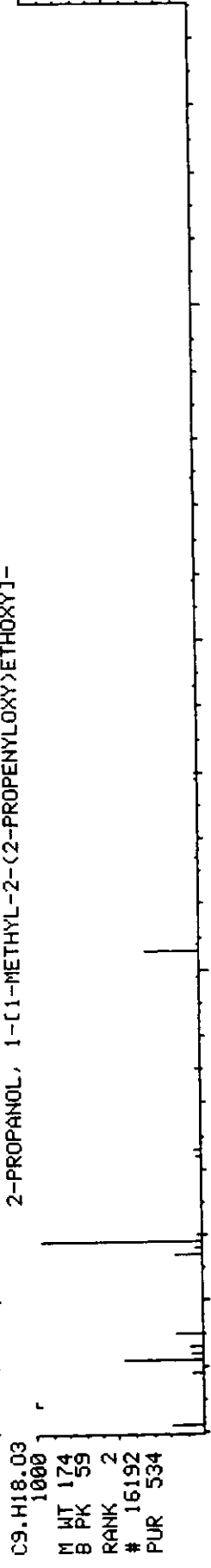
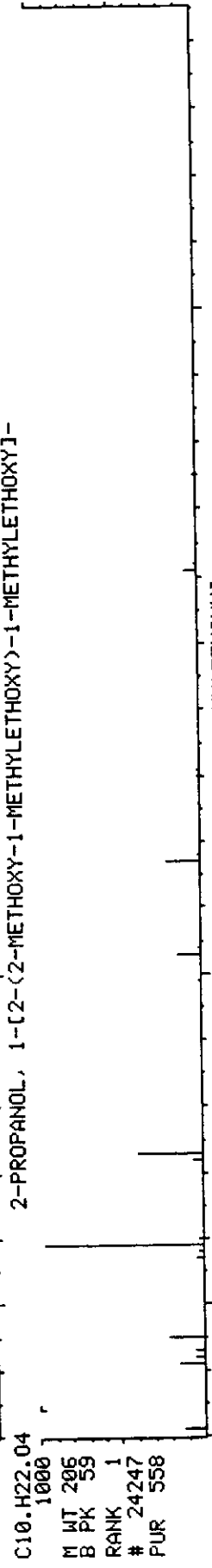
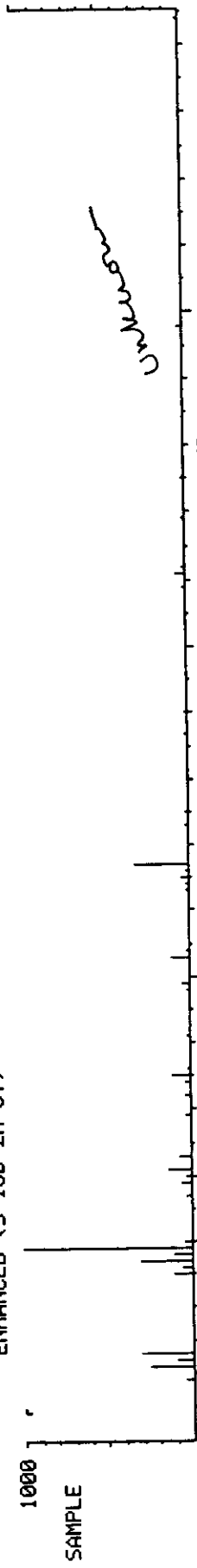
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	20324-33-8
2	—	—	—	—	55956-25-7
3	—	—	—	—	625-08-1
4	—	—	—	—	19550-03-9
5	—	—	—	—	871-92-1
6	—	—	—	—	1638-16-0
7	—	—	—	—	76-09-5
8	—	—	—	—	17612-35-0
9	—	—	—	—	-

DATA: 30068104 #1333
CALI: 30068104 # 3

BASE M/Z: 59
RIC: 7024.

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 20:30:00 + 14:50
SAMPLE: T-MMS-FB 1/35A/1ML
CONDOS.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 158 2N 0T)

INST. ID: F16



Library Search Data: 30068104 #1357 Base m/z: 59
 08/31/98 20:30:00 + 15:06 Cali: 30068104 # 3 RIC: 4152.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 803 matched at least 5 of the 16 largest peaks in the unknown

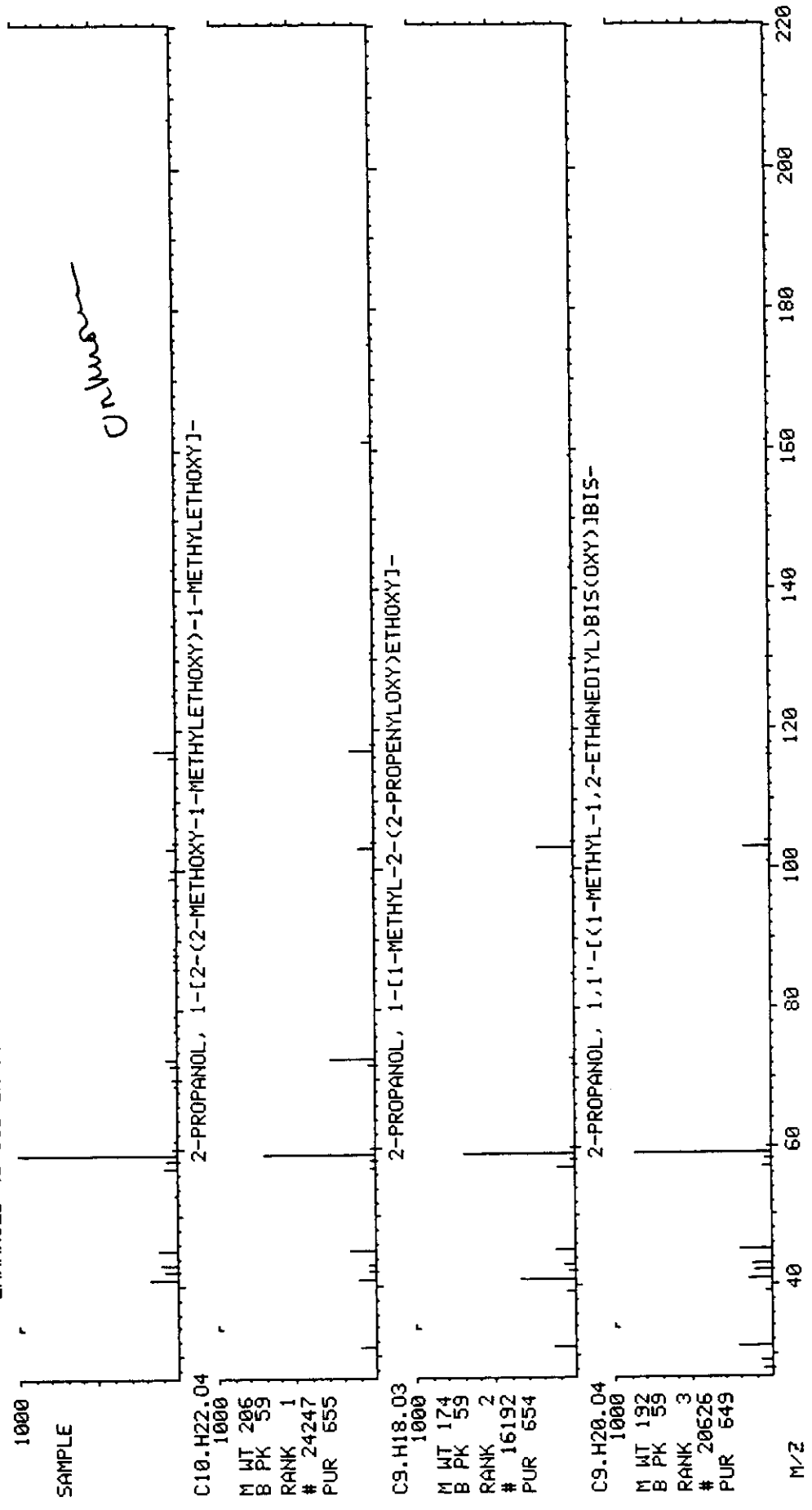
Rank In. Name
 1 24247 2-PROPANOL, 1-(2-(2-METHOXY-1-METHYLETHOXY)-1-METHYLETHOXY)-
 2 16192 2-PROPANOL, 1-[1-METHYL-2-(2-PROPENYLOXY)ETHOXY]-
 3 20626 2-PROPANOL, 1,1'-[(1-METHYL-1,2-ETHANEDIYL)BIS(OXY)]BIS-
 4 9229 2-PROPANOL, 1-(2-METHOXYPROPOXY)-
 5 9228 1-PROPANOL, 2-(2-METHOXYPROPOXY)-
 6 36537 2,5,8,11-TETRAOXATETRADECAN-13-OL, 4,7,10-TRIMETHYL-
 7 9235 1-PROPANOL, 2-(2-METHOXY-1-METHYLETHOXY)-
 8 9234 2-PROPANOL, 1-(2-METHOXY-1-METHYLETHOXY)-
 9 5786 PROPANOIC ACID, 2-HYDROXY-2-METHYL-, ETHYL ESTER

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C10.H22.O4	206	59	655	816	728
2	C9.H18.O3	174	59	654	807	681
3	C9.H20.O4	192	59	649	858	674
4	C7.H16.O3	148	59	643	942	650
5	C7.H16.O3	148	59	624	853	652
6	C13.H28.O5	264	59	617	755	732
7	C7.H16.O3	148	59	585	768	594
8	C7.H16.O3	148	59	584	749	595
9	C6.H12.O3	132	59	581	861	606

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	20324-33-8
2	—	—	—	—	55956-25-7
3	—	—	—	—	1638-16-0
4	—	—	—	—	13429-07-7
5	—	—	—	—	13588-28-8
6	—	—	—	—	20324-34-9
7	—	—	—	—	55956-21-3
8	—	—	—	—	20324-32-7
9	—	—	—	—	80-55-7

DATA: 30068104 #1357
CALI: 30068104 # 3
BASE M/Z: 59
RIC: 4152.

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 20:30:00 + 15:06
SAMPLE: T-MM5-FB 1/35A/1ML INST. ID: F16
CONDS.: UG/ML *1ML *100%/100% *(NA/NA) >1/35A NA M
ENHANCED (S 158 2N 0T)



Library Search Data: 30068104 #1466 Base m/z: 57
 08/31/98 20:30:00 + 16:18 Cal: 30068104 # 3 RIC: 2364.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 603 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 19015 DECANE, 2,5,6-TRIMETHYL-
 2 19056 UNDECANE, 2,7-DIMETHYL-
 3 11612 NONANE, 2,5-DIMETHYL-
 4 19006 DECANE, 6-ETHYL-2-METHYL-
 5 11602 OCTANE, 2,4,6-TRIMETHYL-
 6 19045 DECANE, 2,3,8-TRIMETHYL-
 7 19000 UNDECANE, 5,6-DIMETHYL-
 8 15969 HYDROXYLAMINE, O-DECYL-
 9 19523 1-DECANOL, 2-ETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H28	184	57	707	905	730
2	C13.H28	184	43	700	872	729
3	C11.H24	156	57	699	882	732
4	C13.H28	184	57	695	877	722
5	C11.H24	156	57	693	898	711
6	C13.H28	184	57	692	881	713
7	C13.H28	184	57	678	889	697
8	C10.H23.O.N	173	43	676	842	756
9	C12.H26.O	186	57	674	837	766

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	62108-23-0
2	—	—	—	—	17301-24-5
3	—	—	—	—	17302-27-1
4	—	—	—	—	62108-21-8
5	—	—	—	—	62016-37-9
6	—	—	—	—	62238-14-6
7	—	—	—	—	17615-91-7
8	—	—	—	—	29812-79-1
9	—	—	—	—	21078-65-9

DATA: 30068104 #1466
CALI: 30068104 # 3

BASE M/Z: 57
RIC: 2364.

MID LIBRARY SEARCH (LIBRARYNB)

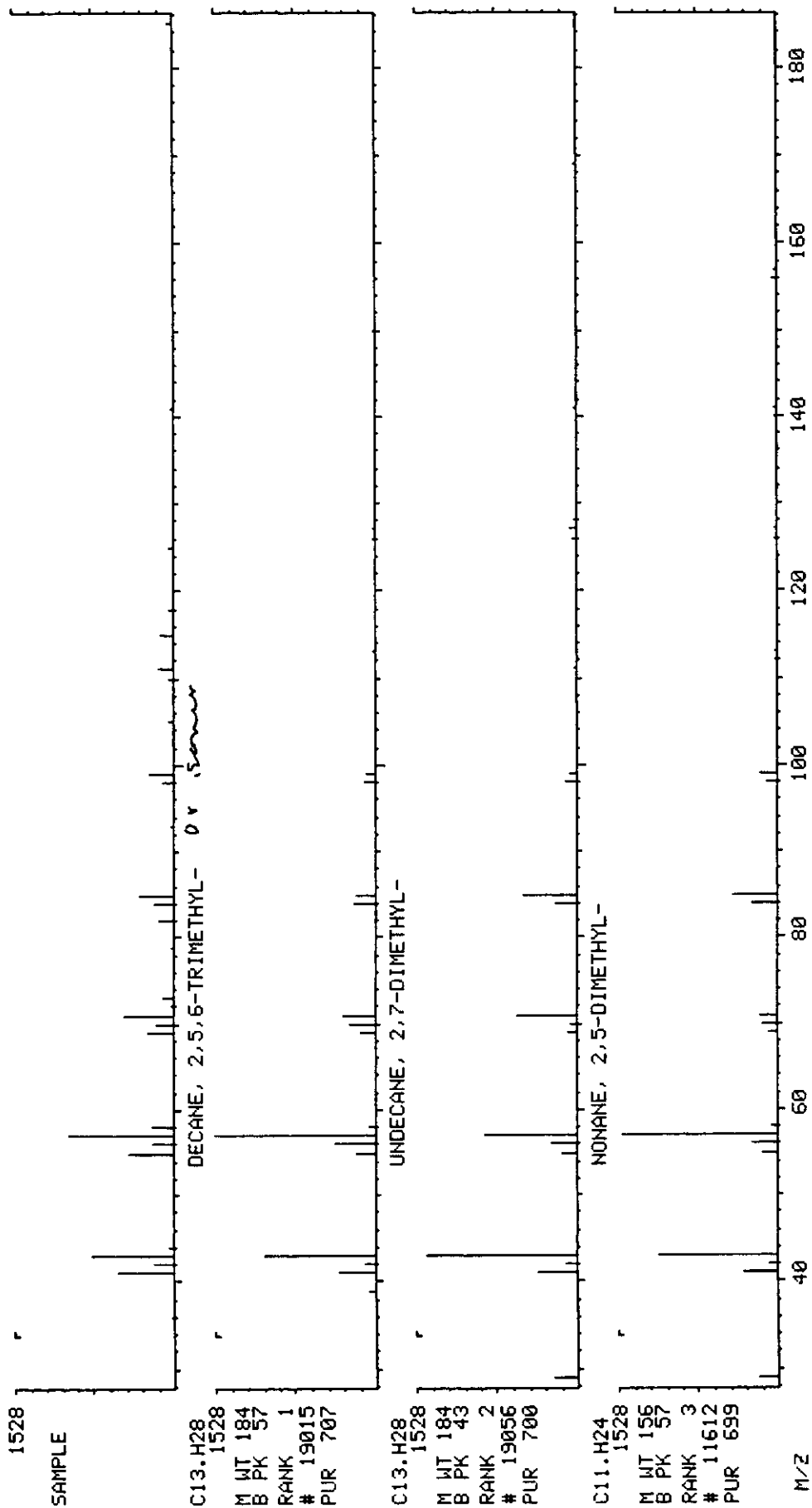
08/31/98 20:30:00 + 16:18

SAMPLE: T-MM5-FB 1/35A/1ML

COND5.: UG/ML *1ML *100%/100% *(NA/NA) 1/35A NA M

ENHANCED (S 158 2N 0T)

INST. ID: F16



Library Search Data: 30068104 #1502 Base m/z: 45
 08/31/98 20:30:00 + 16:42 Cali: 30068104 # 3 RIC: 2424.
 Sample: T-MMS-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

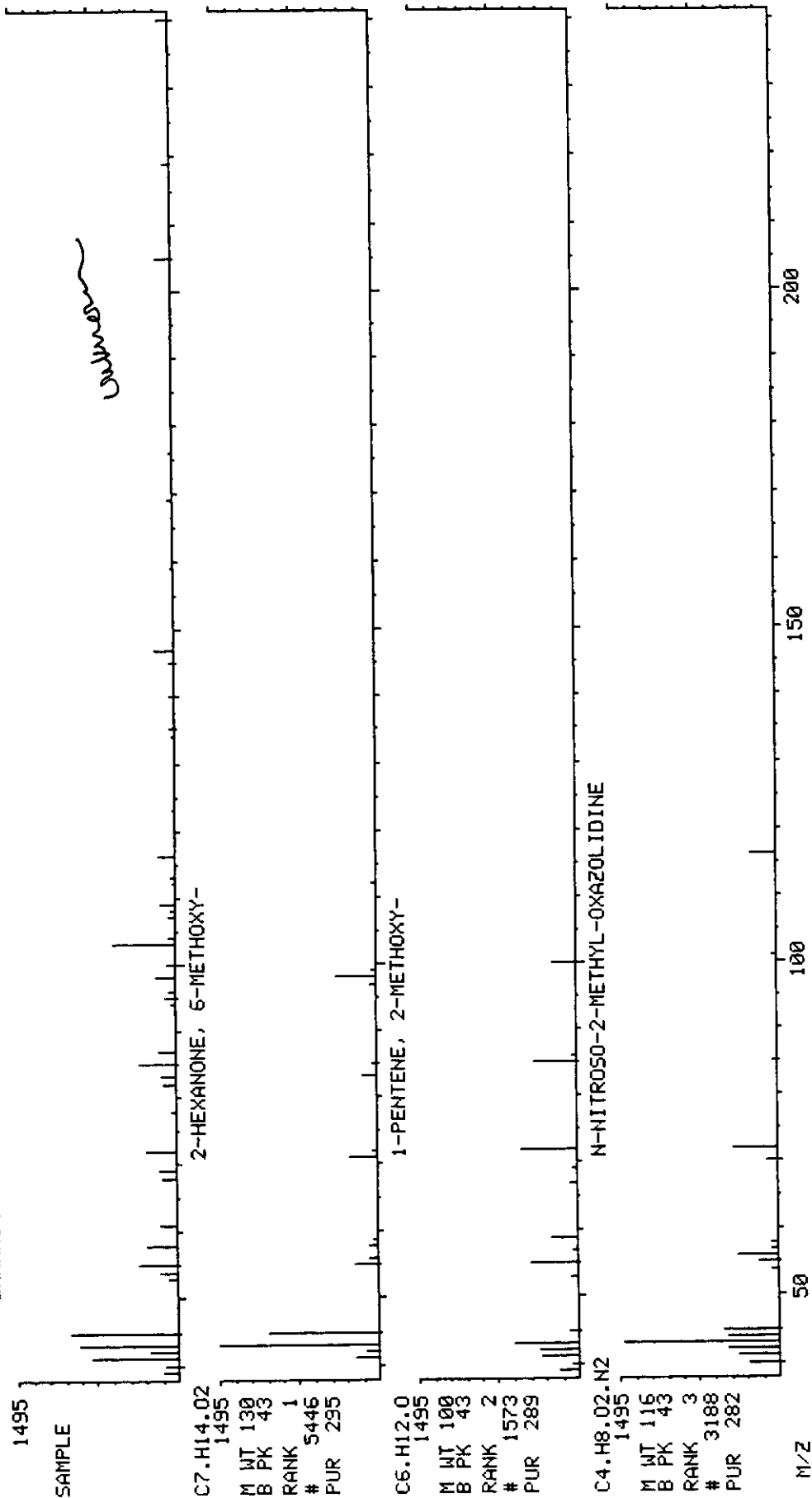
62231 spectra in LIBRARYNB searched for maximum PURITY
 384 matched at least 5 of the 16 largest peaks in the unknown

Rank In.	Name
1	5446 2-HEXANONE, 6-METHOXY-
2	1573 1-PENTENE, 2-METHOXY-
3	3188 N-NITROSO-2-METHYL-OXAZOLIDINE
4	1567 2-PENTENE, 2-METHOXY-
5	6418 2-PENTANOL, 3-CHLORO-4-METHYL-, (R@,S@)-(./-.)-
6	6417 2-PENTANOL, 3-CHLORO-4-METHYL-, (R@,R@)-(./-.)-
7	20219 1-PROPANOL, 2,2-BIS(METHOXYMETHYL)-, ACETATE
8	3339 1-PENTANOL, 2,3-DIMETHYL-
9	1789 2-HEXANOL

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C7.H14.O2	130	43	295	711	348
2	C6.H12.O	100	43	289	735	327
3	C4.H8.O2.N2	116	43	282	679	359
4	C6.H12.O	100	43	281	743	326
5	C6.H13.O.CL	136	45	276	783	309
6	C6.H13.O.CL	136	45	275	780	312
7	C9.H18.O4	190	43	269	564	410
8	C7.H16.O	116	43	268	660	371
9	C6.H14.O	102	45	264	764	293

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	29006-00-6
2	—	—	—	—	53119-70-3
3	—	—	—	—	39884-53-2
4	—	—	—	—	61142-47-0
5	—	—	—	—	74685-48-6
6	—	—	—	—	74685-47-5
7	—	—	—	—	20637-35-8
8	—	—	—	—	10143-23-4
9	—	—	—	—	626-93-7

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:30:00 + 16:42
 SAMPLE: T-NMS-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30068104 #1502
 CALI: 30068104 # 3
 BASE M/Z: 45
 RIC: 2424.



Library Search Data: 30068104 #1509 Base m/z: 59
 08/31/98 20:30:00 + 16:47 Cali: 30068104 # 3 RIC: 8624.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 570 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 3502 BUTANOIC ACID, 3-HYDROXY-3-METHYL-
 2 24247 2-PROPANOL, 1-[2-(2-METHOXY-1-METHYLETHOXY)-1-METHYLETHOXY]-
 3 3522 2,3-BUTANEDIOL, 2,3-DIMETHYL-
 4 5522 2-HEXANOL, 2,3-DIMETHYL-
 5 16192 2-PROPANOL, 1-[1-METHYL-2-(2-PROPENYLOXY)ETHOXY]-
 6 5506 2,3,3-TRIMETHYL-2-PENTANOL
 7 8460 SILANE, OCTYL-
 8 20626 2-PROPANOL, 1,1'-[(1-METHYL-1,2-ETHANEDIYL)BIS(OXY)]BIS-
 9 8805 1,3-DIOXAN-5-OL, 4,4,5-TRIMETHYL-

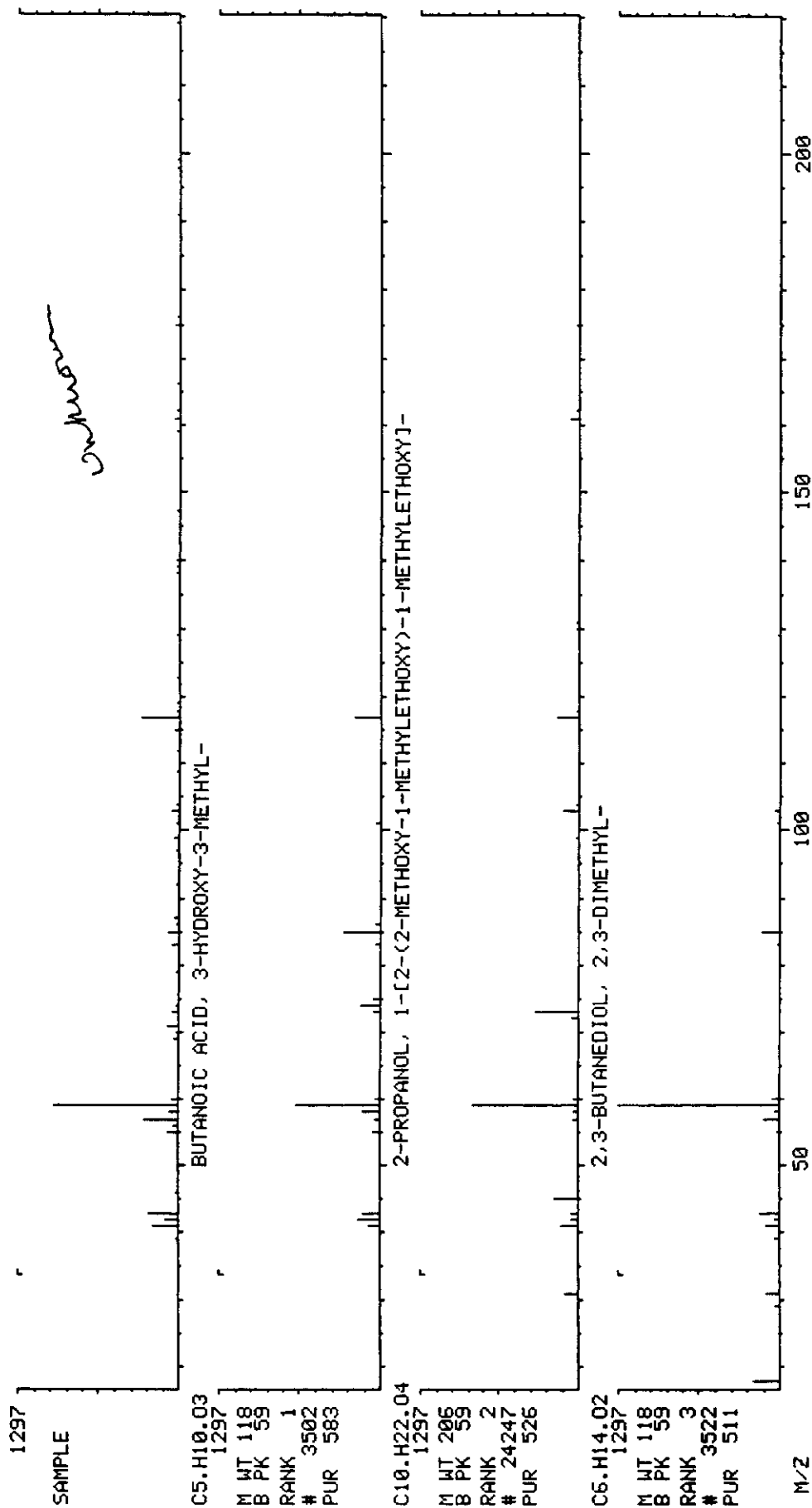
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C5.H10.O3	118	59	583	842	598
2	C10.H22.O4	206	59	526	741	581
3	C6.H14.O2	118	59	511	937	511
4	C8.H18.O	130	59	506	831	514
5	C9.H18.O3	174	59	479	692	513
6	C8.H18.O	130	59	471	931	471
7	C8.H20.S1	144	59	467	848	517
8	C9.H20.O4	192	59	458	714	506
9	C7.H14.O3	146	59	442	903	467

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	625-08-1
2	---	---	---	---	20324-33-8
3	---	---	---	---	76-09-5
4	---	---	---	---	19550-03-9
5	---	---	---	---	55956-25-7
6	---	---	---	---	23171-85-9
7	---	---	---	---	871-92-1
8	---	---	---	---	1638-16-0
9	---	---	---	---	54063-14-8

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:30:00 + 16:47
 SAMPLE: T-MM5-FB 1/35A/1ML
 COND.: UG/ML *1ML *100%/100% *(NA/NA >1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068104 #1509
 CALI: 30068104 # 3

BASE M/Z: 59
 RIC: 8624.



Library Search Data: 30068104 #1528 Base m/z: 59
 08/31/98 20:30:00 + 17:00 Cali: 30068104 # 3 RIC: 7552.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 904 matched at least 6 of the 16 largest peaks in the unknown

Rank In.	Name
1	5522 2-HEXANOL, 2,3-DIMETHYL-
2	5527 2-HEPTANOL, 2-METHYL-
3	16192 2-PROPANOL, 1-[1-METHYL-2-(2-PROPENYLOXY)ETHOXY]-
4	8506 4,5-DIMETHYL-3-HEPTANOL
5	9228 1-PROPANOL, 2-(2-METHOXYPROPOXY)-
6	26482 4-O-ACETYL-2,5-DI-O-METHYL-3,6-DIDEOXY-D-GLUCONITRILE
7	3528 2-METHYL-2,3-PENTANEDIOL
8	3502 BUTANOIC ACID, 3-HYDROXY-3-METHYL-
9	5497 3-ETHYL-2-METHYL-2-PENTANOL

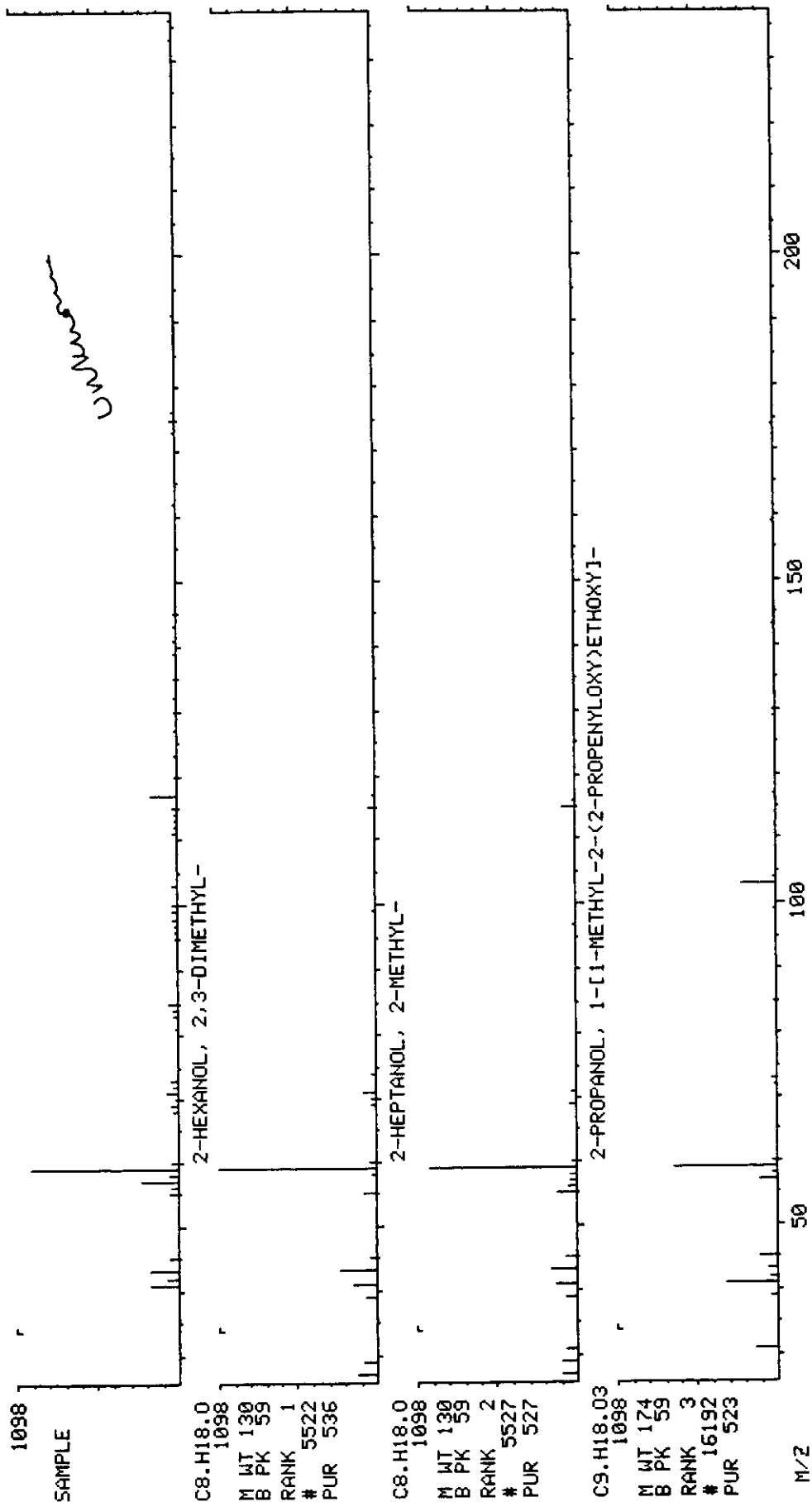
Rank	Formula	M.Wt	B.Pk	Purity	Fit	Rfit
1	C8.H18.O	130	59	536	873	562
2	C8.H18.O	130	59	527	903	544
3	C9.H18.O3	174	59	523	745	541
4	C9.H20.O	144	59	503	846	520
5	C7.H16.O3	148	59	499	790	521
6	C10.H17.O4.N	215	59	496	844	522
7	C6.H14.O2	118	59	496	896	524
8	C5.H10.O3	118	59	496	749	574
9	C8.H18.O	130	59	495	952	508

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	19550-03-9
2	—	—	—	—	625-25-2
3	—	—	—	—	55956-25-7
4	—	—	—	—	- -
5	—	—	—	—	13588-28-8
6	—	—	—	—	- -
7	—	—	—	—	7795-80-4
8	—	—	—	—	625-08-1
9	—	—	—	—	- -

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:30:00 + 17:00
 SAMPLE: T-MM5-FB 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *1ML *1002/1002 *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068104 #1528
 CALI: 30068104 # 3

BASE M/Z: 59
 RIC: 7552.



Library Search Data: 30068104 #1586 Base m/z: 55
 08/31/98 20:30:00 + 17:38 Cali: 30068104 # 3 RIC: 52032.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

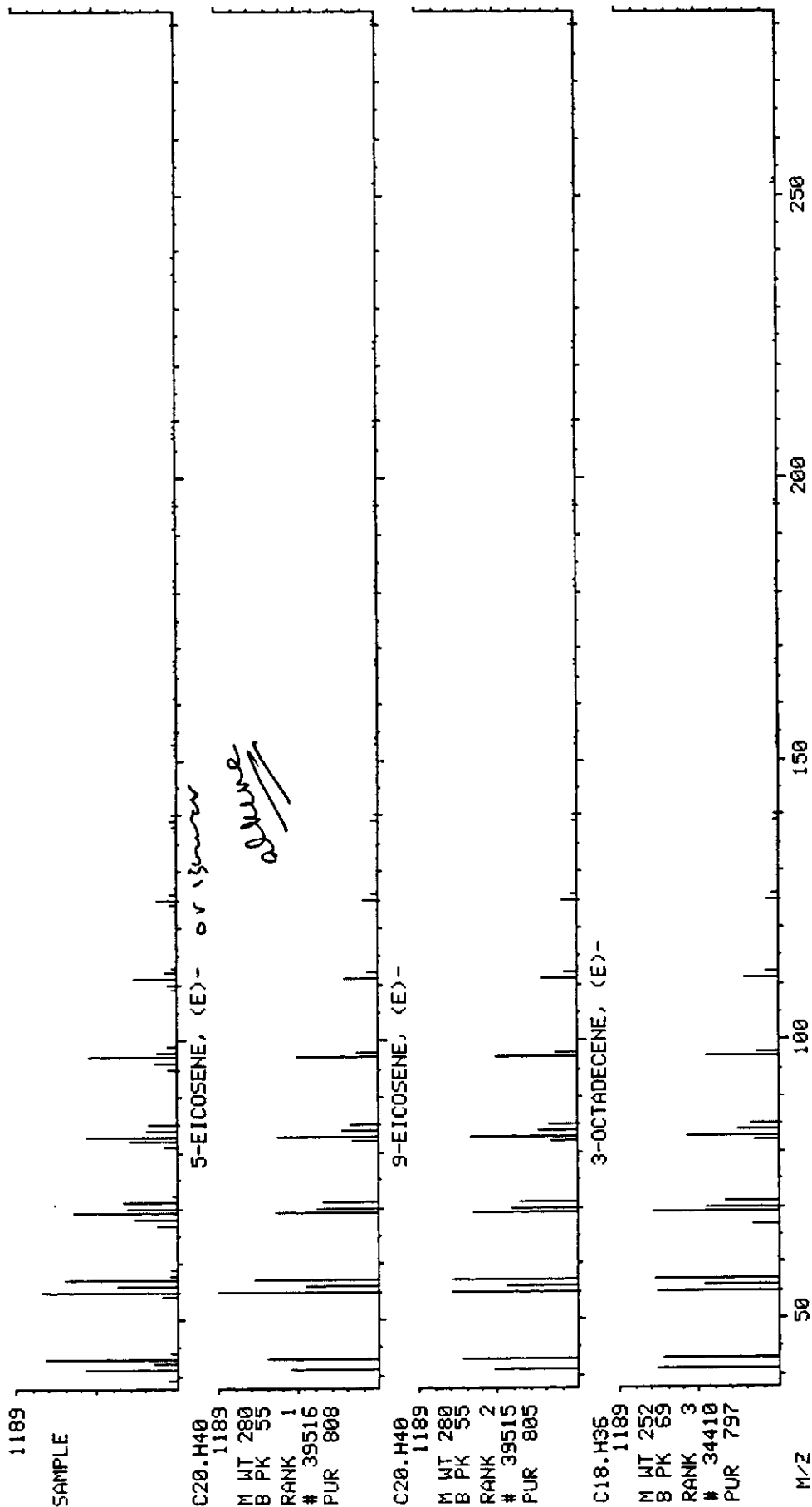
62231 spectra in LIBRARYNB searched for maximum PURITY
 950 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 39516 5-EICOSENE, (E)-
 2 39515 9-EICOSENE, (E)-
 3 34410 3-OCTADECENE, (E)-
 4 37062 1-NONADECENE
 5 34411 5-OCTADECENE, (E)-
 6 39517 3-EICOSENE, (E)-
 7 34418 9-OCTADECENE, (E)-
 8 26416 4-TETRADECANOL
 9 31653 1-HEPTADECENE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C20.H40	280	55	808	983	812
2	C20.H40	280	55	805	982	808
3	C18.H36	252	69	797	962	812
4	C19.H38	266	97	794	971	804
5	C18.H36	252	55	793	961	807
6	C20.H40	280	57	793	985	796
7	C18.H36	252	55	792	960	806
8	C14.H30.O	214	43	792	989	798
9	C17.H34	238	55	792	977	800

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	74685-30-6
2	---	---	---	---	74685-29-3
3	---	---	---	---	7206-19-1
4	---	---	---	---	18435-45-5
5	---	---	---	---	7206-21-5
6	---	---	---	---	74685-33-9
7	---	---	---	---	7206-25-9
8	---	---	---	---	1653-33-4
9	---	---	---	---	6765-39-5

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:30:00 + 17:38
 SAMPLE: T-MM5-FB 1/35A/1ML
 COND5.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068104 #1586
 CALI: 30068104 # 3
 BASE M/2: 55
 RIC: 52032.



Library Search Data: 30068104 #1661 Base m/z: 59
 08/31/98 20:30:00 + 18:29 Cali: 30068104 # 3 RIC: 7296.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 221 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 16192 2-PROPANOL, 1-[1-METHYL-2-(2-PROPENYLOXY)ETHOXY]-
 2 1909 2-PROPANOL, 1-METHOXY-2-METHYL-
 3 20626 2-PROPANOL, 1,1'-[(1-METHYL-1,2-ETHANEDIYL)BIS(OXY)]BIS-
 4 6083 2-PROPANOL, 1,1'-OXYBIS-
 5 3528 2-METHYL-2,3-PENTANEDIOL
 6 5809 1-ETHOXPENTAN-3-OL
 7 16651 DIETHYLENE GLYCOL TERT-BUTYL ETHER METHYL ETHER
 8 8460 SILANE, OCTYL-
 9 6082 1-PROPANOL, 2-(2-HYDROXYPROPOXY)-

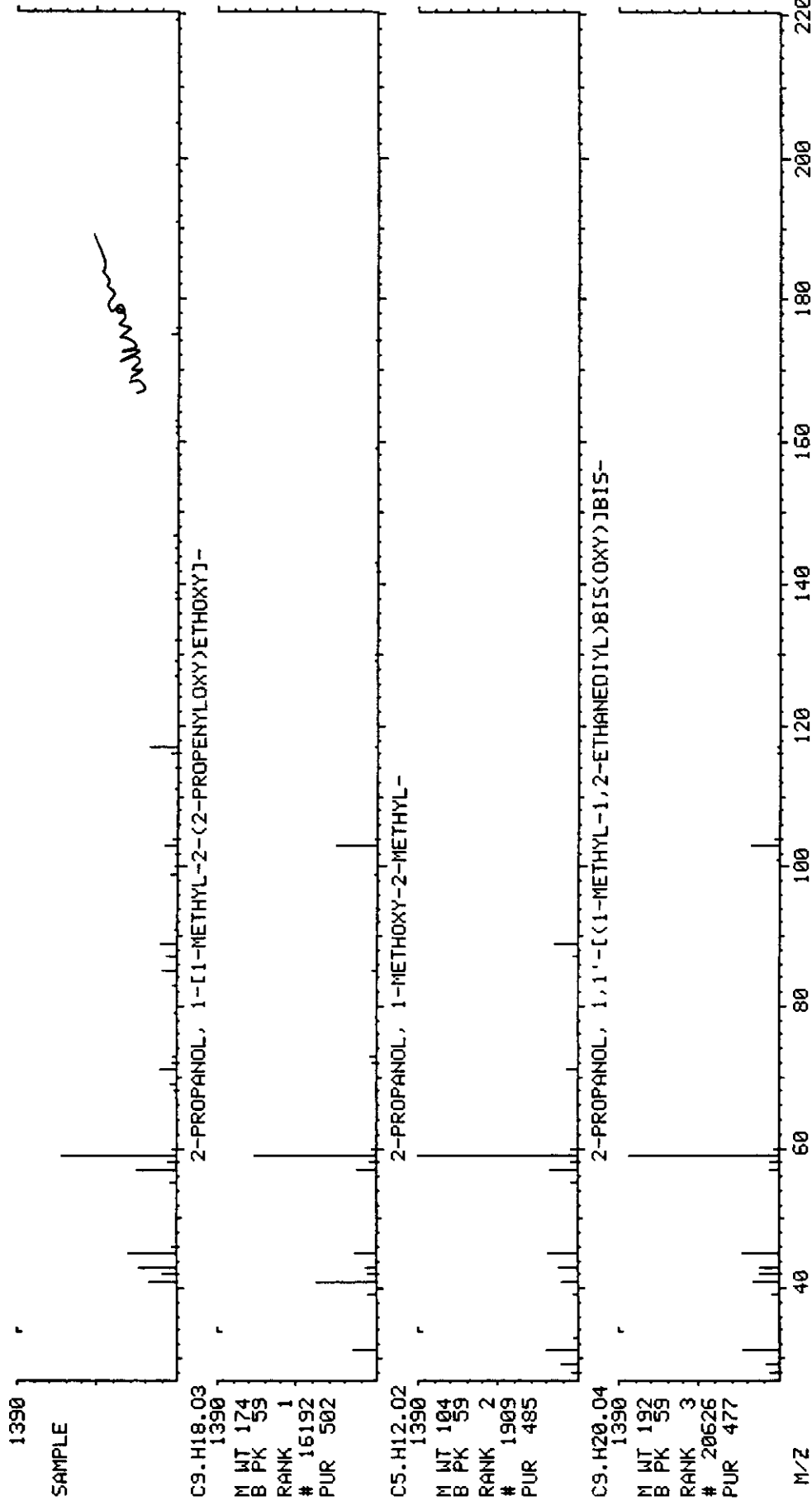
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C9.H18.O3	174	59	502	788	513
2	C5.H12.O2	104	59	485	945	490
3	C9.H20.O4	192	59	477	819	494
4	C6.H14.O3	134	59	452	889	464
5	C6.H14.O2	118	59	445	860	476
6	C7.H16.O2	132	59	440	868	480
7	C9.H20.O3	176	57	438	824	462
8	C8.H20.S1	144	59	430	863	455
9	C6.H14.O3	134	59	421	818	429

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	55956-25-7
2	—	—	—	—	3587-64-2
3	—	—	—	—	1638-16-0
4	—	—	—	—	110-98-5
5	—	—	—	—	7795-80-4
6	—	—	—	—	- -
7	—	—	—	—	52788-79-1
8	—	—	—	—	871-92-1
9	—	—	—	—	106-62-7

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:30:00 + 18:29
 SAMPLE: T-MM5-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (5 158 2N 0T)

DATA: 30068104 #1661
 CALI: 30068104 # 3

BASE M/Z: 59
 RIC: 7296.



Library Search Data: 30068104 #1686 Base m/z: 59
 08/31/98 20:30:00 + 18:45 Cali: 30068104 # 3 RIC: 3724.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 301 matched at least 4 of the 16 largest peaks in the unknown

Rank In. Name
 1 43635 TRIETHYL (S)-2-(CARBOXYMETHOXYMETHYL)OXYDIACETATE
 2 8432 2-HEXANONE, 3-HYDROXY-3,5-DIMETHYL-
 3 42663 2,4,6,8,9-PENTATHIATRICYCLO[3.3.1.1^{3,7}]DECANE-1-THIOL, 3,5,7-TRIMET*
 4 5826 2-PROPANOL, 1-ISOPROPOXY-2-METHYL-
 5 5815 1-TERT-BUTOXY-2-METHOXYETHANE
 6 51383 4-HEXADECEN-2-OL, 1-[(2,2-DIMETHYL-1,3-DIOXOLAN-4-YL)METHOXY]-
 7 60820 2,4,6,8,9-PENTATHIATRICYCLO[3.3.1.1^{3,7}]DECANE, 3,3'-DITHIOBIS[1,5,7*
 8 5778 BUTANOIC ACID, 3-HYDROXY-3-METHYL-, METHYL ESTER
 9 5916 BUTANOIC ACID, 4-NITRO-

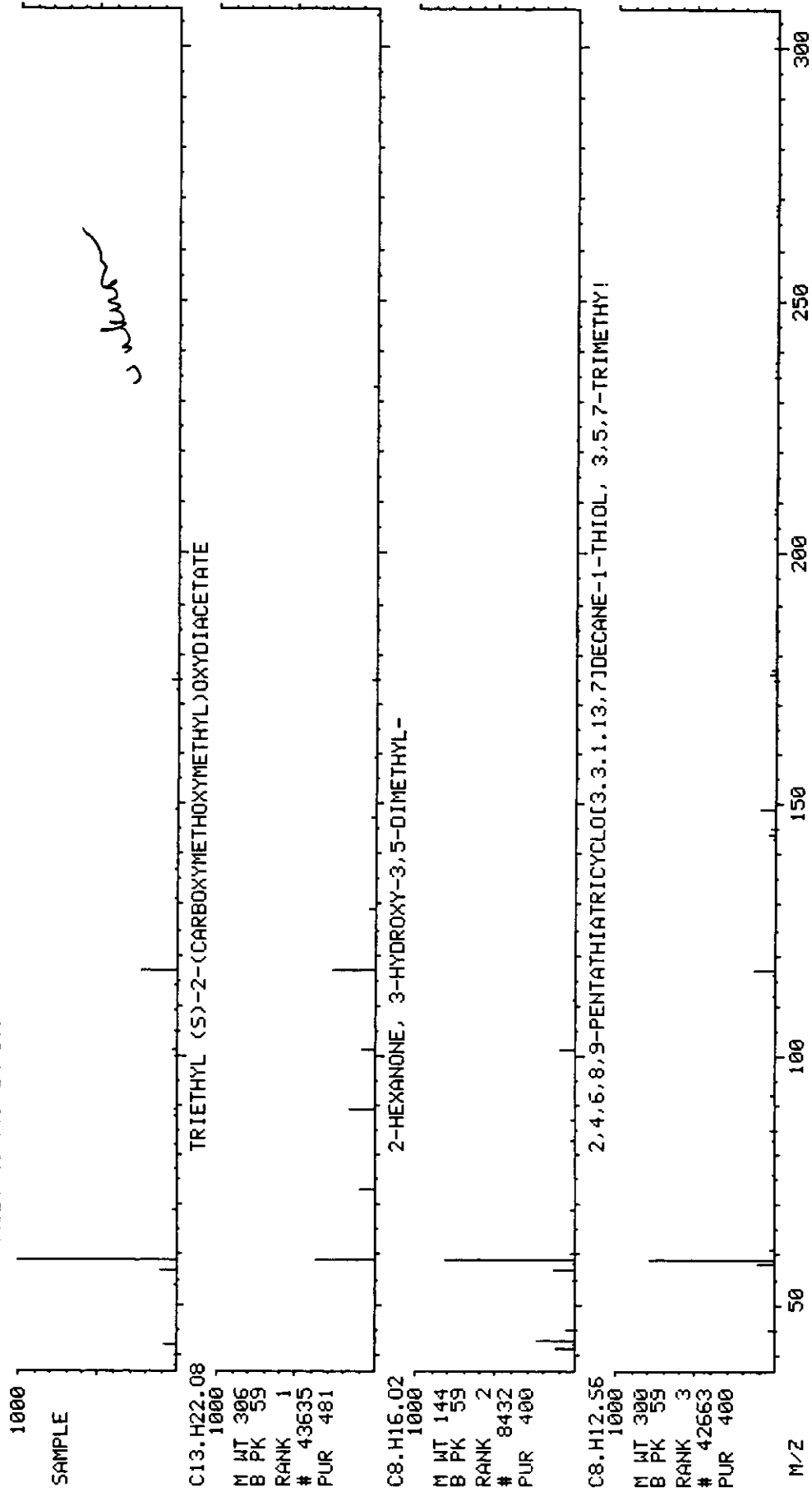
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H22.O8	306	59	481	621	627
2	C8.H16.O2	144	59	400	683	552
3	C8.H12.S6	300	59	400	517	690
4	C7.H16.O2	132	59	384	512	636
5	C7.H16.O2	132	57	372	480	524
6	C22.H42.O4	370	117	347	453	412
7	C16.H22.S12	598	59	330	400	735
8	C6.H12.O3	132	43	328	446	722
9	C4.H7.O4.N	133	59	309	404	586

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	- -
2	—	—	—	—	6321-14-8
3	—	—	—	—	57274-31-4
4	—	—	—	—	3587-75-5
5	—	—	—	—	66728-50-5
6	—	—	—	—	56256-43-0
7	—	—	—	—	57274-52-9
8	—	—	—	—	6149-45-7
9	—	—	—	—	16488-43-0

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:30:00 + 18:45
 SAMPLE: T-MM5-FB 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA) 1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068104 #1686
 CALI: 30068104 # 3

BASE M/Z: 59
 RIC: 3724.



Library Search Data: 30068104 #1704 Base m/z: 57
 08/31/98 20:30:00 + 18:57 Cali: 30068104 # 3 RIC: 7400.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 922 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 26409 2-HEXYL-1-OCTANOL
 2 32418 2-HEXYL-1-DECANOL
 3 19523 1-DECANOL, 2-ETHYL-
 4 32425 1-DECANOL, 2-HEXYL-
 5 34410 3-OCTADECENE, (E)-
 6 34418 9-OCTADECENE, (E)-
 7 34411 5-OCTADECENE, (E)-
 8 58380 DECANEDIOIC ACID, DIDECYL ESTER
 9 19520 1-OCTANOL, 2-BUTYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C14.H30.O	214	57	726	960✓	743
2	C16.H34.O	242	57	715	957✓	742
3	C12.H26.O	186	57	710	944	712
4	C16.H34.O	242	57	708	942	742
5	C18.H36	252	69	704	923	733
6	C18.H36	252	55	701	919	726
7	C18.H36	252	55	699	916	726
8	C30.H58.O4	482	57	696	960	707
9	C12.H26.O	186	57	695	945	718

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	—
2	—	—	—	—	—
3	—	—	—	—	21078-65-9
4	—	—	—	—	2425-77-6
5	—	—	—	—	7206-19-1
6	—	—	—	—	7206-25-9
7	—	—	—	—	7206-21-5
8	—	—	—	—	2432-89-5
9	—	—	—	—	3913-02-8

BASE M/Z: 57
RIC: 7400.

MID LIBRARY SEARCH (LIBRARYNB)

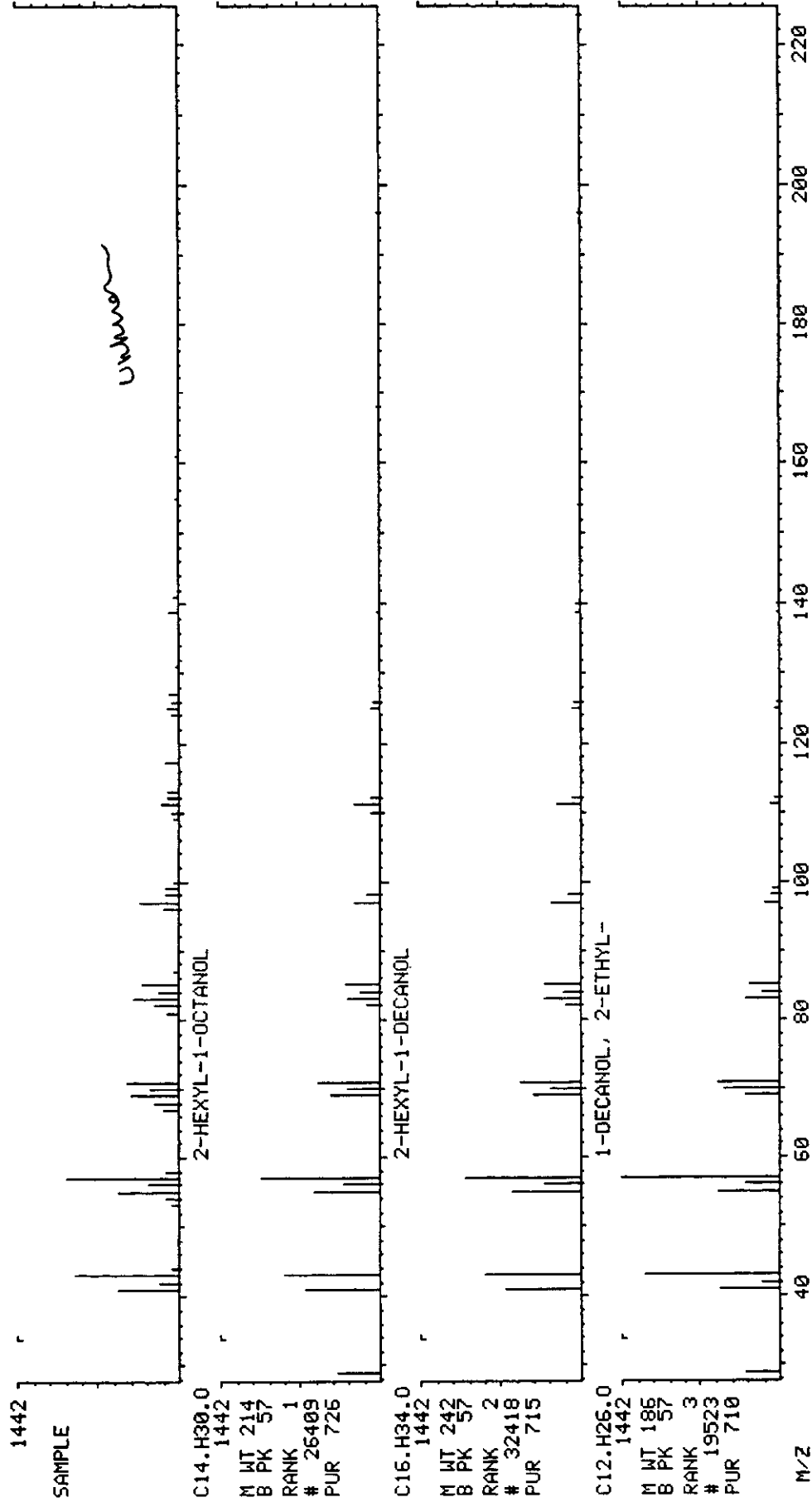
08/31/98 20:30:00 + 18:57

SAMPLE: T-MM5-FB 1/35A/1ML

COND.: UG/MIL *1ML *100%/100% *(NA/NA)/1/35A NA M

INST. ID:

CONDS: 06/11 *11L *10T
ENHANCED (S 15B 2N 0T)



935

Library Search Data: 30068104 #1741 Base m/z: 59
 08/31/98 20:30:00 + 19:22 Cali: 30068104 # 3 RIC: 17064.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 544 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 39622 9-OCTADECENAMIDE, (Z)-
 2 22656 DODECANAMIDE
 3 15482 HEPTANAMIDE, 4-ETHYL-5-METHYL-
 4 11721 NONANAMIDE
 5 3133 PENTANAMIDE, 4-METHYL-
 6 19516 1-DECANOL, 5,9-DIMETHYL-
 7 16267 1,7-OCTANEDIOL, 3,7-DIMETHYL-
 8 7998 TRANS-3,4-EPOXYNONANE
 9 11584 7-OCTEN-2-OL, 2,6-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C18.H35.O.N	281	59	568	866	647
2	C12.H25.O.N	199	59	484	883	523
3	C10.H21.O.N	171	59	472	867	488
4	C9.H19.O.N	157	59	472	867	487
5	C6.H13.O.N	115	59	444	892	456
6	C12.H26.O	186	41	436	747	488
7	C10.H22.O2	174	59	433	887	473
8	C9.H18.O	142	41	419	748	480
9	C10.H20.O	156	59	412	921	433

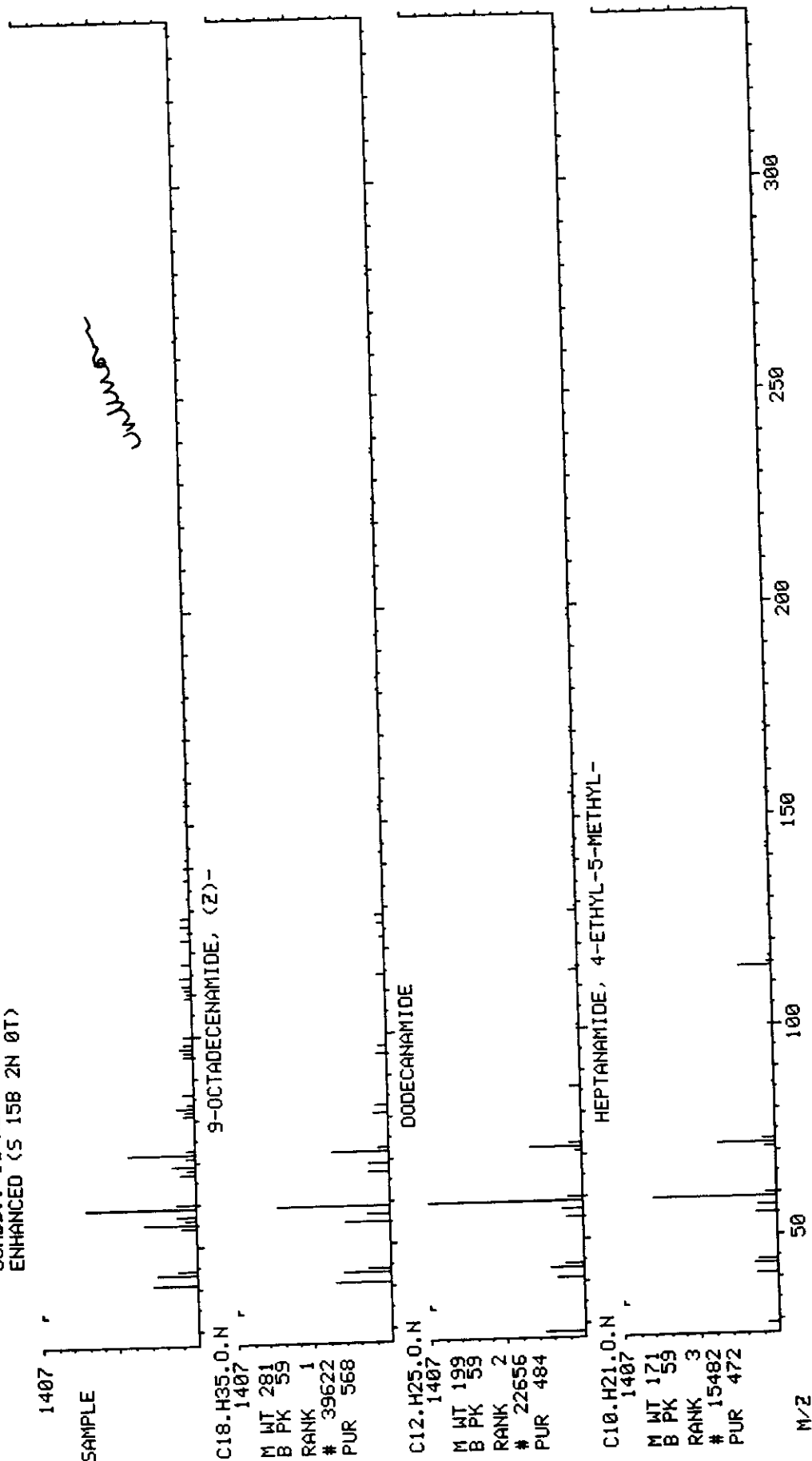
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	301-02-0
2	—	—	—	—	1120-16-7
3	—	—	—	—	54789-40-1
4	—	—	—	—	1120-07-6
5	—	—	—	—	1119-29-5
6	—	—	—	—	91482-38-1
7	—	—	—	—	107-74-4
8	—	—	—	—	—
9	—	—	—	—	18479-58-8

BASE M/Z: 59
RIC: 7064.

DATA: 30068104 #1741
CALI: 30068104 # 3

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 20:30:00 + 19:22
SAMPLE: T-MM5-FB 1/35A/1ML INST. ID: F16
CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)



Library Search Data: 30068104 #1799 Base m/z: 45
 08/31/98 20:30:00 + 20:01 Cali: 30068104 # 3 RIC: 4152.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 996 matched at least 4 of the 16 largest peaks in the unknown

Rank In. Name
 1 9227 1,3,3-TRIMETHOXYBUTANE
 2 41700 1,2,3,5,6-PENTA-O-METHYL-4-O-ACETYL-GLUCITOL
 3 27502 15-CROWN-5
 4 41712 3,6,9,12,15-PENTAOXANONADECAN-1-OL
 5 21103 ETHANOL, 2,2'-[OXYBIS(2,1-ETHANEDIYLOXY)]BIS-
 6 43636 .BETA.-D-MANNOPYRANOSE, 2,4,6-TRI-O-METHYL-, DIACETATE
 7 33841 3,6,9,12-TETRAOXAHEXADECAN-1-OL
 8 36510 1,4,7,10,13,16-HEXAOXACYCLOOCTADECANE
 9 34224 2,5,8,11,14-PENTAOXAHEXADECAN-16-OL

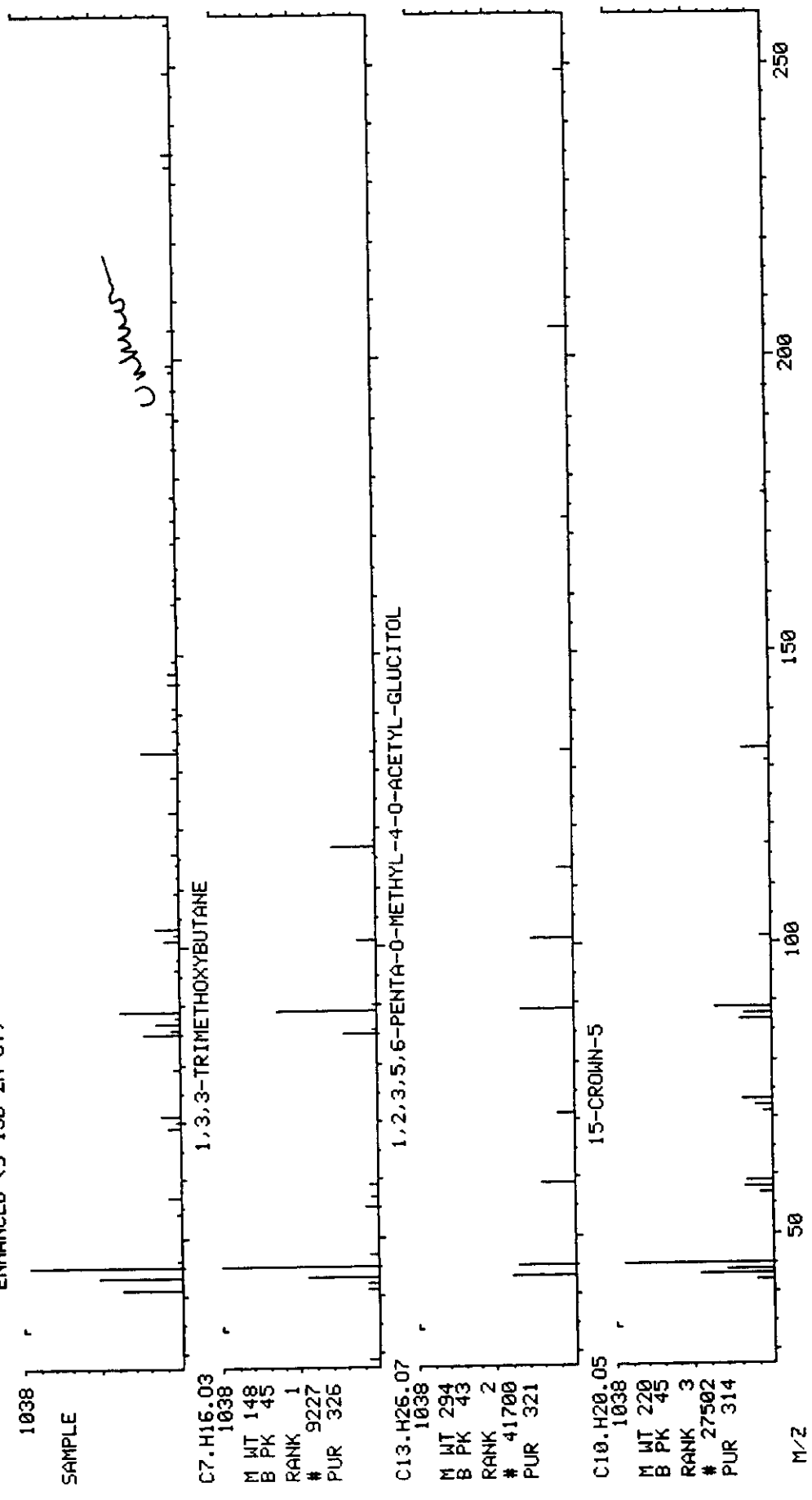
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C7.H16.O3	148	45	326	678	424
2	C13.H26.O7	294	43	321	748	370
3	C10.H20.O5	220	45	314	686	435
4	C14.H30.O6	294	45	304	678	398
5	C8.H18.O5	194	45	294	718	343
6	C13.H22.O8	306	45	293	610	353
7	C12.H26.O5	250	45	292	588	414
8	C12.H24.O6	264	45	281	660	404
9	C11.H24.O6	252	45	277	580	433

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	6607-66-5
2	—	—	—	—	- -
3	—	—	—	—	33100-27-5
4	—	—	—	—	1786-94-3
5	—	—	—	—	112-60-7
6	—	—	—	—	55255-81-7
7	—	—	—	—	1559-34-8
8	—	—	—	—	17455-13-9
9	—	—	—	—	23778-52-1

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:30:00 + 20:01
 SAMPLE: T-MM5-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068104 #1799
 CALI: 30068104 # 3

BASE M/Z: 45
 RIC: 4152.



Library Search Data: 30068104 #1804 Base m/z: 59
08/31/98 20:30:00 + 20:04 Cali: 30068104 # 3 RIC: 6872.
Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
221 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
1 9228 1-PROPANOL, 2-(2-METHOXYPROPOXY)-
2 5826 2-PROPANOL, 1-ISOPROPOXY-2-METHYL-
3 1909 2-PROPANOL, 1-METHOXY-2-METHYL-
4 5786 PROPANOIC ACID, 2-HYDROXY-2-METHYL-, ETHYL ESTER
5 1905 2-BUTANOL, 3-METHOXY-
6 1919 1-BUTANOL, 3-METHOXY-
7 3502 BUTANOIC ACID, 3-HYDROXY-3-METHYL-
8 3528 2-METHYL-2,3-PENTANEDIOL
9 36510 1,4,7,10,13,16-HEXAOXACYCLOOCTADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C7.H16.O3	148	59	474	865	479
2	C7.H16.O2	132	59	419	676	492
3	C5.H12.O2	104	59	419	930	441
4	C6.H12.O3	132	59	412	760	422
5	C5.H12.O2	104	59	404	860	418
6	C5.H12.O2	104	59	403	877	429
7	C5.H10.O3	118	59	402	801	439
8	C6.H14.O2	118	59	401	860	430
9	C12.H24.O6	264	45	398	608	450

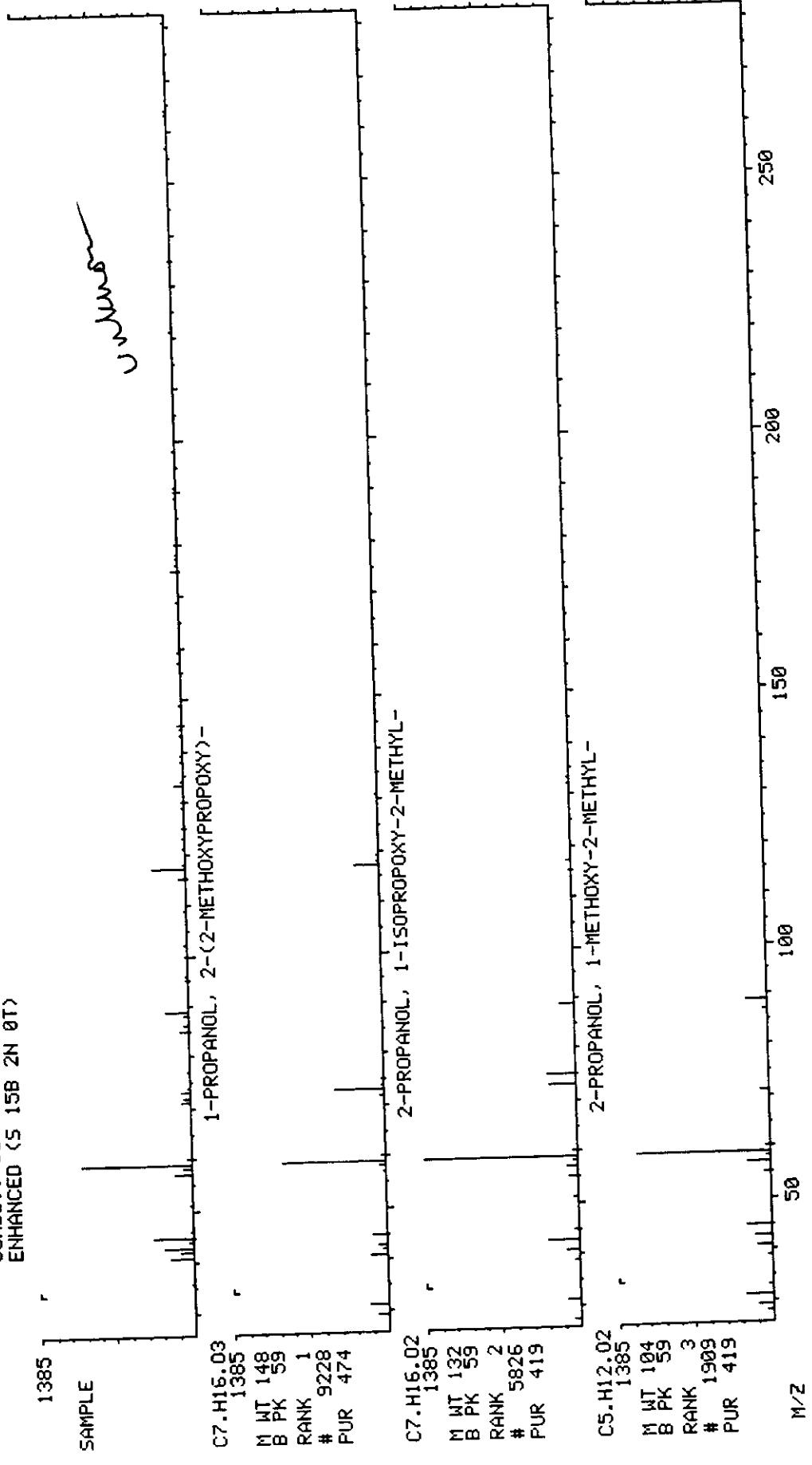
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	13588-28-8
2	—	—	—	—	3587-75-5
3	—	—	—	—	3587-64-2
4	—	—	—	—	80-55-7
5	—	—	—	—	53778-72-6
6	—	—	—	—	2517-43-3
7	—	—	—	—	625-08-1
8	—	—	—	—	7795-80-4
9	—	—	—	—	17455-13-9

BASE M/Z: 59
RIC: 6872.

DATA: 30058104 #1804
CALI: 30068104 # 3

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 20:30:00 + 20:04
SAMPLE: T-MM5-FB 1/35A/1ML
COND5.: UG/ML *1ML *100%/100% *(NA/NA) >1/35A NA M
ENHANCED (5 15B 2N 0T)

INST. ID: F16
NA M



Library Search Data: 30068104 #1914 Base m/z: 43
 08/31/98 20:30:00 + 21:17 Cali: 30068104 # 3 RIC: 22720.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 818 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 26416 4-TETRADECANOL
 2 37062 1-NONADECENE
 3 39515 9-EICOSENE, (E)-
 4 39516 5-EICOSENE, (E)-
 5 31653 1-HEPTADECENE
 6 34410 3-OCTADECENE, (E)-
 7 34418 9-OCTADECENE, (E)-
 8 58701 17-PENTATRIACONTENE
 9 34411 5-OCTADECENE, (E)-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C14.H30.O	214	43	754	982	758
2	C19.H38	266	97	754	963	768
3	C20.H40	280	57	752	958	764
4	C20.H40	280	55	750	956	764
5	C17.H34	238	55	750	961	754
6	C18.H36	252	69	747	945	761
7	C18.H36	252	55	745	942	758
8	C35.H70	490	43	745	935	785
9	C18.H36	252	55	745	942	758

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	1653-33-4
2	---	---	---	---	18435-45-5
3	---	---	---	---	74685-29-3
4	---	---	---	---	74685-30-6
5	---	---	---	---	6765-39-5
6	---	---	---	---	7206-19-1
7	---	---	---	---	7206-25-9
8	---	---	---	---	6971-40-0
9	---	---	---	---	7206-21-5

BASE M/Z: 43
RIC: 22720.

MTD LIBRARY SEARCH (LIBRARYNB)

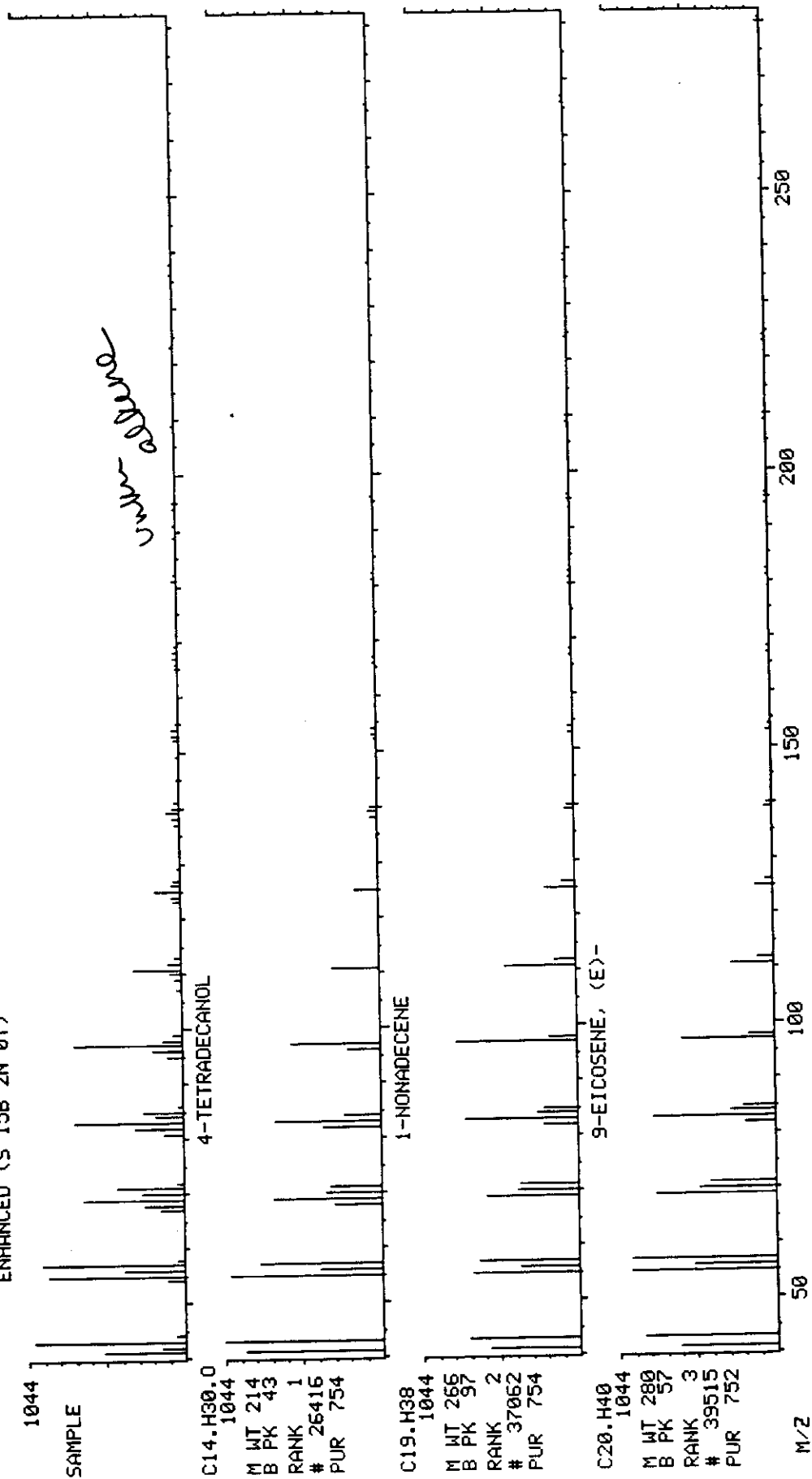
08/31/98 20:30:00 + 21:17

SAMPLE: T-MM5-FB 1/35A/1ML

08/31/20 20:33:33	INST. ID:	F16
SAMPLE: T-MM5-FB	1/35A/1ML	NA M
CONDS.: 116/1	*1ML	
	*100%/100%	
	*(NA/NA)/1/35A	

INST. ID:

CONDUITS: 00/11C *11C *11C
ENHANCED (5 15B 2N 0T)



Library Search Data: 30068104 #1932 Base m/z: 59
 08/31/98 20:30:00 + 21:29 Cali: 30068104 # 3 RIC: 7480.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 816 matched at least 5 of the 16 largest peaks in the unknown

Rank In. Name
 1 1909 2-PROPANOL, 1-METHOXY-2-METHYL-
 2 3528 2-METHYL-2,3-PENTANEDIOL
 3 27502 15-CROWN-5
 4 5825 2-PROPANOL, 1-(2-METHYLPROPOXY)-
 5 5809 1-ETHOXYPENTAN-3-OL
 6 41712 3,6,9,12,15-PENTAOXANONADECAN-1-OL
 7 1760 2-PENTANOL, 2-METHYL-
 8 8813 ETHENE, (2-ETHOXY-1-METHOXYETHOXY)-
 9 34224 2,5,8,11,14-PENTAOXAHEXADECAN-16-OL

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C5.H12.O2	104	59	435	910	447
2	C6.H14.O2	118	59	418	848	435
3	C10.H20.O5	220	45	414	719	462
4	C7.H16.O2	132	57	400	777	441
5	C7.H16.O2	132	59	400	892	425
6	C14.H30.O6	294	45	393	727	445
7	C6.H14.O	102	59	391	858	397
8	C7.H14.O3	146	59	390	826	401
9	C11.H24.O6	252	45	388	840	424

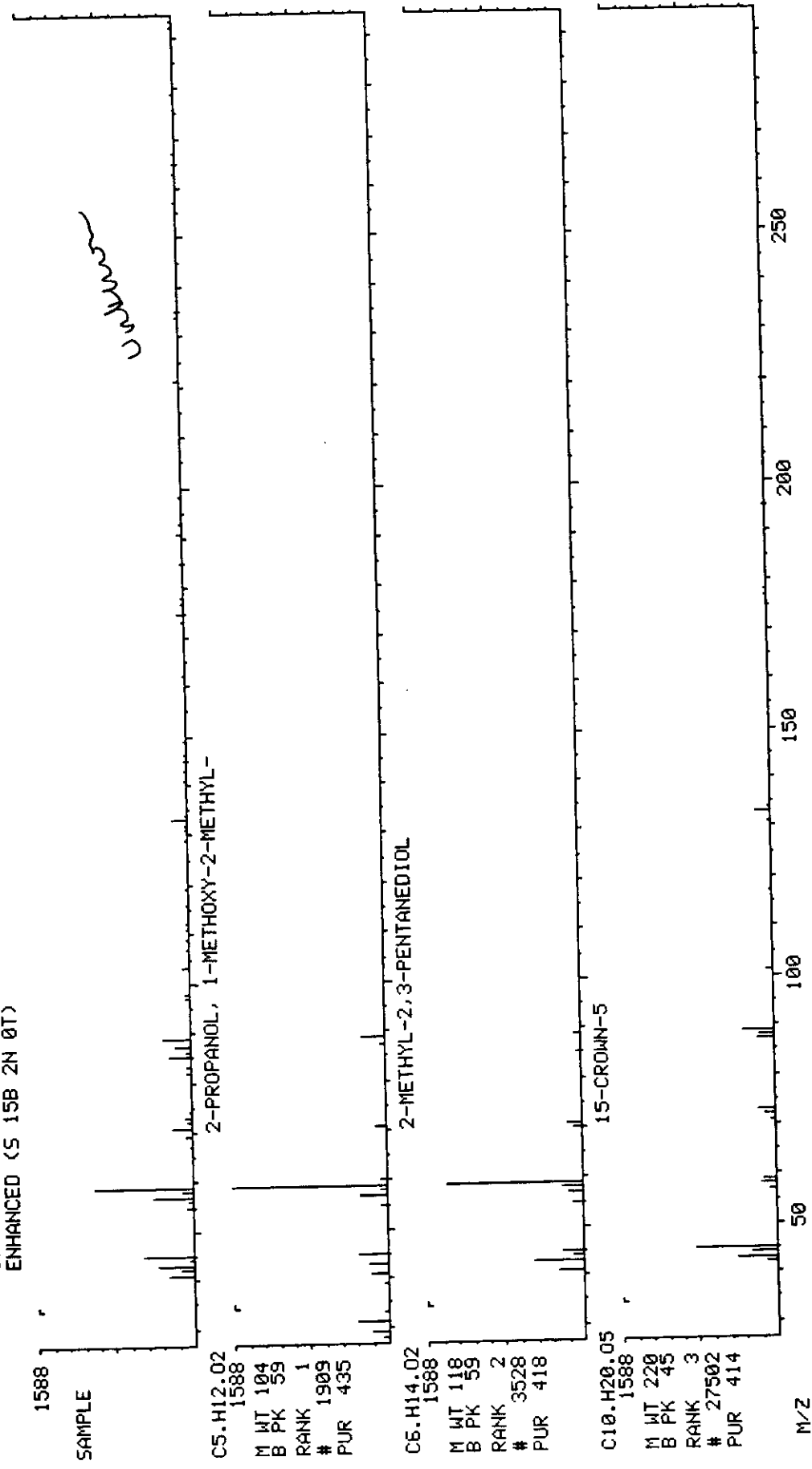
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	3587-64-2
2	—	—	—	—	7795-80-4
3	—	—	—	—	33100-27-5
4	—	—	—	—	23436-19-3
5	—	—	—	—	- -
6	—	—	—	—	1786-94-3
7	—	—	—	—	590-36-3
8	—	—	—	—	54063-18-2
9	—	—	—	—	23778-52-1

BASE M/Z: 59
RIC: 7480.

DATA: 30068104 #1932
CALI: 30068104 # 3

MID LIBRARY SEARCH <LIBRARYNB>
08/31/98 20:30:00 + 21:29
SAMPLE: T-MM5-FB 1/35A/1ML
CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



CS
CI

Library Search Data: 30068104 #2012 Base m/z: 57
 08/31/98 20:30:00 + 22:23 Cali: 30068104 # 3 RIC: 11760.
 Sample: T-MM5-F8 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 818 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 39515 9-EICOSENE, (E)-
 2 39516 5-EICOSENE, (E)-
 3 26416 4-TETRADECANOL
 4 34410 3-OCTADECENE, (E)-
 5 34418 9-OCTADECENE, (E)-
 6 37062 1-NONADECENE
 7 37049 2-OCTADECENAL
 8 34411 5-OCTADECENE, (E)-
 9 31653 1-HEPTADECENE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C20.H40	280	55	692	945	708
2	C20.H40	280	55	691	943	708
3	C14.H30.O	214	43	691	973	701
4	C18.H36	252	69	690	941	707
5	C18.H36	252	55	689	939	704
6	C19.H38	266	97	687	943	710
7	C18.H34.O	266	43	687	909	731
8	C18.H36	252	55	687	937	704
9	C17.H34	238	55	685	945	700

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	74685-29-3
2	—	—	—	—	74685-30-6
3	—	—	—	—	1653-33-4
4	—	—	—	—	7206-19-1
5	—	—	—	—	7206-25-9
6	—	—	—	—	18435-45-5
7	—	—	—	—	56554-96-2
8	—	—	—	—	7206-21-5
9	—	—	—	—	6765-39-5

BASE M/Z: 57
RIC: 11760.

DATA: 30068104 #2012
CALI: 30068104 # 3

MID LIBRARY SEARCH (LIBRARYNB)

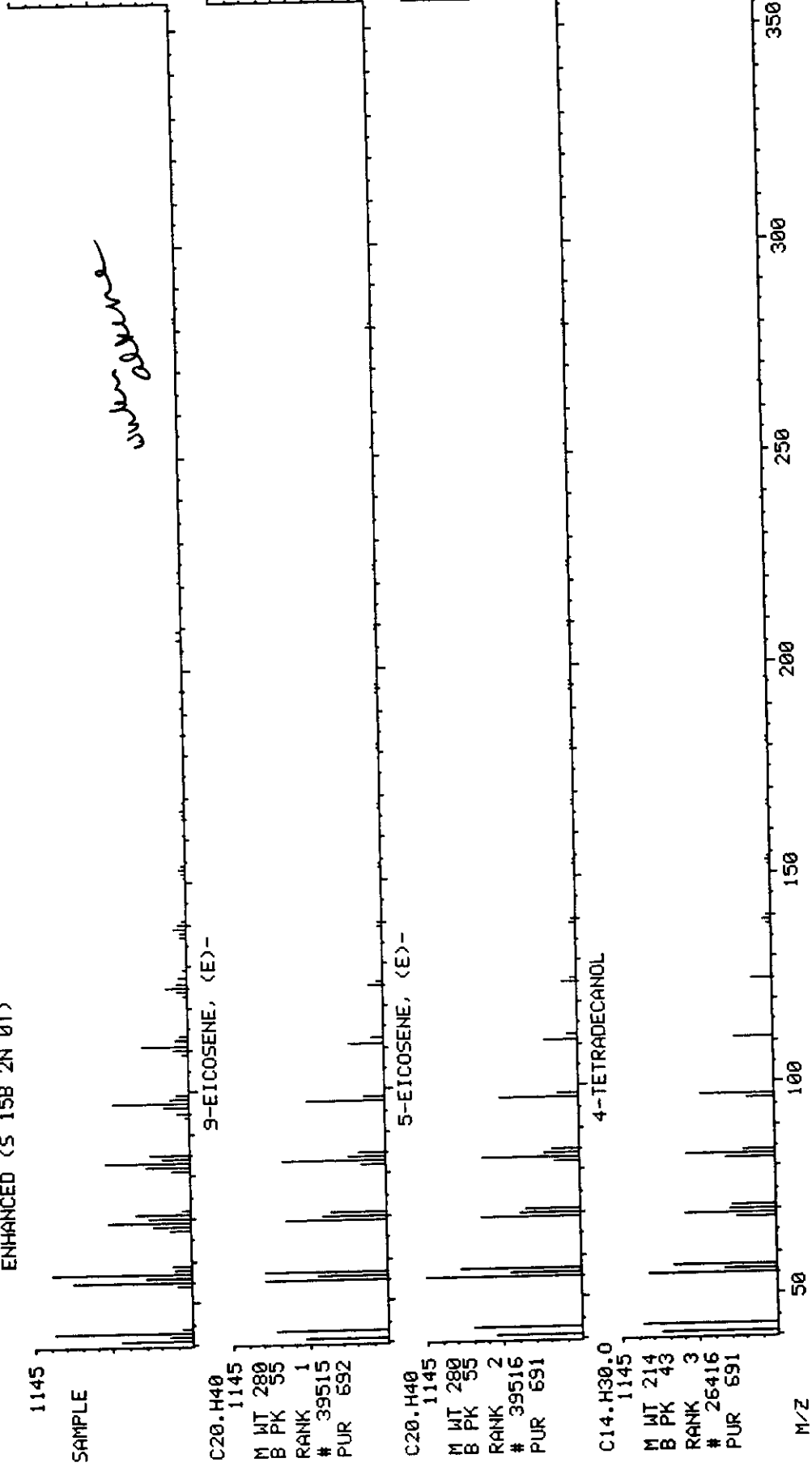
08/31/98 20:30:00 + 22:23

SAMPLE: T-MMS-FB 1/35A/1ML

CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA) 1/35A NA M

ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068104 #2049 Base m/z: 59
 08/31/98 20:30:00 + 22:47 Cali: 30068104 # 3 RIC: 4808.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 932 matched at least 5 of the 16 largest peaks in the unknown

Rank In. Name
 1 12385 BUTYRIC ACID, 4-ISOPROPOXY-, METHYL ESTER
 2 16199 BUTYRIC ACID, 4-BUTOXY-, METHYL ESTER
 3 8822 BUTANOIC ACID, 4-ETHOXY-, METHYL ESTER
 4 9227 1,3,3-TRIMETHOXYBUTANE
 5 1909 2-PROPANOL, 1-METHOXY-2-METHYL-
 6 8425 2-HEPTANONE, 3-HYDROXY-3-METHYL-
 7 5825 2-PROPANOL, 1-(2-METHYLPROPOXY)-
 8 27502 15-CROWN-5
 9 1760 2-PENTANOL, 2-METHYL-

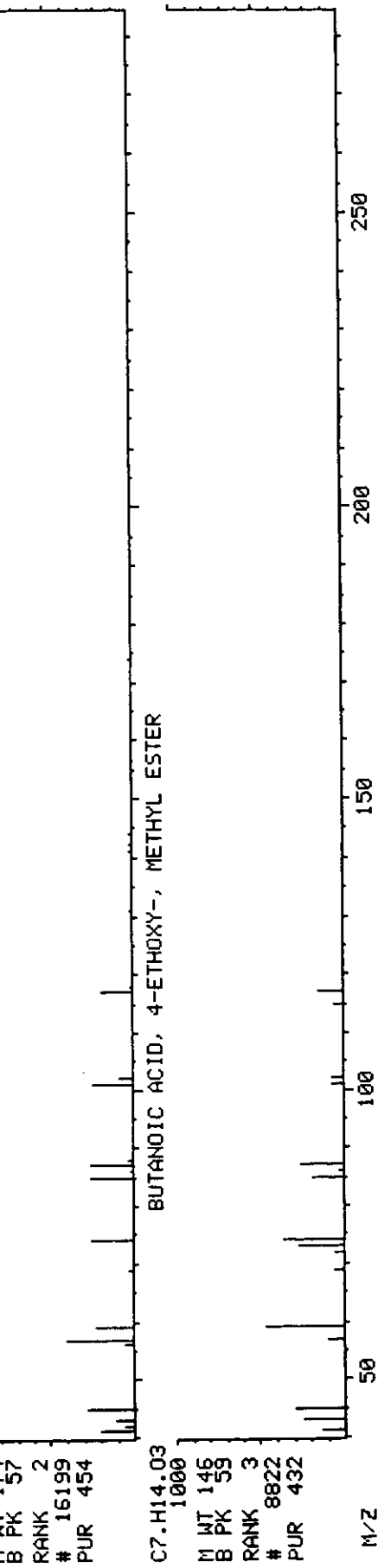
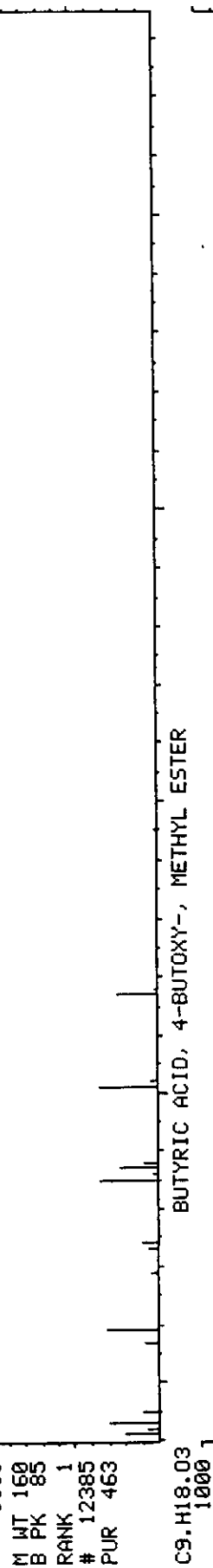
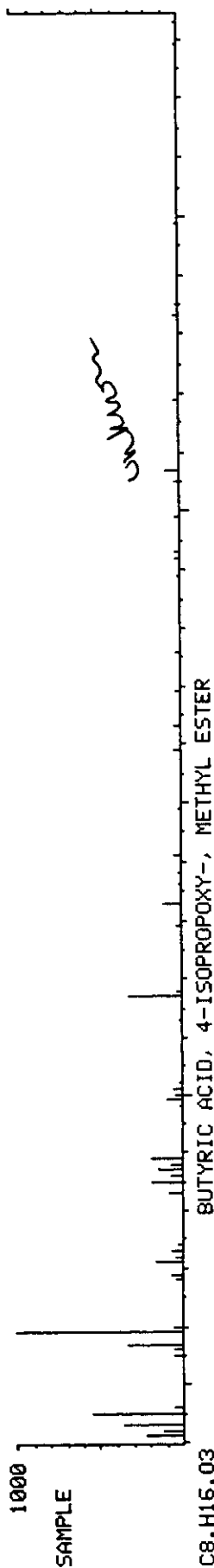
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C8.H16.O3	160	85	463	814	464
2	C9.H18.O3	174	57	454	787	464
3	C7.H14.O3	146	59	432	802	448
4	C7.H16.O3	148	45	408	712	420
5	C5.H12.O2	104	59	388	913	399
6	C8.H16.O2	144	59	382	793	385
7	C7.H16.O2	132	57	379	804	408
8	C10.H20.O5	220	45	369	643	429
9	C6.H14.O	102	59	367	877	372

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	29006-05-1
2	—	—	—	—	29006-06-2
3	—	—	—	—	29006-04-0
4	—	—	—	—	6607-66-5
5	—	—	—	—	3587-64-2
6	—	—	—	—	13757-91-0
7	—	—	—	—	23436-19-3
8	—	—	—	—	33100-27-5
9	—	—	—	—	590-36-3

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 20:30:00 + 22:47
 SAMPLE: T-MM5-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068104 #2049
 CALI: 30068104 # 3

BASE M/Z: 59
 RIC: 4808.



Library Search Data: 30068104 #2056 Base m/z: 59
 08/31/98 20:30:00 + 22:52 Cali: 30068104 # 3 RIC: 5416.
 Sample: T-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 338 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 27502 15-CROWN-5
 2 41712 3,6,9,12,15-PENTAOXANONADECAN-1-OL
 3 1909 2-PROPANOL, 1-METHOXY-2-METHYL-
 4 9227 1,3,3-TRIMETHOXYBUTANE
 5 36510 1,4,7,10,13,16-HEXAOXACYCLOOCTADECANE
 6 34224 2,5,8,11,14-PENTAOXAHEXADECAN-16-OL
 7 12385 BUTYRIC ACID, 4-ISOPROPOXY-, METHYL ESTER
 8 5825 2-PROPANOL, 1-(2-METHYLPROPOXY)-
 9 3528 2-METHYL-2,3-PENTANEDIOL

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C10.H20.O5	220	45	429	717	481
2	C14.H30.O6	294	45	416	727	470
3	C5.H12.O2	104	59	406	906	410
4	C7.H16.O3	148	45	401	715	414
5	C12.H24.O6	264	45	396	725	441
6	C11.H24.O6	252	45	378	834	407
7	C8.H16.O3	160	85	377	751	416
8	C7.H16.O2	132	57	373	777	418
9	C6.H14.O2	118	59	373	812	383

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	33100-27-5
2	---	---	---	---	1786-94-3
3	---	---	---	---	3587-64-2
4	---	---	---	---	6607-66-5
5	---	---	---	---	17455-13-9
6	---	---	---	---	23778-52-1
7	---	---	---	---	29006-05-1
8	---	---	---	---	23436-19-3
9	---	---	---	---	7795-80-4

DATA: 30068104 #2056
CALI: 30068104 # 3

BASE M/Z: 59
RIC: 5416.

MID LIBRARY SEARCH (LIBRARYNB)

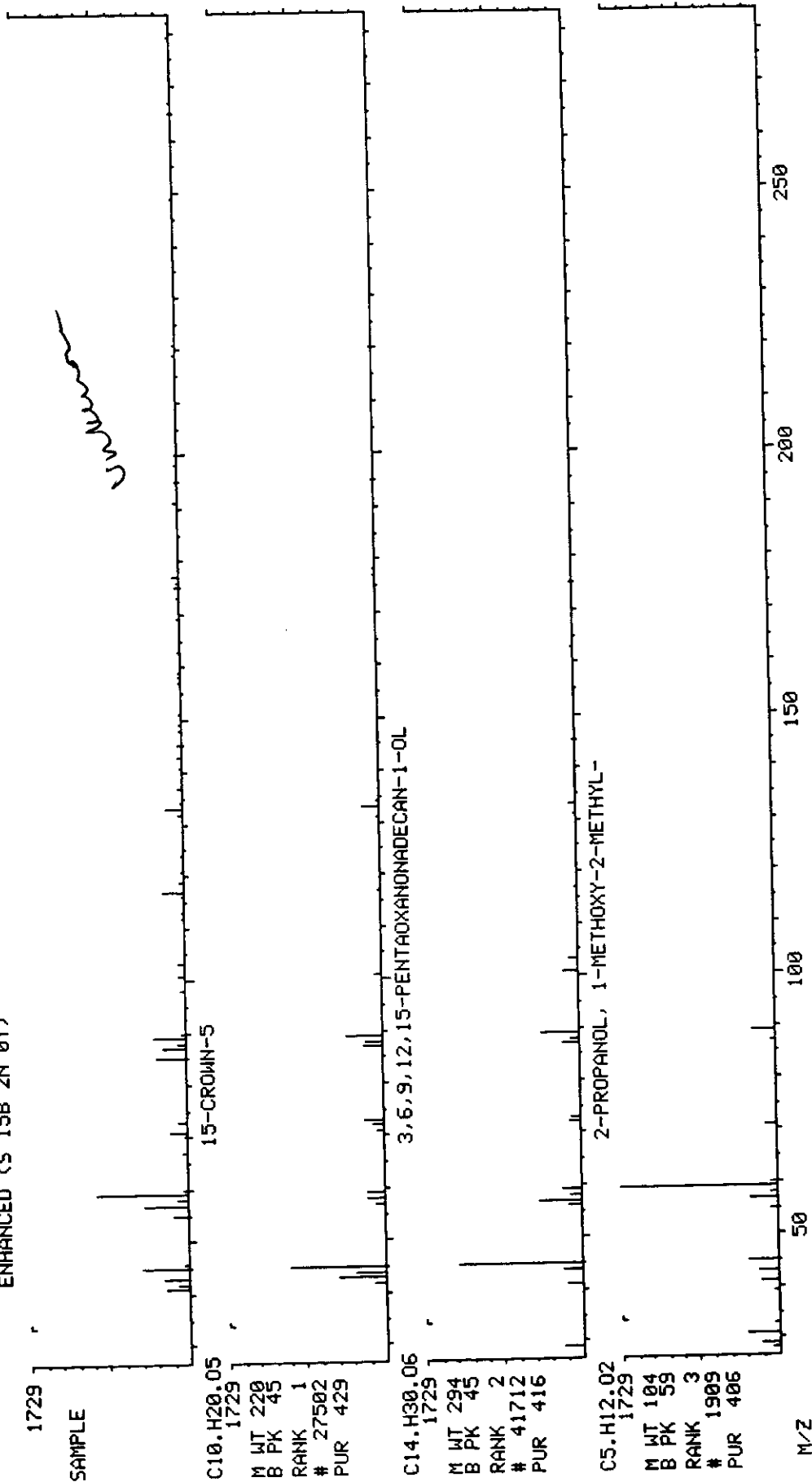
08/31/98 20:30:00 + 22:52

SAMPLE: T-MM5-FB 1/35A/1ML

INST. ID: F16

CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M

ENHANCED (S 15B 2N 0T)



TIC SELECTION REPORT

DATA FILE: 30068104

THE FOLLOWING PEAKS WERE REJECTED BECAUSE
AT LEAST 40 % OF THEIR SIZE WAS ACCOUNTED FOR BY
TARGET COMPOUNDS ELUTING WITHIN 4 SCANS OF THE
PEAK TOP.

SCAN	SIZE	AMOUNT
199	232884.	28.865
308	294104.	36.453
341	217648.	26.977
421	128124.	15.881
529	322720.	40.000
722	288940.	36.262
828	318724.	40.000
968	253508.	31.756
1086	319320.	40.000
1383	309316.	54.242
1383	311820.	54.681
1551	228100.	40.000
1782	194188.	40.000

TOTAL NUMBER OF UNIDENTIFIED PEAKS WITH SIZE
GREATER THAN 10 % OF THE CLOSEST INTERNAL STANDARD
THAT DOES NOT HAVE INTERFERENCES = 23

INTERNAL STANDARDS THAT HAVE RIC SIZE LESS THAN
50 % OR GREATER THAN 200 % OF THE ESTIMATED RIC SIZE
ARE CONSIDERED TO HAVE INTERFERENCES AND WILL NOT BE USED
FOR QUANTITATION.

#	INTERNAL STANDARD	PERCENT OF ESTIMATED		SCAN
		RIC SIZE	RIC SIZE	
1	C140 NAPHTHALENE-D8	322720.	156	529
2	C150 ACENAPHTHENE-D10	318724.	130	828
3	C160 PHENANTHRENE-D10	319320.	121	1086
4	C170 CHRYSENE-D12	228100.	130	1551
5	C175 PERYLENE-D12	194188.	123	1782

* INDICATES INTERFERENCE

SIZE = AREA

Semivolatile Organics
Method 0010/8270

Client Name: Pacific Environmental Services
Client ID: T-MM5-4-F, FH, XAD, COND, BH
LAB ID: 300681-0005-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 26 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
Phenol	ND	ug/Sample	30	R
bis(2-Chloroethyl)ether	ND	ug/Sample	30	
2-Chlorophenol	ND	ug/Sample	30	
1,3-Dichlorobenzene	ND	ug/Sample	30	
1,4-Dichlorobenzene	ND	ug/Sample	30	
Benzyl alcohol	ND	ug/Sample	30	
1,2-Dichlorobenzene	ND	ug/Sample	30	
2-Methylphenol	ND	ug/Sample	30	
2,2'-Oxybis(1-chloropropane)	ND	ug/Sample	30	
3/4-Methylphenol	ND	ug/Sample	30	
N-Nitroso-di-n-propylamine	ND	ug/Sample	30	
Hexachloroethane	ND	ug/Sample	30	
Nitrobenzene	ND	ug/Sample	30	
Isophorone	ND	ug/Sample	30	
2-Nitrophenol	ND	ug/Sample	30	
2,4-Dimethylphenol	ND	ug/Sample	30	
Benzoic acid	ND	ug/Sample	150	
bis(2-Chloroethoxy)-methane	ND	ug/Sample	30	
2,4-Dichlorophenol	ND	ug/Sample	30	
1,2,4-Trichlorobenzene	ND	ug/Sample	30	
Naphthalene	ND	ug/Sample	30	
4-Chloroaniline	ND	ug/Sample	30	
Hexachlorobutadiene	ND	ug/Sample	30	
4-Chloro-3-methylphenol	ND	ug/Sample	30	
2-Methylnaphthalene	ND	ug/Sample	30	
Hexachlorocyclopentadiene	ND	ug/Sample	30	
2,4,6-Trichlorophenol	ND	ug/Sample	30	
2,4,5-Trichlorophenol	ND	ug/Sample	150	
2-Chloronaphthalene	ND	ug/Sample	30	
2-Nitroaniline	ND	ug/Sample	30	
Dimethyl phthalate	ND	ug/Sample	30	
Acenaphthylene	ND	ug/Sample	150	
3-Nitroaniline	ND	ug/Sample	30	
Acenaphthene	ND	ug/Sample	150	
2,4-Dinitrophenol	ND	ug/Sample	150	
4-Nitrophenol	ND	ug/Sample	150	
Dibenzofuran	ND	ug/Sample	30	

Note R = Reporting limit(s) raised due to sample volume limitations.
ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.
Rev 230787

Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: T-MM5-4-F, FH, XAD, COND, BH
LAB ID: 300681-0005-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 26 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
2,4-Dinitrotoluene	ND	ug/Sample	30	
2,6-Dinitrotoluene	ND	ug/Sample	30	
Diethyl phthalate	ND	ug/Sample	30	
4-Chlorophenyl phenyl ether	ND	ug/Sample	30	
Fluorene	ND	ug/Sample	30	
4-Nitroaniline	ND	ug/Sample	150	
4,6-Dinitro-2-methylphenol	ND	ug/Sample	150	
N-Nitrosodiphenylamine	ND	ug/Sample	30	
4-Bromophenyl phenyl ether	ND	ug/Sample	30	
Hexachlorobenzene	ND	ug/Sample	30	
Pentachlorophenol	ND	ug/Sample	150	
Phenanthrene	ND	ug/Sample	30	
Anthracene	ND	ug/Sample	30	
Di-n-butyl phthalate	ND	ug/Sample	30	
Fluoranthene	ND	ug/Sample	30	
Pyrene	ND	ug/Sample	30	
Butyl benzyl phthalate	ND	ug/Sample	30	
3,3'-Dichlorobenzidine	ND	ug/Sample	60	
Benzo(a)anthracene	ND	ug/Sample	30	
bis(2-Ethylhexyl)-phthalate	ND	ug/Sample	30	
Chrysene	ND	ug/Sample	30	
Di-n-octyl phthalate	ND	ug/Sample	30	
Benzo(b)fluoranthene	ND	ug/Sample	30	
Benzo(k)fluoranthene	ND	ug/Sample	30	
Benzo(a)pyrene	ND	ug/Sample	30	
Indeno(1,2,3-cd)pyrene	ND	ug/Sample	30	
Dibenz(a,h)anthracene	ND	ug/Sample	30	
Benzo(g,h,i)perylene	ND	ug/Sample	30	
Acetophenone	ND	ug/Sample	30	
4-Aminobiphenyl	ND	ug/Sample	150	
Aniline	ND	ug/Sample	30	
Benzidine	ND	ug/Sample	300	
3,3'-Dimethylbenzidine	ND	ug/Sample	60	
N-Nitrosodimethylamine	ND	ug/Sample	30	
N-Nitrosomorpholine	ND	ug/Sample	30	
Pentachloronitrobenzene (PCNB)	ND	ug/Sample	150	
o-Toluidine	ND	ug/Sample	60	

ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.
Rev 230787

954

Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: T-MM5-4-F, FH, XAD, COND, BH
LAB ID: 300681-0005-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 26 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
2-Methoxybenzenamine	ND	ug/Sample	--	K
Biphenyl	ND	ug/Sample	--	K
Chloroacetophenone	ND	ug/Sample	--	K
Cumene	ND	ug/Sample	--	K
DBCP (1,2-Dibromo-3-chloropropane)	ND	ug/Sample	--	K
Benzo(e)pyrene	ND	ug/Sample	--	K
N-N-Diethylaniline	ND	ug/Sample	--	K
Dimethylaniline	ND	ug/Sample	--	K
3,3'-Dimethoxybenzidine	ND	ug/Sample	--	K
Hydroquinone	ND	ug/Sample	--	K
4,4'-Methyl-bis(2-chloroaniline)	ND	ug/Sample	--	K
4-Nitrodiphenyl	ND	ug/Sample	--	K
Trifluralin	ND	ug/Sample	--	K

Surrogate	Recovery	Acceptable Range
Nitrobenzene-d5	62 %	45 - 107
2-Fluorobiphenyl	86 %	62 - 110
Terphenyl-d14	110 %	58 - 135
Phenol-d5	61 %	43 - 130
2-Fluorophenol	54 %	36 - 111
2,4,6-Tribromophenol	65 %	58 - 131

Note K = Identified by mass spectrum only; quantitation based on 1:1 response with internal standard.
ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.
Rev 230787

955

Semivolatiles Library Search (20 Compound TID)
Method 8270

Client Name: Pacific Environmental Services
Client ID: T-MM5-4-F, FH, XAD, COND, BH
LAB ID: 300681-0005-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 26 JUL 98
Prepared: NA

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Result	Units	Reporting Limit	Qualifier
Benzaldehyde	43	ug/Sample	--	0
Pentadecane	61	ug/Sample	--	0
Nonadecane	66	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	41	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	120	ug/Sample	--	0
Nonadecane	45	ug/Sample	--	0
Nonadecane	36	ug/Sample	--	0
Unknown	36	ug/Sample	--	
Unknown	97	ug/Sample	--	
Unknown	40	ug/Sample	--	
Tricosane	37	ug/Sample	--	0
4-Hydroxy-4-methyl-2-pentanone	320	ug/Sample	--	0
Unknown	63	ug/Sample	--	
Unknown	58	ug/Sample	--	
5-Eicosenene, (E) -	120	ug/Sample	--	0
Unknown	42	ug/Sample	--	
n-Nonane	350	ug/Sample	--	0
Unknown	63	ug/Sample	--	
3-Octadecene, (E)-	38	ug/Sample	--	0
Unknown	47	ug/Sample	--	

Note 0 = Or structurally similar compound (isomer).
NA = Not Applicable

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.
Rev 230787

950

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068105
Std Id: ST16980831

Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 21:00
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Surrogate Spike Recoveries
8270-G, LIMS 10/10/96

Surrogate	Surrogate Amount (ug)		% Recovery	
	Spiked	Measured	Measured	QC Limits
CS20 NITROBENZENE-D5	50.00	31.20	62.4	45 107
CS25 2-FLUOROBIPHENY	50.00	43.00	86.0	62 110
CS30 TERPHENYL-D14	50.00	55.00	110.	58 135
CS45 PHENOL-D5	100.0	61.10	61.1	43 130
CS50 2-FLUOROPHENOL	100.0	53.90	53.9	36 111
CS55 2,4,6-TRIBROMOP	100.0	65.00	65.0	58 131

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
HEXACHLOROBENZENE-C13		ND	UG/A	30.0
C310 N-NITROSODIMETHYLAMINE		ND	UG/A	30.0
PYRIDINE		ND	UG/A	60.0
2-PICOLINE		ND	UG/A	30.0
N-NITROSOMETHYLETHYLAMINE	153	30.0 ND	UG/A	30.0
METHYLMETHANESULFONATE		ND	UG/A	30.0
N-NITROSODIETHYLAMINE		ND	UG/A	30.0
ETHYLMETHANE SULFONATE		ND	UG/A	30.0
PENTACHLOROETHANE		ND	UG/A	30.0
C320 ANILINE		ND	UG/A	30.0
C315 PHENOL		ND	UG/A	30.0
C325 BIS(2-CHLOROETHYL)ETHER		ND	UG/A	30.0
C330 2-CHLOROPHENOL		ND	UG/A	30.0
C335 1,3-DICHLOROBENZENE		ND	UG/A	30.0
C340 1,4-DICHLOROBENZENE		ND	UG/A	30.0
C345 BENZYL ALCOHOL		ND	UG/A	30.0
C350 1,2-DICHLOROBENZENE		ND	UG/A	30.0
C355 2-METHYLPHENOL		ND	UG/A	30.0
C360 2,2'-OXYBIS(1-CLPROPAN)		ND	UG/A	30.0
C361 ACETOPHENONE		ND	UG/A	30.0
N-NITROSOPYRROLIDINE		ND	UG/A	30.0
N-NITROSOMORPHOLINE		ND	UG/A	30.0
3-METHYL PHENOL		ND	UG/A	30.0
C365 4-METHYLPHENOL		ND	UG/A	30.0
C370 N-NITROSO-DI-N-PROPYLAM		ND	UG/A	30.0
O-TOLUIDINE		ND	UG/A	30.0
C375 HEXACHLOROETHANE		ND	UG/A	30.0
C410 NITROBENZENE		ND	UG/A	30.0
N-NITROSOPIPERIDINE		ND	UG/A	30.0

Reviewed by: 144/17/96

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068105
Std Id: ST16980831

Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 21:00
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C415 ISOPHORONE		ND	UG/A	30.0
C420 2-NITROPHENOL		ND	UG/A	30.0
C425 2,4-DIMETHYLPHENOL		ND	UG/A	30.0
C435 BIS(2-CHLOROETHOXY)METH		ND	UG/A	30.0
C440 2,4-DICHLOROPHENOL		ND	UG/A	30.0
C430 BENZOIC ACID		ND	UG/A	150.0
AA-DIMETHYLPHENETHYLAMINE		ND	UG/A	30.0
C445 1,2,4-TRICHLOROBENZENE		ND	UG/A	30.0
C450 NAPHTHALENE		ND	UG/A	30.0
C455 4-CHLOROANILINE		ND	UG/A	30.0
2,6-DICHLOROPHENOL		ND	UG/A	30.0
HEXACHLOROPROPENE		ND	UG/A	30.0
C460 HEXACHLOROBUTADIENE		ND	UG/A	30.0
P-PHENYLENE DIAMINE		ND	UG/A	30.0
N-NITROSODI-N-BUTYLAMINE		ND	UG/A	30.0
C465 4-CHLORO-3-METHYLPHENO		ND	UG/A	30.0
SAFROLE		ND	UG/A	30.0
C470 2-METHYLNAPHTHALENE		ND	UG/A	30.0
1,2,4,5-TETRACHLOROBENZENE		ND	UG/A	30.0
ISOSAFROLE (#1)		ND	UG/A	60.0
C510 HEXACHLOROCYCLOPENTADI		ND	UG/A	30.0
C515 2,4,6-TRICHLOROPHENOL		ND	UG/A	30.0
C520 2,4,5-TRICHLOROPHENOL		ND	UG/A	30.0
ISOSAFROLE (#2)		ND	UG/A	60.0
C525 2-CHLORONAPHTHALENE		ND	UG/A	30.0
1-CHLORONAPHTHALENE		ND	UG/A	15.0
C530 2-NITROANILINE		ND	UG/A	150.0
1,4-NAPHTHOQUINONE		ND	UG/A	30.0
C535 DIMETHYLPHTHALATE		ND	UG/A	30.0
1,3-DINITROBENZENE		ND	UG/A	30.0
C540 ACENAPHTHYLENE		ND	UG/A	30.0
C543 2,6-DINITROTOLUENE		ND	UG/A	30.0
C545 3-NITROANILINE		ND	UG/A	150.0
C550 ACENAPHTHENE		ND	UG/A	30.0
C555 2,4-DINITROPHENOL		ND	UG/A	150.0
C565 DIBENZOFURAN		ND	UG/A	30.0
C560 4-NITROPHENOL		ND	UG/A	150.0
PENTACHLOROBENZENE		ND	UG/A	30.0
C570 2,4-DINITROTOLUENE		ND	UG/A	30.0
1-NAPHTHYLAMINE		ND	UG/A	30.0
2-NAPHTHYLAMINE		ND	UG/A	30.0
2,3,4,6-TETRACHLOROPHENOL		ND	UG/A	60.0
C580 DIETHYLPHTHALATE		ND	UG/A	30.0
C590 FLUORENE		ND	UG/A	30.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068105
Std Id: ST16980831

Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 21:00
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C585 4-CHLOROPHENYL-PHENYLE		ND	UG/A	30.0
5-NITRO-O-TOLUIDINE		ND	UG/A	30.0
C595 4-NITROANALINE		ND	UG/A	150.0
C610 4,6-DINITRO-2-METHYLPH		ND	UG/A	150.0
C615 N-NITROSODIPHENYLAMINE		ND	UG/A	30.0
C620 AZOBENZENE		ND	UG/A	30.0
SYM-TRINITROBENZENE		ND	UG/A	30.0
C625 4-BROMOPHENYL-PHENYLET		ND	UG/A	30.0
PHENACETIN		ND	UG/A	30.0
DIALATE	234	ND	UG/A	30.0
C630 HEXACHLOROBENZENE		ND	UG/A	30.0
4-AMINOBIIPHENYL		ND	UG/A	30.0
C635 PENTACHLOROPHENOL		ND	UG/A	150.0
PRONAMIDE		ND	UG/A	30.0
PENTACHLORONITROBENZENE		ND	UG/A	150.0
C640 PHENANTHRENE		ND	UG/A	30.0
C645 ANTHRACENE		ND	UG/A	30.0
2SECBUTYL-4,6-DINITROPHENOL		ND	UG/A	30.0
C647 CARBAZOLE		ND	UG/A	30.0
C650 DI-N-BUTYLPHTHALATE		ND	UG/A	30.0
4-NITROQUINOLINE-1-OXIDE		ND	UG/A	30.0
METHAPYRILENE		ND	UG/A	30.0
ISODRIN		ND	UG/A	30.0
C655 FLUORANTHENE		ND	UG/A	30.0
CHLOROBENZILATE		ND	UG/A	30.0
C710 BENZIDINE		ND	UG/A	300.0
C715 PYRENE		ND	UG/A	30.0
ARAMITE (#1)		ND	UG/A	30.0
ARAMITE (#2)		ND	UG/A	30.0
P-DIMETHYLAMINOAZOBENZENE		ND	UG/A	30.0
3,3'-DIMETHYLBENZIDINE		ND	UG/A	30.0
KEPONE		ND	UG/A	150.0
C720 BUTYLBENZYLPHTHALATE		ND	UG/A	30.0
2-ACETYLAMINOFLUORENE		ND	UG/A	30.0
C730 BENZO(A)ANTHRACENE		ND	UG/A	30.0
C725 3,3'-DICHLOBENZIDINE		ND	UG/A	60.0
C740 CHRYSENE		ND	UG/A	30.0
C745 BIS(2-ETHYLHEXYL)PHTHA		ND	UG/A	30.0
3-METHYLCHOLANTHRENE		ND	UG/A	30.0
C760 DI-N-OCTYL PHTHALATE		ND	UG/A	30.0
C765 BENZO(B)FLUORANTHENE		ND	UG/A	30.0
7,12-DIMETHYLBENZANTHRACENE		ND	UG/A	30.0
C770 BENZO(K)FLUORANTHENE		ND	UG/A	30.0
HEXACHLOROPHENE		ND	UG/A	30.0

QUANTERRA GC/MS

Target Compound Data Summary Sheet

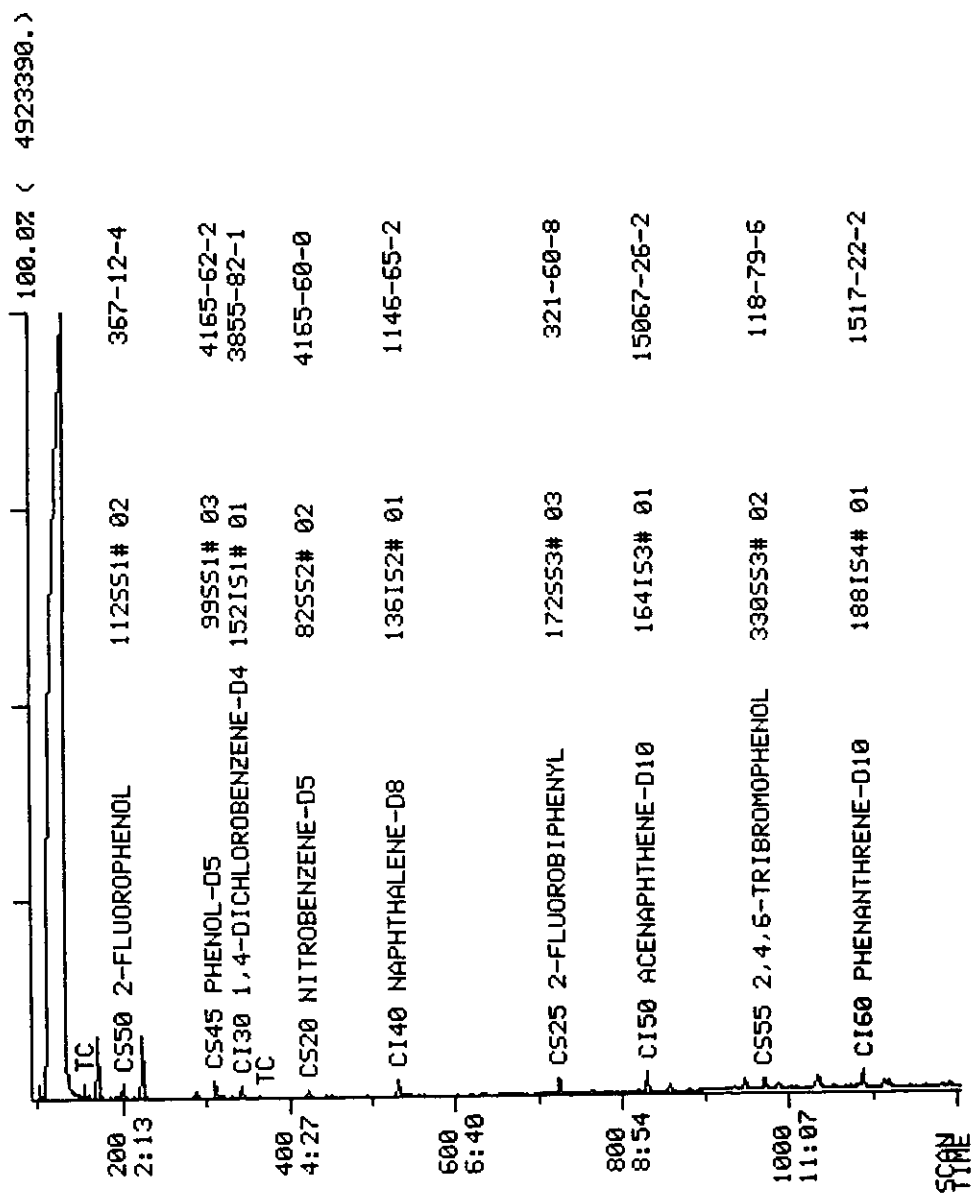
Data File: 30068105
Std Id: ST16980831

Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 21:00
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

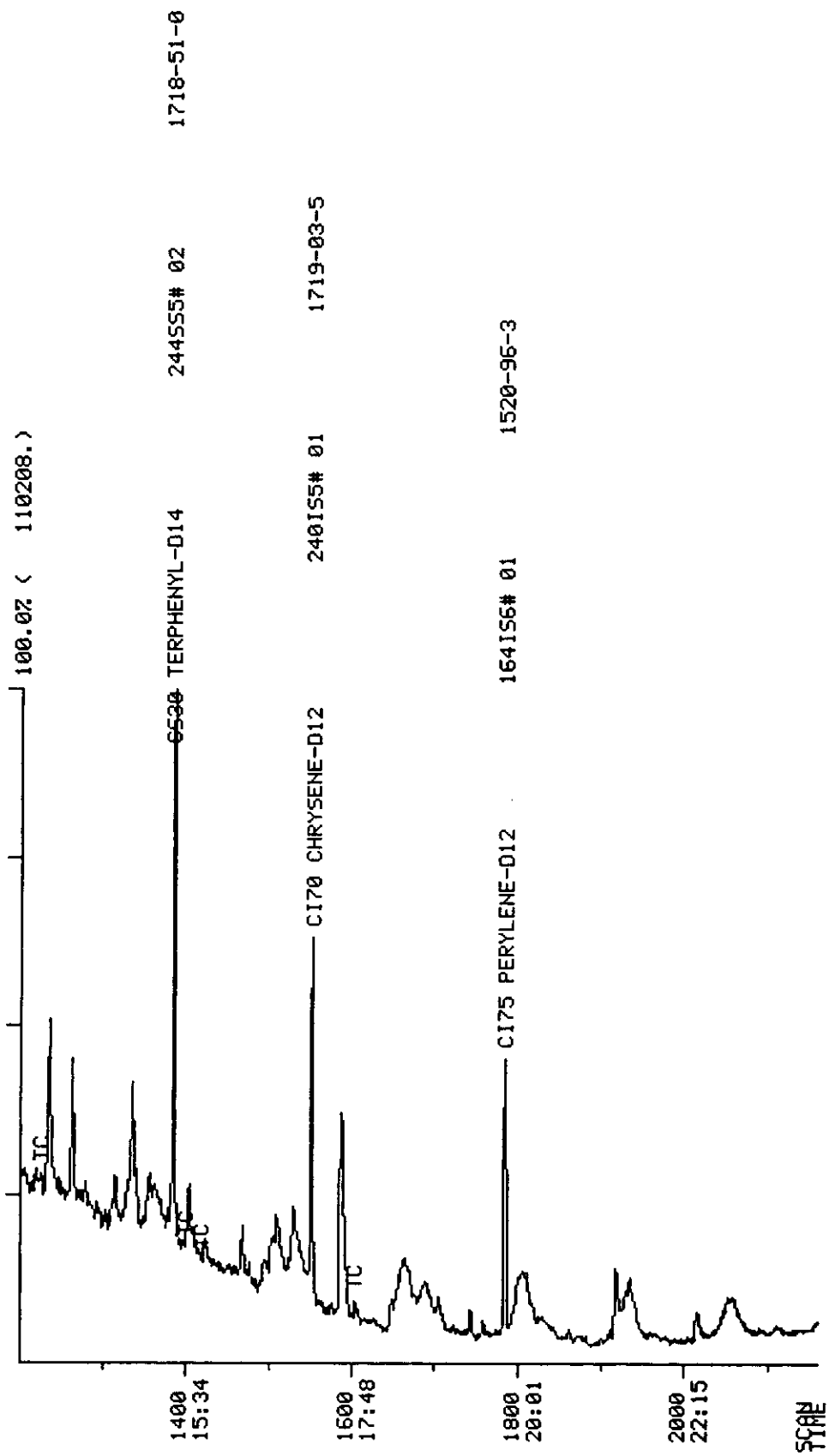
Parameter	Scan	Result	Units	Reporting Limit
=====				
C775 BENZO(A)PYRENE		ND	UG/A	30.0
C780 INDENO(1,2,3-CD)PYRENE		ND	UG/A	30.0
C785 DIBENZ(A,H)ANTHRACENE		ND	UG/A	30.0
C790 BENZO(G,H,I)PERYLENE		ND	UG/A	30.0

DATA FROM FILE: 30068105 SCANS 93 TO 1203 ACQUIRED: 08/31/98 21:00:00
 CALI: 30068105 #3
 SAMPLE: T-MM5-4 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M



5
 (7)
 1.1

DATA FROM FILE: 30068105 SCANS 1203 TO 2158 ACQUIRED: 08/31/98 21:00:00
 CALI: 30068105 #3
 SAMPLE: T-HMS-4 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M



ADDITIONAL STANDARD USED FOR THIS REPORT IS: ST16980831A

QUANTERRA QUANTITATION SUMMARY

File: 30068105

Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Analyst: DAT Instrument Id: F16 Analyzed: 08/31/98 21:00

Compounds with amounts less than 0.20 reported as NOT FOUND

Library							Units: UG/ML		
No	Entry	Name	Mass	Meth	Scan	Ref	Area	RRF(L)	Amount
1	S1#	1 C130 1,4-DICHLOROBENZENE	152	A	BB	340 1	35775.	1.000	40.000
2	S2#	1 C140 NAPHTHALENE-D8	136	A	BB	529 2	135276.	1.000	40.000
3	S3#	1 C150 ACENAPHTHENE-D10	164	A	BB	828 3	72725.	1.000	40.000
4	S4#	1 C160 PHENANTHRENE-D10	188	A	BB	1086 4	117535.	1.000	40.000
5	S5#	1 C170 CHRYSENE-D12	240	A	BB	1551 5	55639.	1.000	40.000
6	S6#	1 C175 PERYLENE-D12	264	A	BB	1782 6	42978.	1.000	40.000
7	S2#	2 CS20 NITROBENZENE-D5	82	A	BB	421 2	49476.	0.470	31.156
8	S3#	3 CS25 2-FLUOROBIPHENYL	172	A	BB	722 3	96577.	1.234	43.036
9	S5#	2 CS30 TERPHENYL-D14	244	A	BB	1383 5	76073.	0.995	54.953
10	S1#	3 CS45 PHENOL-D5	99	A	BB	308 1	110653.	2.025	61.081
11	S1#	2 CS50 2-FLUOROPHENOL	112	A	BB	198 1	65971.	1.368	53.910
12	S3#	2 CS55 2,4,6-TRIBROMOPHENO	330	A	BV	968 3	27845.	0.236	64.989
13	S4#	4 HEXACHLOROBENZENE-C13	294			4 NOT FOUND			
14	S1#	15 C310 N-NITROSODIMETHYLAM	74			1 NOT FOUND			
15	S1#	20 PYRIDINE	79			1 NOT FOUND			
16	S1#	30 2-PICOLINE	93			1 NOT FOUND			
17	S1#	40 N-NITROSOMETHYLETHYLAMIN	42	A	BV	153 1	7174.	0.811	9.890
18	S1#	60 METHYLMETHANESULFONATE	80			1 NOT FOUND			
19	S1#	70 N-NITROSODIETHYLAMINE	102	A	BB	212 1	332.	0.521	0.712
20	S1#	85 ETHYLMETHANE SULFONATE	79	A	BB	253 1	338.	1.108	0.341
21	S1#	95 PENTACHLOROETHANE	117			1 NOT FOUND			
22	S1#	100 C320 ANILINE	93	A	BB	301 1	408.	1.846	0.247
23	S1#	105 C315 PHENOL	94	A	BV	309 1	1292.	1.968	0.734
24	S1#	110 C325 BIS(2-CHLOROETHYL)E	93			1 NOT FOUND			
25	S1#	115 C330 2-CHLOROPHENOL	128			1 NOT FOUND			
26	S1#	125 C335 1,3-DICHLOROBENZENE	146			1 NOT FOUND			
27	S1#	130 C340 1,4-DICHLOROBENZENE	146	A	BB	342 1	812.	1.617	0.561
28	S1#	145 C345 BENZYL ALCOHOL	108	A	BB	370 1	4056.	0.854	5.308
29	S1#	150 C350 1,2-DICHLOROBENZENE	146			1 NOT FOUND			
30	S1#	160 C355 2-METHYLPHENOL	108			1 NOT FOUND			
31	S1#	165 C360 2,2'-OXYBIS(1-CLPRO	45			1 NOT FOUND			
32	S1#	170 C361 ACETOPHENONE	105	A	BB	400 1	1325.	1.934	0.766
33	S1#	175 N-NITROSOPYRROLIDINE	100			1 NOT FOUND			
34	S1#	180 N-NITROSOMORPHOLINE	56			1 NOT FOUND			
35	S1#	182 3-METHYL PHENOL	108	A	BB	422 1	825.	2.252	0.410
36	S1#	185 C365 4-METHYLPHENOL	108			1 NOT FOUND			
37	S1#	190 C370 N-NITROSO-DI-N-PROP	70			1 NOT FOUND			
38	S1#	195 O-TOLUIDINE	106			1 NOT FOUND			
39	S1#	200 C375 HEXACHLOROETHANE	117			1 NOT FOUND			
40	S2#	10 C410 NITROBENZENE	77			2 NOT FOUND			
41	S2#	15 N-NITROSOPIPERIDINE	42	A	BB	449 2	348.	0.320	0.321
42	S2#	20 C415 ISOPHORONE	82			2 NOT FOUND			
43	S2#	25 C420 2-NITROPHENOL	139			2 NOT FOUND			
44	S2#	30 C425 2,4-DIMETHYLPHENOL	107			2 NOT FOUND			
45	S2#	35 C435 BIS(2-CHLOROETHOXY)	93			2 NOT FOUND			

46 S2# 40 C440 2,4-DICHLOROPHENOL	162	2	NOT FOUND
47 S2# 45 C430 BENZOIC ACID	122	2	NOT FOUND
48 S2# 50 AA-DIMETHYLPHENETHYLAMIN	58	2	NOT FOUND
49 S2# 55 C445 1,2,4-TRICHLOROBENZ	180	2	NOT FOUND

50	S2# 60	C450 NAPHTHALENE	128 A BB	532 2	2438. 1.022	0.705
51	S2# 80	C455 4-CHLOROANALINE	127	2	NOT FOUND	
52	S2# 85	2,6-DICHLOROPHENOL	162	2	NOT FOUND	
53	S2# 90	HEXACHLOROPROPENE	213	2	NOT FOUND	
54	S2# 95	C460 HEXACHLOROBUTADIENE	225	2	NOT FOUND	
55	S2#115	P-PHENYLENE DIAMINE	108	2	NOT FOUND	
56	S2#120	N-NITROSODI-N-BUTYLAMINE	84	2	NOT FOUND	
57	S2#130	C465 4-CHLORO-3-METHYLPH	107 A BB	649 2	496. 0.338	0.434
58	S2#140	SAFROLE	162	2	NOT FOUND	
59	S2#145	C470 2-METHYLNAPHTHALENE	142 A BB	651 2	2124. 0.751	0.836
60	S3# 10	1,2,4,5-TETRACHLOROBENZE	214	3	NOT FOUND	
61	S3# 15	ISOSAFROLE (#1)	162	3	NOT FOUND	
62	S3# 20	C510 HEXACHLOROCYCLOPENT	237	3	NOT FOUND	
63	S3# 25	C515 2,4,6-TRICHLOROPHEN	196	3	NOT FOUND	
64	S3# 30	C520 2,4,5-TRICHLOROPHEN	196	3	NOT FOUND	
65	S3# 35	ISOSAFROLE (#2)	104	3	NOT FOUND	
66	S3# 40	C525 2-CHLORONAPHTHALENE	162	3	NOT FOUND	
67	S3# 42	1-CHLORONAPHTHALENE	162	3	NOT FOUND	
68	S3# 45	C530 2-NITROANALINE	65	3	NOT FOUND	
69	S3# 50	1,4-NAPHTHOQUINONE	158	3	NOT FOUND	
70	S3# 55	C535 DIMETHYLPHTHALATE	163 A BB	804 3	549. 1.341	0.225
71	S3# 60	1,3-DINITROBENZENE	168	3	NOT FOUND	
72	S3# 65	C540 ACENAPHTHYLENE	152	3	NOT FOUND	
73	S3# 70	C543 2,6-DINITROTOLUENE	165	3	NOT FOUND	
74	S3# 75	C545 3-NITROANILINE	138	3	NOT FOUND	
75	S3# 80	C550 ACENAPHTHENE	153	3	NOT FOUND	
76	S3# 85	C555 2,4-DINITROPHENOL	184	3	NOT FOUND	
77	S3# 90	C565 DIBENZOFURAN	168	3	NOT FOUND	
78	S3# 95	C560 4-NITROPHENOL	109	3	NOT FOUND	
79	S3#100	PENTACHLOROBENZENE	250	3	NOT FOUND	
80	S3#105	C570 2,4-DINITROTOLUENE	165	3	NOT FOUND	
81	S3#110	1-NAPHTHYLAMINE	143	3	NOT FOUND	
82	S3#115	2-NAPHTHYLAMINE	143	3	NOT FOUND	
83	S3#120	2,3,4,6-TETRACHLOROPHENO	232	3	NOT FOUND	
84	S3#130	C580 DIETHYLPHTHALATE	149	3	NOT FOUND	
85	S3#135	C590 FLUORENE	166 A BB	925 3	880. 1.298	0.373
86	S3#140	C585 4-CHLOROPHENYL-PHEN	204	3	NOT FOUND	
87	S3#145	5-NITRO-O-TOLUIDINE	152 A VB	940 3	425. 0.356	0.658
88	S3#150	C595 4-NITROANALINE	138	3	NOT FOUND	
89	S4# 10	C610 4,6-DINITRO-2-METHY	198	4	NOT FOUND	
90	S4# 15	C615 N-NITROSODIPHENYLAM	169 A BB	956 4	479. 0.551	0.296
91	S4# 20	C620 AZOBENZENE	77	4	NOT FOUND	
92	S4# 25	SYM-TRINITROBENZENE	75 A VV	1023 4	404. 0.162	0.851
93	S4# 30	C625 4-BROMOPHENYL-PHENY	248	4	NOT FOUND	
94	S4# 35	PHENACETIN	108	4	NOT FOUND	
95	S4# 37	DIALATE	234	4	NOT FOUND	
96	S4# 40	C630 HEXACHLOROBENZENE	284	4	NOT FOUND	
97	S4# 45	4-AMINOBIIPHENYL	169	4	NOT FOUND	
98	S4# 50	C635 PENTACHLOROPHENOL	266	4	NOT FOUND	
99	S4# 55	PRONAMIDE	173	4	NOT FOUND	
100	S4# 60	PENTACHLORONITROBENZENE	237	4	NOT FOUND	
101	S4# 65	C640 PHENANTHRENE	178 A BB	1090 4	3844. 1.033	1.267
102	S4# 70	C645 ANTHRACENE	178	4	NOT FOUND	
103	S4# 75	2SECBUTYL-4,6-DINITROPHE	211	4	NOT FOUND	
104	S4# 80	C647 CARBAZOLE	167	4	NOT FOUND	
105	S4# 85	C650 DI-N-BUTYLPHTHALATE	149 A BB	1226 4	3511. 1.247	0.959
106	S4#100	4-NITROQUINOLINE-1-OXIDE	190	4	NOT FOUND	
107	S4#105	METHAPYRILENE	58 A BB	1264 4	552. 0.305	0.617
108	S4#106	ISODRIN	193	4	NOT FOUND	
109	S4#110	C655 FLUORANTHENE	202	4	NOT FOUND	
110	S4#120	CHLOROBENZILATE	139	4	NOT FOUND	

111	S5# 10	C710 BENZIDINE	184	5	NOT FOUND		
112	S5# 15	C715 PYRENE	202	A BB 1335	5	572.	1.323 0.311
113	S5# 20	ARAMITE (#1)	185	A BV 1398	5	96.	0.061 1.125
114	S5# 25	ARAMITE (#2)	185	A VB 1417	5	346.	0.087 2.871
115	S5# 30	P-DIMETHYLAMINOAZOBENZEN	120	5	NOT FOUND		
116	S5# 35	3,3'-DIMETHYLBENZIDINE	212	5	NOT FOUND		
117	S5# 37	KEPONE	272	5	NOT FOUND		
118	S5# 40	C720 BUTYLBENZYLPHTHALAT	149	A BB 1487	5	234.	0.685 0.246
119	S5# 45	2-ACETYLAMINOFLUORENE	181	5	NOT FOUND		
120	S5# 50	C730 BENZO(A)ANTHRACENE	228	A BB 1552	5	438.	1.125 0.280
121	S5# 55	C725 3,3'-DICHLOROBENZID	252	5	NOT FOUND		
122	S5# 60	C740 CHRYSENE	228	5	NOT FOUND		
123	S5# 65	C745 BIS(2-ETHYLHEXYL)PH	149	A BB 1603	5	2203.	0.980 1.616
124	S5# 85	3-METHYLCHOLANTHRENE	268	5	NOT FOUND		
125	S6# 10	C760 DI-N-OCTYL PHTHALAT	149	6	NOT FOUND		
126	S6# 15	C765 BENZO(B)FLUORANTHEN	252	6	NOT FOUND		
127	S6# 20	7,12-DIMETHYLBENZANTHRAC	256	6	NOT FOUND		
128	S6# 25	C770 BENZO(K)FLUORANTHEN	252	6	NOT FOUND		
129	S6# 30	HEXACHLOROPHENE	196	6	NOT FOUND		
130	S6# 35	C775 BENZO(A)PYRENE	252	6	NOT FOUND		
131	S6# 55	C780 INDENO(1,2,3-CD)PYR	276	6	NOT FOUND		
132	S6# 60	C785 DIBENZ(A,H)ANTHRACE	278	6	NOT FOUND		
133	S6# 65	C790 BENZO(G,H,I)PERYLEN	276	6	NOT FOUND		

BNA, S LIBRARIES

ANALYST: _____ DATE: _____

QUANTERRA Internal Standard Check

Standard Filename: ST16980831

Analyzed: 08/31/98 12:31

Standard	IS# 1			IS# 2			IS# 3		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	28649.		3.78	100432.		5.90	54347.		9.21
Upper Limit	57297.=+200		4.28	200864.=+200		6.40	108695.=+200		9.71
Lower Limit	14324.=+ 50		3.28	50216.=+ 50		5.40	27174.=+ 50		8.71
Filename									
1 30068105	35775.= 124		3.78	135276.= 134		5.88	72725.= 133		9.21

IS# 1 = C130 1,4-DICHLOROBENZENE-D4

IS# 2 = C140 NAPHTHALENE-D8

IS# 3 = C150 ACENAPHTHENE-D10

* - indicates an I.S. is outside QC limit(s)

Standard	IS# 4			IS# 5			IS# 6		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	98520.		12.08	62289.		17.26	50750.		19.83
Upper Limit	197039.=+200		12.58	124578.=+200		17.76	101500.=+200		20.33
Lower Limit	49260.=+ 50		11.58	31145.=+ 50		16.76	25375.=+ 50		19.33
Filename									
1 30068105	117535.= 119		12.08	55639.= 89		17.25	42978.= 84		19.82

IS# 4 = C160 PHENANTHRENE-D10

IS# 5 = C170 CHRYSENE-D12

IS# 6 = C175 PERYLENE-D12

* - indicates an I.S. is outside QC limit(s)

Data Reduced by: *an* Date: *9/17/98*
Data Reviewed by: *V* Date: *9/17/98*

Data File: 30068105

QUANTERRA GC/MS TIC REPORT (Part 1)

Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
Analyst: DAT Date Analyzed: 08/31/98 21:00
Run Factor: 3.00

Concentration

in Sample

yes
JOA
SCAN (UG/A) CAS #
1 170 320. ~~123-42-2~~
~~2-PENTANONE, 4-HYDROXY-4-METHYL-~~ *5700*

yes
JOA
2 222 350. 111-84-2
NONANE or *Bomer*

3 285 43. 100-52-7
BENZALDEHYDE or *Bomer*

X 761 31. 629-59-4
TETRADECANE or *Bomer*

5 856 61. 629-62-9
PENTADECANE

6 946 66. 629-92-5
NONADECANE

7 986 41. 54105-67-8
HEPTADECANE, 2,6-DIMETHYL-

8 1032 120. 54105-67-8
HEPTADECANE, 2,6-DIMETHYL-

9 1113 45. 629-92-5
NONADECANE

10 1117 36. 629-92-5
NONADECANE

~~11~~ 1190 29. 629-92-5
NONADECANE or isomer

12 1236 36. 57-10-3
HEXADECANOIC ACID

unknown. 5700

13 1334 97. 00-00-0
UNKNOWN

14 1360 40. 00-00-0
UNKNOWN

15 1402 37. 638-67-5
TRICOSANE or isomer

~~16~~ 1467 31. 00-00-0
UNKNOWN alkane

5746400

17 1508 63. 00-00-0
UNKNOWN

5700

18 1529 58. 00-00-0
UNKNOWN

19 1587 120. 74685-30-6
5-EICOSENE, (E)- or isomer

20 1662 42. 00-00-0
UNKNOWN

5700

~~21~~ 1688 32. 00-00-0
UNKNOWN

22 1804 63. 00-00-0
UNKNOWN

23 1916	38.	7206-19-1
3-OCTADECENE, (E)-		<i>or Bomer</i>

24 1933	47.	00-00-0
UNKNOWN		

25 2057	29.	00-00-0
UNKNOWN		

DATA FILE: 30068105

QUANTERRA GC/MS TIC REPORT (Part 2)

CONCENTRATION = AREA(TIC)*CONC(IS)/AREA(IS)

#	FIT	PURITY	INT.		RRT	AREA	HEIGHT	AMOUNT		LIB	LIB #
			STD.	RT				AS ANALYZED	(UG/ML)		
1	956	919	1	1:53	0.321	915820.	389859.	108.038	NB	3241.	
2	986	938	1	2:28	0.420	976832.	382976.	115.236	NB	5159.	
3	975	926	1	3:10	0.539	120816.	35462.	14.253	NB	2017.	
4	960	762	2	8:27	0.919	86560.	25764.	10.201	NB	22530.	
5	965	792	2	9:31	1.034	171391.	43553.	20.199	NB	25997.	
6	963	834	2	10:31	1.143	185824.	58432.	21.900	NB	37465.	
7	957	814	3	10:58	0.908	134784.	36544.	13.810	NB	37462.	
8	972	850	3	11:28	0.950	391552.	70400.	40.118	NB	37462.	
9	939	785	3	12:22	1.025	144896.	41984.	14.846	NB	37465.	
10	941	811	3	12:25	1.029	116416.	38400.	11.928	NB	37465.	
11	971	746	3	13:14	1.096	92992.	30528.	9.528	NB	37465.	
12	975	682	3	13:44	1.138	118352.	27355.	12.126	NB	35182.	
13	881	596	4	14:50	0.860	144384.	22400.	32.404	UK	1.	
14	725	270	4	15:07	0.877	58976.	5928.	13.236	UK	1.	
15	890	654	4	15:35	0.904	55232.	10688.	12.396	NB	46161.	
16	891	588	4	16:19	0.946	45440.	8408.	10.198	UK	1.	
17	772	423	4	16:46	0.972	93120.	9152.	20.899	UK	1.	
18	913	506	4	17:00	0.986	85616.	10128.	19.214	UK	1.	
19	974	771	4	17:39	1.023	177728.	31680.	39.887	NB	39516.	
20	916	471	4	18:29	1.072	62400.	6960.	14.004	UK	1.	
21	646	240	5	18:46	0.947	37296.	3821.	10.683	UK	1.	
22	799	436	5	20:03	1.012	73392.	5788.	21.023	UK	1.	
23	930	699	5	21:18	1.075	44560.	9603.	12.764	NB	34410.	
24	677	407	5	21:30	1.085	54336.	5944.	15.565	UK	1.	
25	704	420	5	22:52	1.154	33762.	4298.	9.671	UK	1.	

Library Search Data: 30068105 # 170 Base m/z: 43
 08/31/98 21:00:00 + 1:53 Cali: 30068105 # 3 RIC: 278528.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 617 matched at least 6 of the 16 largest peaks in the unknown

Rank In.	Name
1	3241 2-PENTANONE, 4-HYDROXY-4-METHYL-
2	5791 1,3-DIOXOLANE-2-METHANOL, 2,4-DIMETHYL-
3	3366 2-PENTANOL, 2,4-DIMETHYL-
4	3288 ACETIC ACID, 1,1-DIMETHYLETHYL ESTER
5	919 TERT-BUTYL HYDROPEROXIDE
6	5594 2-PROPANOL, 2-NITROSO-, ACETATE (ESTER)
7	3349 2-HEXANOL, 2-METHYL-
8	8432 2-HEXANONE, 3-HYDROXY-3,5-DIMETHYL-
9	8425 2-HEPTANONE, 3-HYDROXY-3-METHYL-

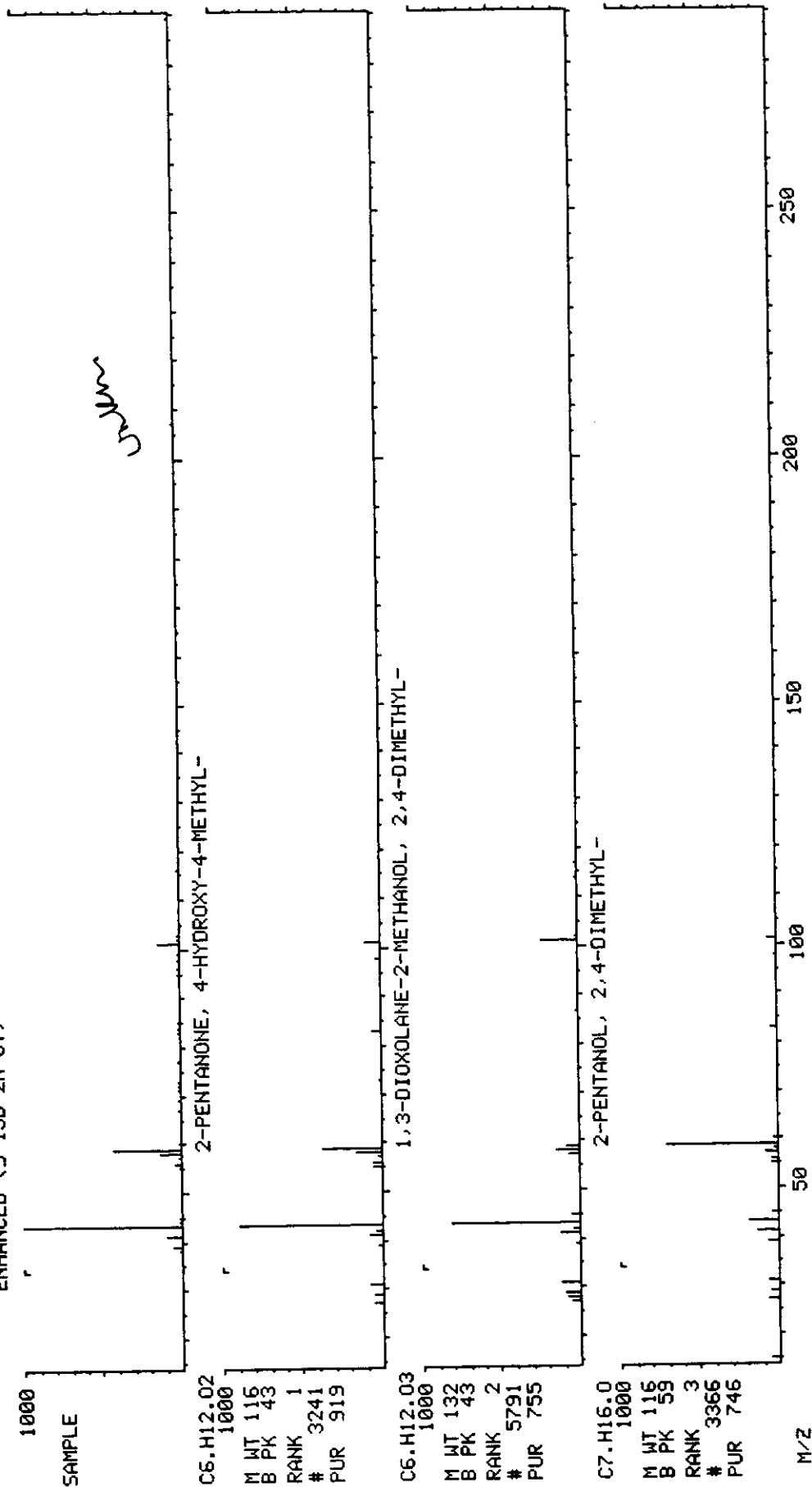
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C6.H12.O2	116	43	919	956	932
2	C6.H12.O3	132	43	755	828	798
3	C7.H16.O	116	59	746	768	746
4	C6.H12.O2	116	43	729	777	742
5	C4.H10.O2	90	59	714	866	736
6	C5.H9.O3.N	131	43	711	824	763
7	C7.H16.O	116	59	708	741	708
8	C8.H16.O2	144	59	667	729	694
9	C8.H16.O2	144	59	661	715	677

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	123-42-2
2	—	—	—	—	53951-43-2
3	—	—	—	—	625-06-9
4	—	—	—	—	540-88-5
5	—	—	—	—	75-91-2
6	—	—	—	—	6931-04-0
7	—	—	—	—	625-23-0
8	—	—	—	—	6321-14-8
9	—	—	—	—	13757-91-0

DATA: 30068105 # 170
CALI: 30068105 # 3

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 21:00:00 + 1:53
SAMPLE: T-MMS-4 1/35A/1ML
COND.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 158 2N 0T)

BASE M/Z: 43
RIC: 278528.



Library Search Data: 30068105 # 222 Base m/z: 43
 08/31/98 21:00:00 + 2:28 Cali: 30068105 # 3 RIC: 282112.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (\$ 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 921 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 5159 NONANE
 2 5151 HEPTANE, 3,4-DIMETHYL-
 3 5141 HEPTANE, 2,4-DIMETHYL-
 4 3086 HEXANE, 2,4-DIMETHYL-
 5 3081 OCTANE
 6 11607 UNDECANE
 7 19015 DECANE, 2,5,6-TRIMETHYL-
 8 5154 HEXANE, 4-ETHYL-2-METHYL-
 9 19028 DECANE, 2,4,6-TRIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C9.H20	128	43	938	986	938
2	C9.H20	128	43	912	940	912
3	C9.H20	128	43	892	944	893
4	C8.H18	114	43	884	956	894
5	C8.H18	114	43	870	945	893
6	C11.H24	156	43	867	926	905
7	C13.H28	184	57	866	906	876
8	C9.H20	128	57	856	907	856
9	C13.H28	184	43	853	884	898

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	111-84-2
2	—	—	—	—	922-28-1
3	—	—	—	—	2213-23-2
4	—	—	—	—	589-43-5
5	—	—	—	—	111-65-9
6	—	—	—	—	1120-21-4
7	—	—	—	—	62108-23-0
8	—	—	—	—	3074-75-7
9	—	—	—	—	62108-27-4

BASE M/Z: 43
RIC: 282112.

DATA: 30068105 # 222
CALI: 30068105 # 3

MID LIBRARY SEARCH (LIBRARYNB)

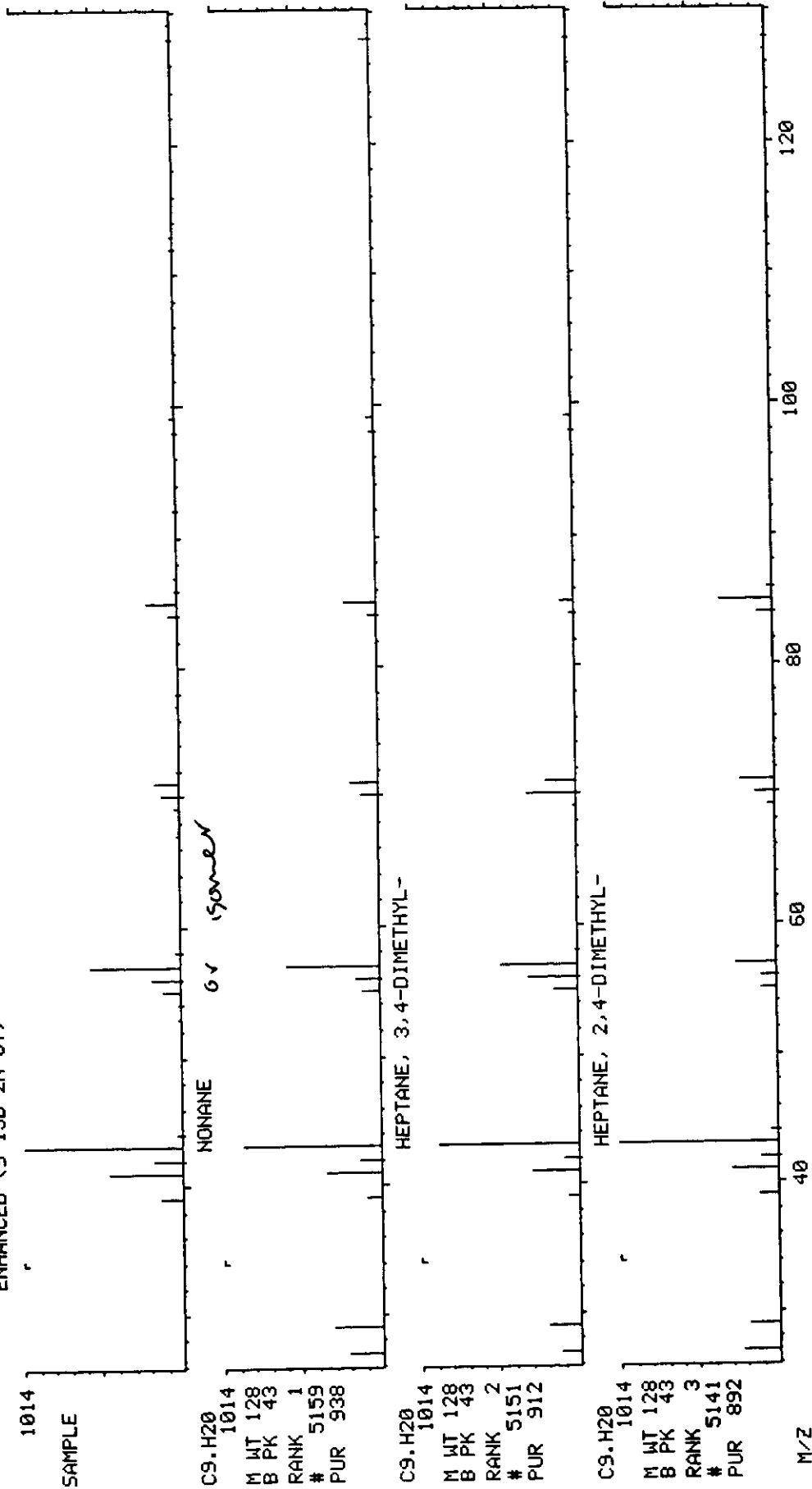
08/31/98 21:00:00 + 2:28

SAMPLE: T-MM5-4 1/35A/1ML

INST. ID: F16

CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M

ENHANCED (S 158 2N 0T)



Library Search Data: 30068105 # 285 Base m/z: 77
 08/31/98 21:00:00 + 3:10 Cali: 30068105 # 3 RIC: 28832.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

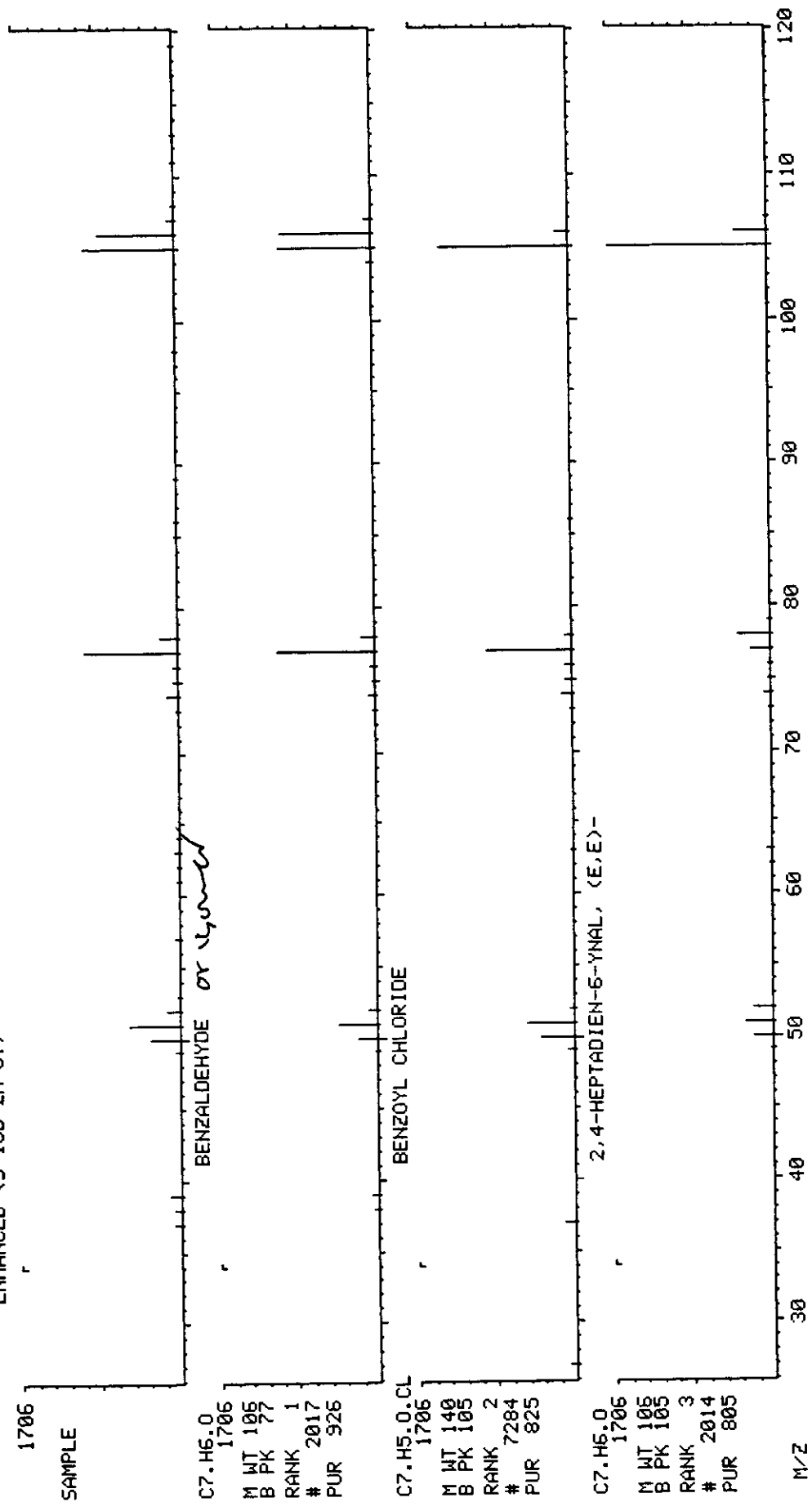
62231 spectra in LIBRARYNB searched for maximum PURITY
 416 matched at least 5 of the 16 largest peaks in the unknown

Rank In.	Name
1	2017 BENZALDEHYDE
2	7284 BENZOYL CHLORIDE
3	2014 2,4-HEPTADIEN-6-YNAL, (E,E)-
4	6466 ETHANONE, 2-HYDROXY-1-PHENYL-
5	13112 BENZOYL ISOTHIOCYANATE
6	6857 BENZENECARBOETHIOIC ACID
7	9645 BENZOYLFORMIC ACID
8	6111 1,2-BENZENEDICARBOXALDEHYDE
9	13359 BENZENEACETIC ACID, .ALPHA.-OXO-, METHYL ESTER

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C7.H6.O	106	77	926	975	941
2	C7.H5.O.CL	140	105	825	889	825
3	C7.H6.O	106	105	805	854	810
4	C8.H8.O2	136	105	793	842	812
5	C8.H5.O.N.S	163	105	784	835	790
6	C7.H6.O.S	138	77	779	865	805
7	C8.H6.O3	150	105	775	849	800
8	C8.H6.O2	134	105	773	825	908
9	C9.H8.O3	164	105	773	846	791

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	100-52-7
2	—	—	—	—	98-88-4
3	—	—	—	—	7200-04-6
4	—	—	—	—	582-24-1
5	—	—	—	—	532-55-8
6	—	—	—	—	98-91-9
7	—	—	—	—	611-73-4
8	—	—	—	—	643-79-8
9	—	—	—	—	15206-55-0

MID LIBRARY SEARCH <LIBRARYNB>
 08/31/98 21:00:00 + 3:10
 SAMPLE: T-MM5-4 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
 ENHANCED <S 15B 2N 0T>
 DATA: 30068105 # 285
 CALI: 30068105 # 3
 BASE M/Z: 77
 RIC: 28832.



Library Search Data: 30068105 # 761 Base m/z: 57
 08/31/98 21:00:00 + 8:28 Cali: 30068105 # 3 RIC: 22976.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 611 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 22530 TETRADECANE
 2 19006 DECANE, 6-ETHYL-2-METHYL-
 3 11607 UNDECANE
 4 19007 DODECANE, 3-METHYL-
 5 19016 UNDECANE, 4,7-DIMETHYL-
 6 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 7 15358 UNDECANE, 3-METHYL-
 8 11612 NONANE, 2,5-DIMETHYL-
 9 25997 PENTADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C14.H30	198	43	762	960	773
2	C13.H28	184	57	755	975	755
3	C11.H24	156	43	737	980	737
4	C13.H28	184	57	734	927	738
5	C13.H28	184	43	727	938	727
6	C21.H44	296	57	727	927	760
7	C12.H26	170	57	724	934	728
8	C11.H24	156	57	724	887	724
9	C15.H32	212	57	722	932	753

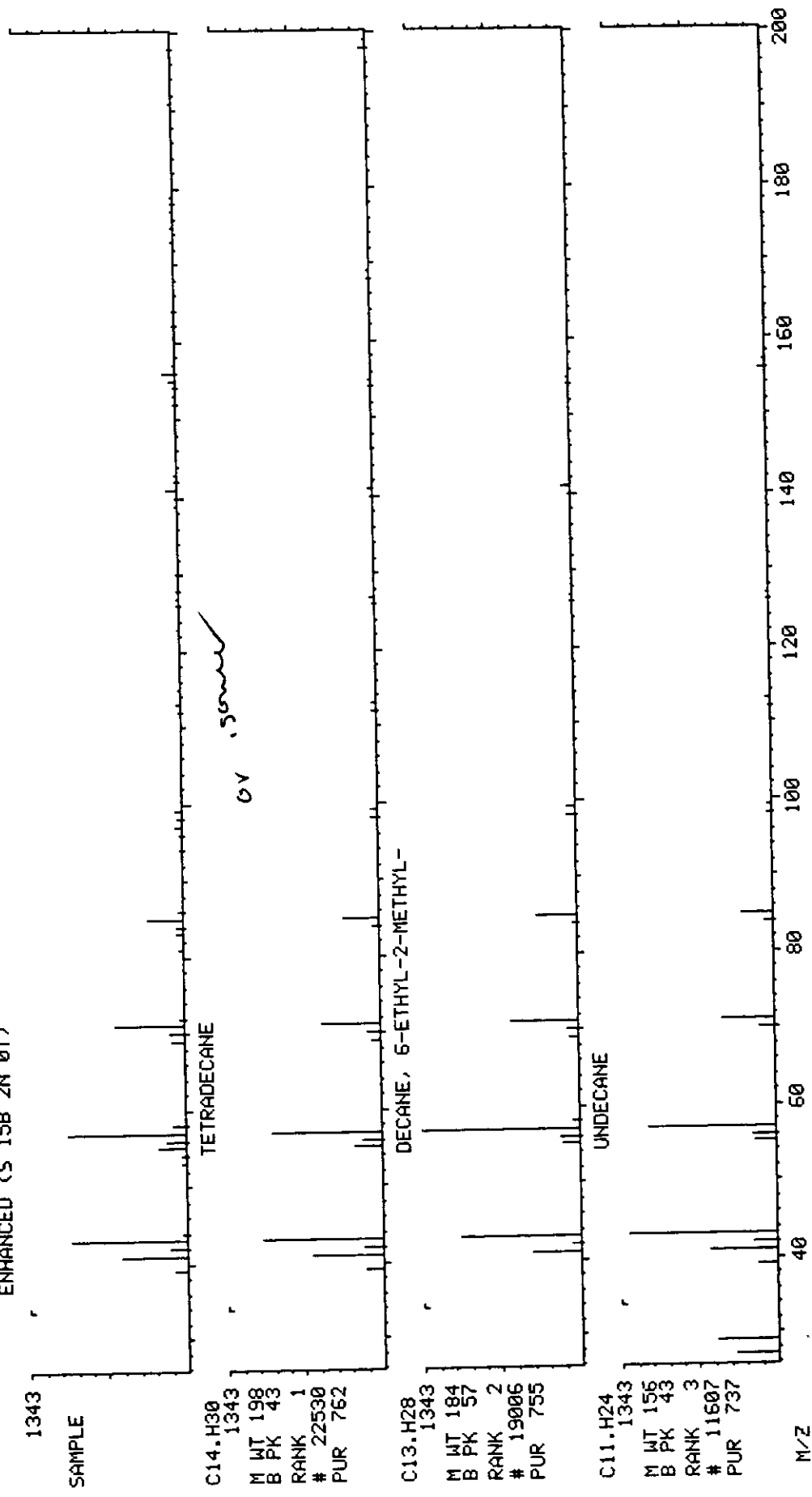
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	629-59-4
2	---	---	---	---	62108-21-8
3	---	---	---	---	1120-21-4
4	---	---	---	---	17312-57-1
5	---	---	---	---	17301-32-5
6	---	---	---	---	18344-37-1
7	---	---	---	---	1002-43-3
8	---	---	---	---	17302-27-1
9	---	---	---	---	629-62-9

DATA: 30058105 # 761
 CALI: 30058105 # 3

BASE M/Z: 57
 RIC: 22976.

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 21:00:00 + 8:28
 SAMPLE: T-MM5-4 1/35A/1ML INST. ID: F15
 CONDS.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)



Library Search Data: 30068105 # 856 Base m/z: 57
 08/31/98 21:00:00 + 9:31 Cali: 30068105 # 3 RIC: 35840.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 792 matched at least 7 of the 16 largest peaks in the unknown

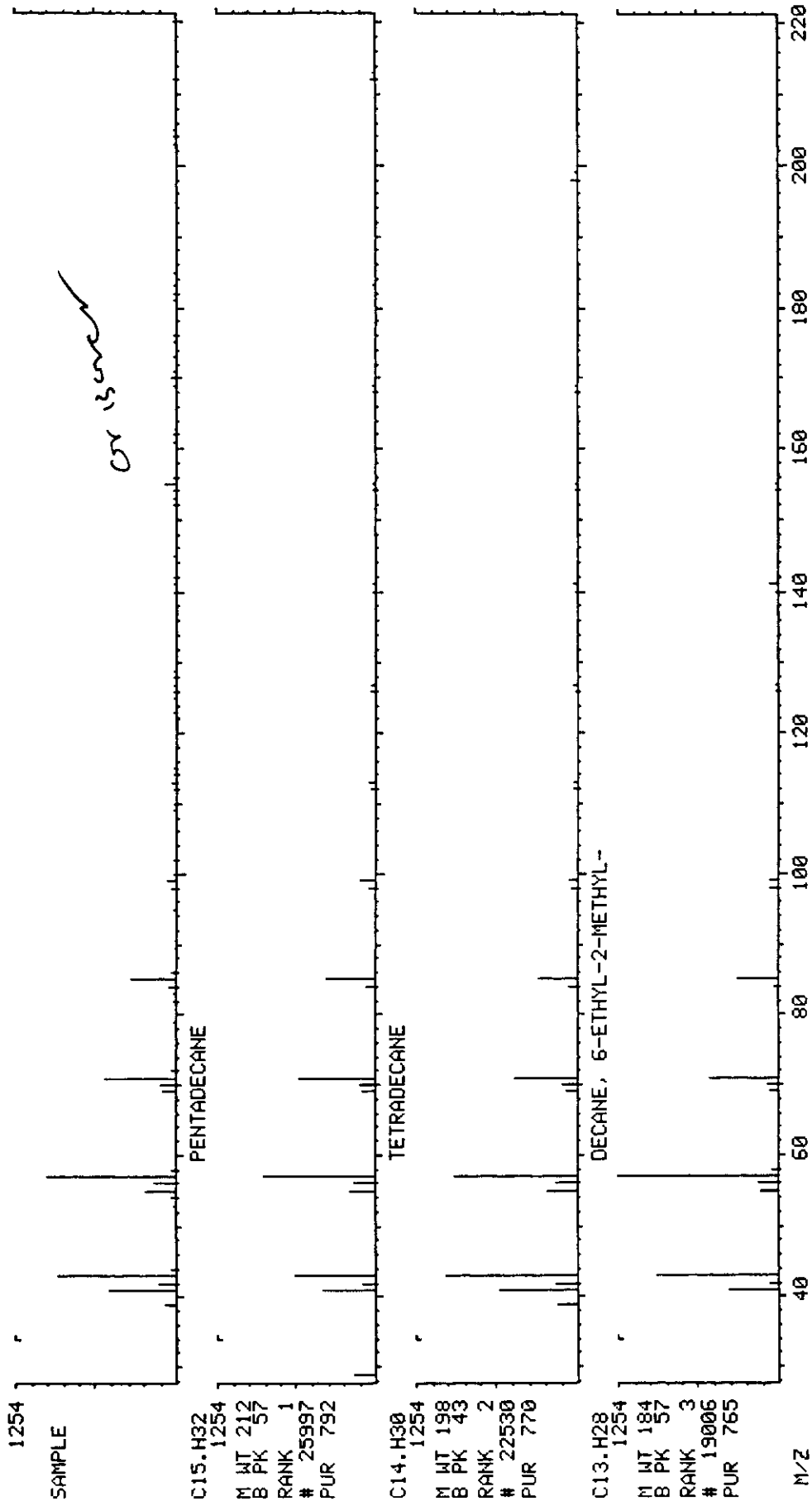
Rank In. Name
 1 25997 PENTADECANE
 2 22530 TETRADECANE
 3 19006 DECANE, 6-ETHYL-2-METHYL-
 4 15343 DODECANE
 5 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 6 19007 DODECANE, 3-METHYL-
 7 39681 1-iodoundecane
 8 15352 UNDECANE, 2-METHYL-
 9 18990 UNDECANE, 2,9-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C15.H32	212	57	792	965	798
2	C14.H30	198	43	770	939	804
3	C13.H28	184	57	765	966	765
4	C12.H26	170	57	764	989	764
5	C21.H44	296	57	759	930	783
6	C13.H28	184	57	755	954	763
7	C11.H23.1	282	57	755	971	760
8	C12.H26	170	43	754	939	762
9	C13.H28	184	57	754	957	756

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-62-9
2	—	—	—	—	629-59-4
3	—	—	—	—	62108-21-8
4	—	—	—	—	112-40-3
5	—	—	—	—	18344-37-1
6	—	—	—	—	17312-57-1
7	—	—	—	—	4282-44-4
8	—	—	—	—	7045-71-8
9	—	—	—	—	17301-26-7

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:00:00 + 9:31
 SAMPLE: T-MM5-4 1/35A/1ML
 CONDS.: UG/ML *1002/1002 *(NA/NA >1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068105 # 856
 CALI: 30068105 # 3
 BASE M/Z: 57
 RIC: 35840.



Library Search Data: 30068105 # 946 Base m/z: 57
 08/31/98 21:00:00 + 10:31 Cali: 30068105 # 3 RIC: 52608.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

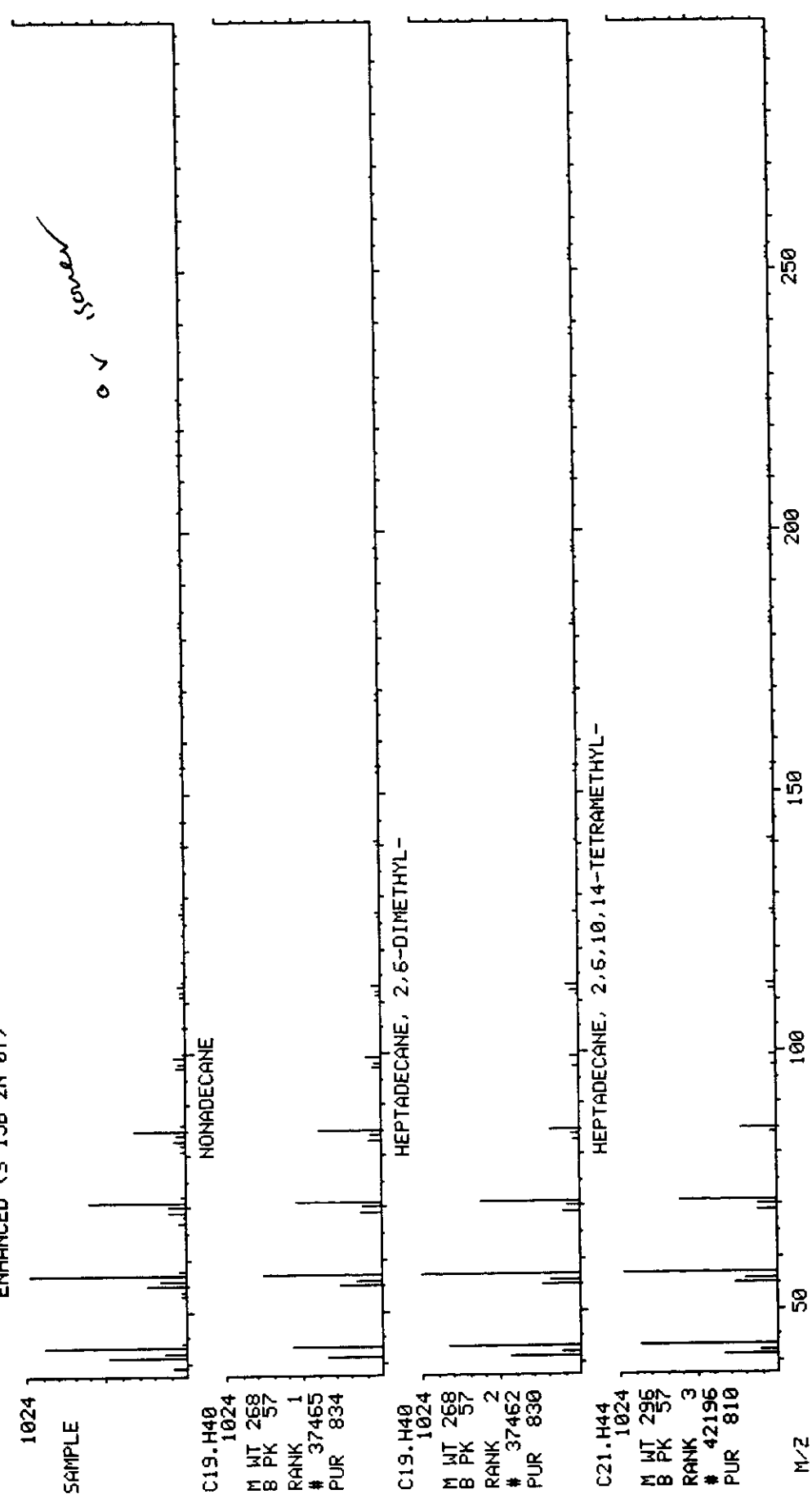
62231 spectra in LIBRARYNB searched for maximum PURITY
 360 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 25997 PENTADECANE
 5 15969 HYDROXYLAMINE, O-DECYL-
 6 25994 DODECANE, 2,6,11-TRIMETHYL-
 7 26001 DODECANE, 2,7,10-TRIMETHYL-
 8 19016 UNDECANE, 4,7-DIMETHYL-
 9 22535 DODECANE, 4,6-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	834	963	848
2	C19.H40	268	57	830	954	844
3	C21.H44	296	57	810	955	827
4	C15.H32	212	57	801	957	819
5	C10.H23.O.N	173	43	797	962	799
6	C15.H32	212	57	794	971	795
7	C15.H32	212	57	791	972	792
8	C13.H28	184	43	790	965	791
9	C14.H30	198	57	789	956	793

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	18344-37-1
4	—	—	—	—	629-62-9
5	—	—	—	—	29812-79-1
6	—	—	—	—	31295-56-4
7	—	—	—	—	74645-98-0
8	—	—	—	—	17301-32-5
9	—	—	—	—	61141-72-8

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:00:00 + 10:31
 SAMPLE: T-MM5-4 1/35A/1ML
 COND.: UG/ML *1ML *1002/1002 *(NA/NA) >1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068105 # 946
 CALI: 30068105 # 3
 BASE M/Z: 57
 RIC: 52608.



Library Search Data: 30068105 # 986 Base m/z: 57
 08/31/98 21:00:00 + 10:58 Cali: 30068105 # 3 RIC: 33280.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYN8 searched for maximum PURITY
 644 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 37465 NONADECANE
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 19054 UNDECANE, 2,6-DIMETHYL-
 5 22530 TETRADECANE
 6 25997 PENTADECANE
 7 19026 DECANE, 2,6,8-TRIMETHYL-
 8 26001 DODECANE, 2,7,10-TRIMETHYL-
 9 22535 DODECANE, 4,6-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	814	957	825
2	C19.H40	268	57	790	933	820
3	C21.H44	296	57	788	946	806
4	C13.H28	184	57	787	947	798
5	C14.H30	198	43	781	941	793
6	C15.H32	212	57	773	939	805
7	C13.H28	184	57	771	953	776
8	C15.H32	212	57	770	970	772
9	C14.H30	198	57	766	953	769

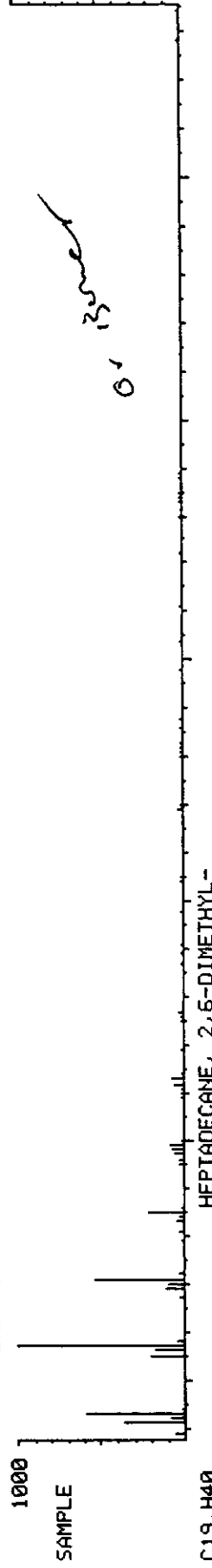
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	629-92-5
3	—	—	—	—	18344-37-1
4	—	—	—	—	17301-23-4
5	—	—	—	—	629-59-4
6	—	—	—	—	629-62-9
7	—	—	—	—	62108-26-3
8	—	—	—	—	74645-98-0
9	—	—	—	—	61141-72-8

DATA: 30068105 # 986
CALI: 30068105 # 3

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 21:00:00 + 10:58

SAMPLE: T-MM5-4 1/35A/1ML
COND.: UG/ML *1ML *100Z/100Z * (NA/NA) >1/35A NA M

INST. ID: F16
ENHANCED (S 158 2N 0T)



Library Search Data: 30068105 #1032 Base m/z: 57
 08/31/98 21:00:00 + 11:29 Cali: 30068105 # 3 RIC: 62400.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 360 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 3 37465 NONADECANE
 4 32059 HEPTADECANE
 5 46161 TRICOSANE
 6 25991 DODECANE, 2,6,10-TRIMETHYL-
 7 26001 DODECANE, 2,7,10-TRIMETHYL-
 8 25997 PENTADECANE
 9 22535 DODECANE, 4,6-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	850	972	861
2	C21.H44	296	57	822	963	838
3	C19.H40	268	57	816	948	849
4	C17.H36	240	57	788	925	821
5	C23.H48	324	43	785	901	854
6	C15.H32	212	57	783	958	795
7	C15.H32	212	57	781	959	795
8	C15.H32	212	57	778	938	819
9	C14.H30	198	57	778	954	788

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	18344-37-1
3	—	—	—	—	629-92-5
4	—	—	—	—	629-78-7
5	—	—	—	—	638-67-5
6	—	—	—	—	3891-98-3
7	—	—	—	—	74645-98-0
8	—	—	—	—	629-62-9
9	—	—	—	—	61141-72-8

DATA: 30068105 #1032
CALI: 30068105 # 3

BASE M/Z: 57
RIC: 62400.

MID LIBRARY SEARCH (LIBRARYNB)

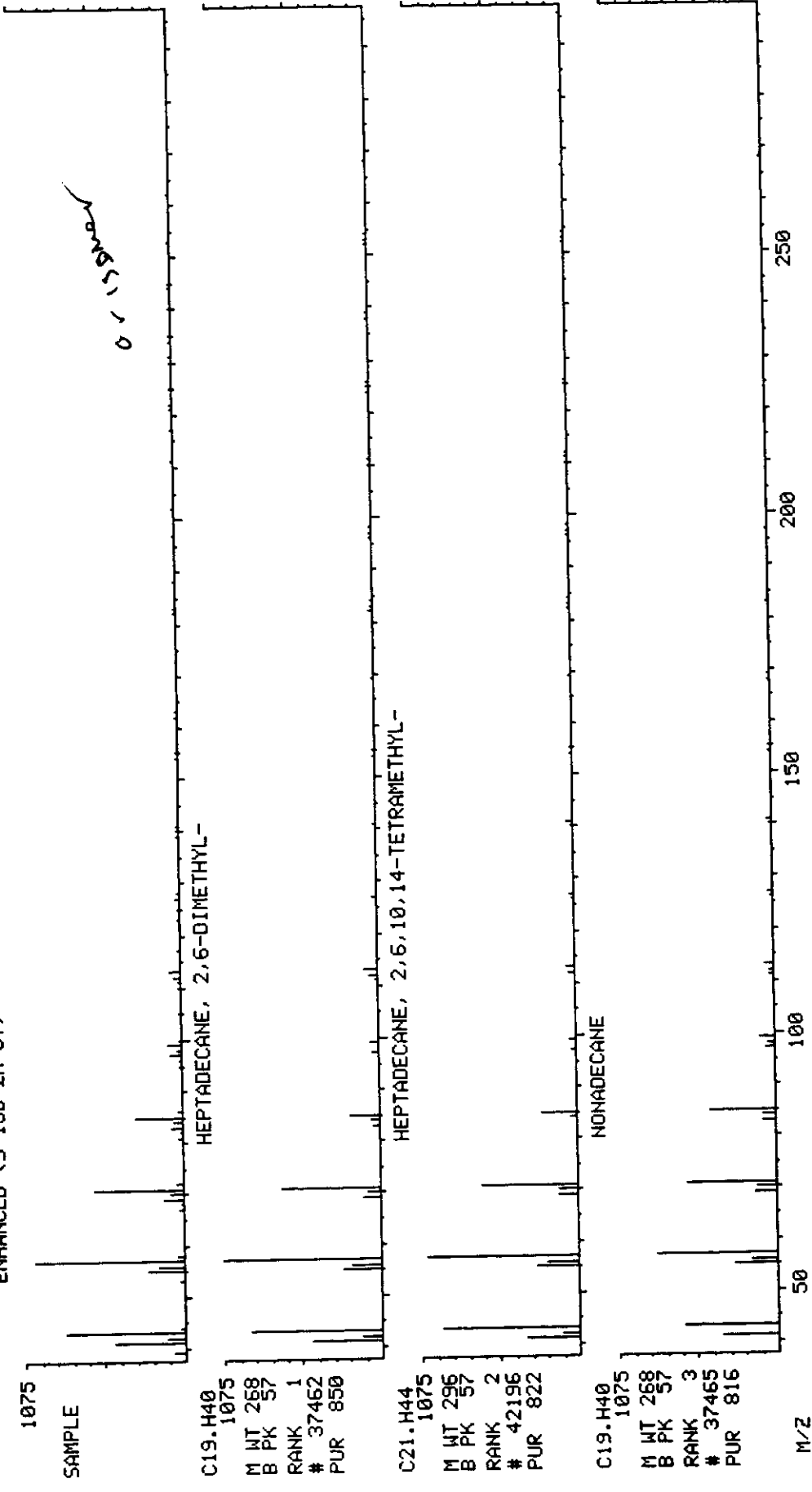
08/31/98 21:00:00 + 11:29

SAMPLE: T-MM5-4 1/35A/1ML

COND.: UG/ML *1ML *100%/100% *(NA/NA >1/35A NA M

ENHANCED (S 158 2N 0T)

INST. ID: F16



Library Search Data: 30068105 #1113 Base m/z: 57
 08/31/98 21:00:00 + 12:23 Cali: 30068105 # 3 RIC: 34816.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 312 matched at least 8 of the 16 largest peaks in the unknown

Rank In.	Name
1	37465 NONADECANE
2	42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
3	25997 PENTADECANE
4	19016 UNDECANE, 4,7-DIMETHYL-
5	37462 HEPTADECANE, 2,6-DIMETHYL-
6	15969 HYDROXYLAMINE, O-DECYL-
7	15352 UNDECANE, 2-METHYL-
8	18998 UNDECANE, 3,7-DIMETHYL-
9	46161 TRICOSANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	785	939	829
2	C21.H44	296	57	783	940	811
3	C15.H32	212	57	777	951	809
4	C13.H28	184	43	777	958	777
5	C19.H40	268	57	775	924	815
6	C10.H23.O.N	173	43	775	958	785
7	C12.H26	170	43	767	955	773
8	C13.H28	184	43	766	954	770
9	C23.H48	324	43	759	893	844

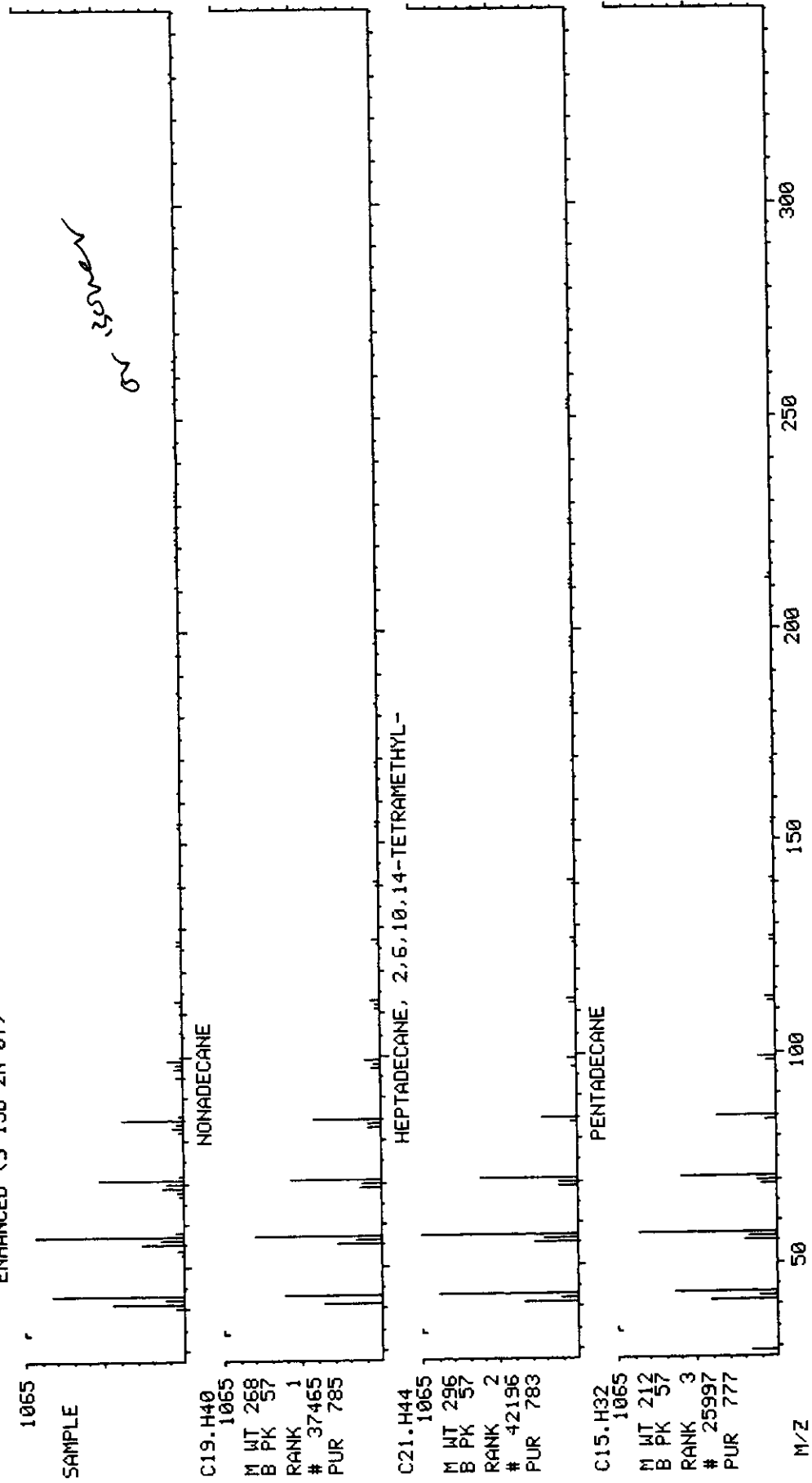
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	18344-37-1
3	—	—	—	—	629-62-9
4	—	—	—	—	17301-32-5
5	—	—	—	—	54105-67-8
6	—	—	—	—	29812-79-1
7	—	—	—	—	7045-71-8
8	—	—	—	—	17301-29-0
9	—	—	—	—	638-67-5

DATA: 30068105 #1113
CALI: 30068105 # 3

BASE M/Z: 57
RIC: 34816.

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 21:00:00 + 12:23
SAMPLE: T-HMS-4 1/35A/1ML
CONDS.: UG/ML *1ML *1002/1002 *(NA/NA) >1/35A NA M
ENHANCED (S 158 2N 0T)

INST. ID: F16



Library Search Data: 30068105 #1117 Base m/z: 57
 08/31/98 21:00:00 + 12:25 Cali: 30068105 # 3 RIC: 29952.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 774 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 25991 DODECANE, 2,6,10-TRIMETHYL-
 5 26001 DODECANE, 2,7,10-TRIMETHYL-
 6 37466 PENTADECANE, 2,6,10,14-TETRAMETHYL-
 7 29264 TRIDECAE, 5-PROPYL-
 8 32059 HEPTADECANE
 9 46161 TRICOSANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	811	941	838
2	C19.H40	268	57	808	932	819
3	C21.H44	296	57	788	923	809
4	C15.H32	212	57	782	958	784
5	C15.H32	212	57	780	959	780
6	C19.H40	268	71	779	924	798
7	C16.H34	226	57	775	928	777
8	C17.H36	240	57	770	909	813
9	C23.H48	324	43	770	888	835

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	18344-37-1
4	—	—	—	—	3891-98-3
5	—	—	—	—	74645-98-0
6	—	—	—	—	1921-70-6
7	—	—	—	—	55045-11-9
8	—	—	—	—	629-78-7
9	—	—	—	—	638-67-5

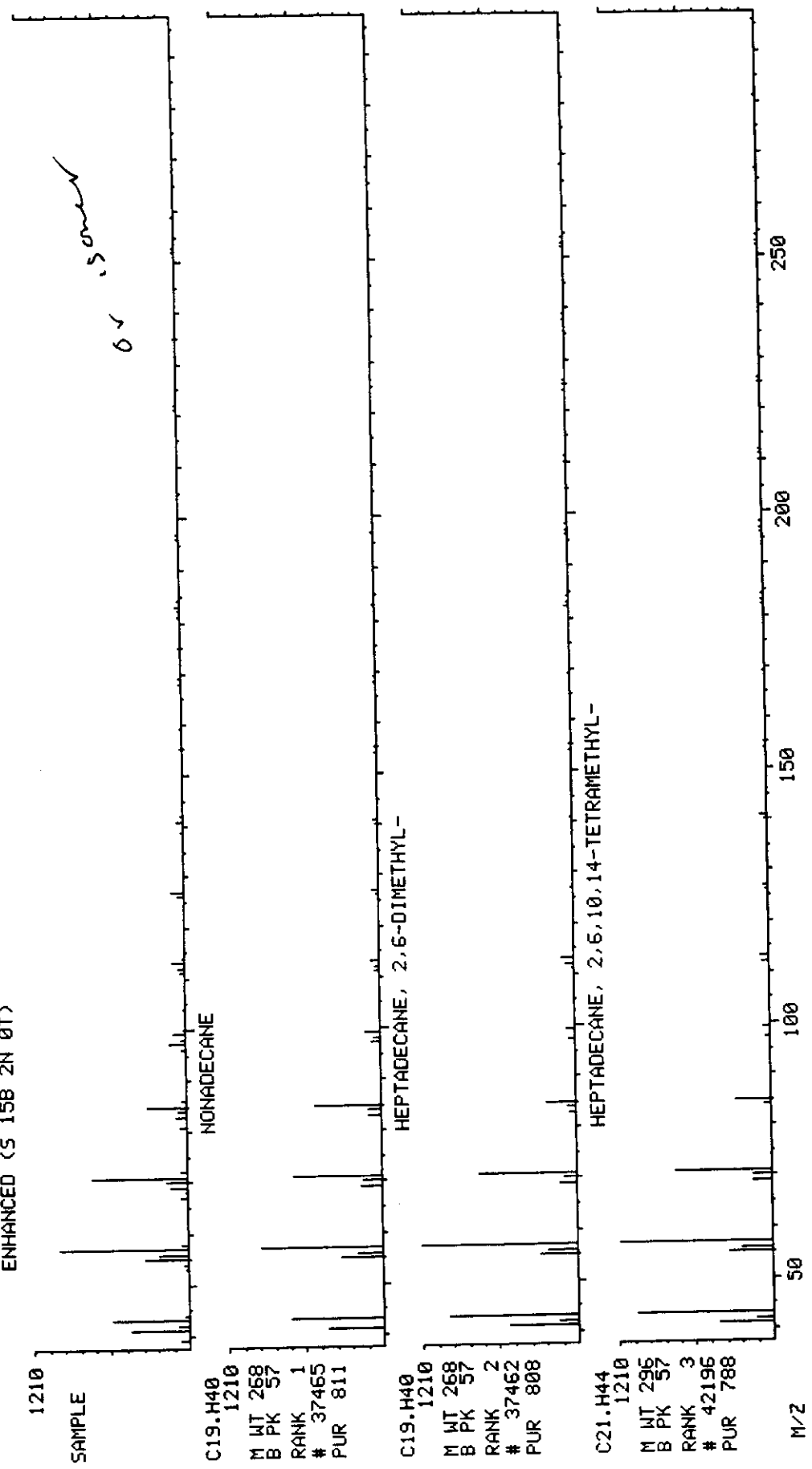
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DATA: 30068105 #1117
CALI: 30068105 # 3

BASE M/Z: 57
RIC: 29952.

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 21:00:00 + 12:25
SAMPLE: T-MM5-4 1/35A/1ML
CONDS.: UG/ML *1ML *1002/1002 *(NA/NA)/1/35A NA M
ENHANCED (S 158 2N 0T)

INST. ID: F16



Library Search Data: 30068105 #1190 Base m/z: 57
 08/31/98 21:00:00 + 13:14 Cali: 30068105 # 3 RIC: 26400.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N QT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 294 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 46161 TRICOSANE
 4 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 5 15969 HYDROXYLAMINE, O-DECYL-
 6 25997 PENTADECANE
 7 15353 2,6-DIMETHYLDECANE
 8 19016 UNDECANE, 4,7-DIMETHYL-
 9 32059 HEPTADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	746	971	757
2	C19.H40	268	57	721	941	743
3	C23.H48	324	43	701	902	770
4	C21.H44	296	57	700	946	722
5	C10.H23.O.N	173	43	699	961	702
6	C15.H32	212	57	698	954	722
7	C12.H26	170	43	689	932	689
8	C13.H28	184	43	684	950	684
9	C17.H36	240	57	676	914	714

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	638-67-5
4	—	—	—	—	18344-37-1
5	—	—	—	—	29812-79-1
6	—	—	—	—	629-62-9
7	—	—	—	—	13150-81-7
8	—	—	—	—	17301-32-5
9	—	—	—	—	629-78-7

BASE M/Z: 57
RIC: 26400.

DATA: 30068105 #1190
CALI: 30068105 # 3

MID LIBRARY SEARCH (LIBRARYNB)

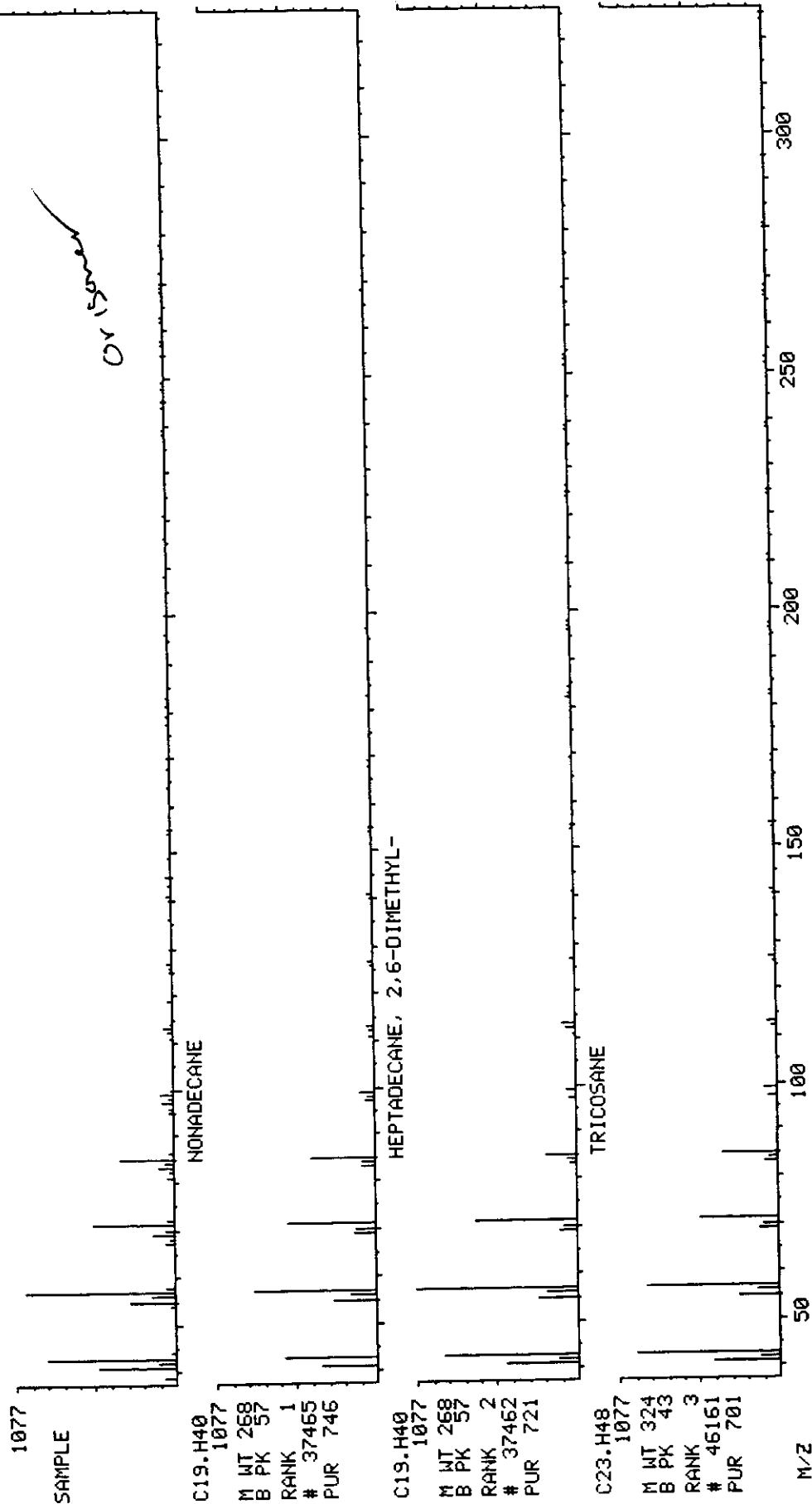
08/31/98 21:00:00 + 13:14

SAMPLE: T-MM5-4 1/35A/1ML

COND5.: UG/ML *1ML *100Z/100Z *KNA/NA >1/35A NA M

ENHANCED (S 158 2N 0T)

INST. ID: F16



Library Search Data: 30068105 #1236 Base m/z: 73
 08/31/98 21:00:00 + 13:45 Cali: 30068105 # 3 RIC: 23072.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 747 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 35182 HEXADECANOIC ACID
 2 37913 GLYCINE, N-METHYL-N-(1-OXODODECYL)-
 3 40669 DODECANAMIDE, N,N-BIS(2-HYDROXYETHYL)-
 4 22949 DODECANOIC ACID
 5 26351 TRIDECANOIC ACID
 6 29642 TETRADECANOIC ACID
 7 40184 OCTADECANOIC ACID
 8 19469 UNDECANOIC ACID
 9 15784 DECANOIC ACID

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C16.H32.O2	256	43	682	975	688
2	C15.H29.O3.N	271	73	588	974	588
3	C16.H33.O3.N	287	73	567	959	573
4	C12.H24.O2	200	60	542	916	566
5	C13.H26.O2	214	73	542	887	558
6	C14.H28.O2	228	73	532	871	589
7	C18.H36.O2	284	43	514	753	665
8	C11.H22.O2	186	60	506	912	531
9	C10.H20.O2	172	60	504	939	506

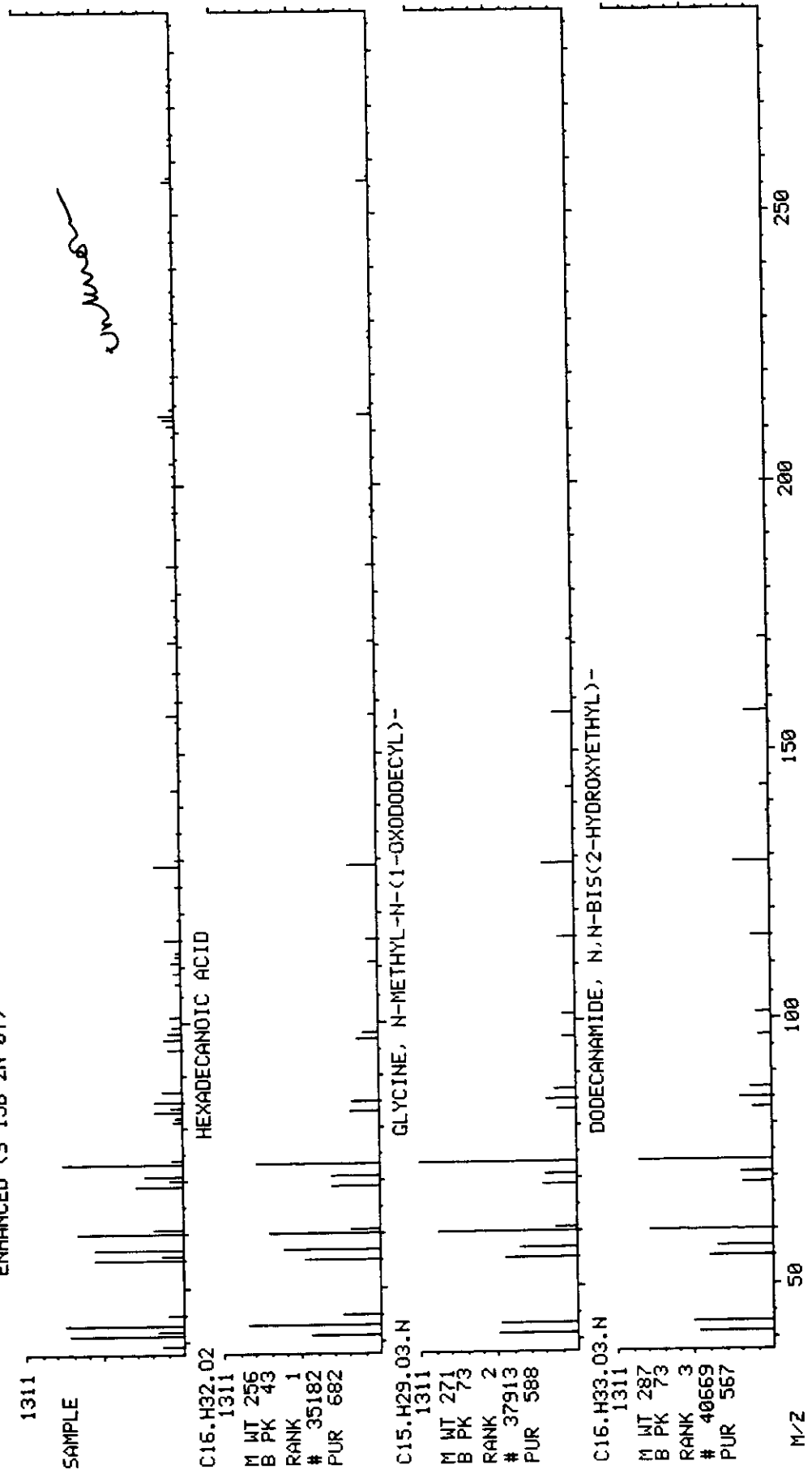
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	57-10-3
2	—	—	—	—	97-78-9
3	—	—	—	—	120-40-1
4	—	—	—	—	143-07-7
5	—	—	—	—	638-53-9
6	—	—	—	—	544-63-8
7	—	—	—	—	57-11-4
8	—	—	—	—	112-37-8
9	—	—	—	—	334-48-5

DATA: 30068105 #1235
CALI: 30068105 # 3

BASE M/Z: 73
RIC: 23072.

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 21:00:00 + 13:45
SAMPLE: T-MMS-4 1/35A/1ML
COND.: UG/ML *1ML *100%/100% *(NA/NA) >1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068105 #1334 Base m/z: 57
 08/31/98 21:00:00 + 14:50 Cali: 30068105 # 3 RIC: 18240.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (\$ 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 439 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 37456 2-METHYLOCTADECANE
 3 37465 NONADECANE
 4 19523 1-DECANOL, 2-ETHYL-
 5 14793 1-UNDECENE, 4-METHYL-
 6 37458 OCTADECANE, 2-METHYL-
 7 29263 HEXADECANE
 8 46161 TRICOSANE
 9 42481 OXIRANE, [(HEXADECYLOXY)METHYL]-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	596	881	632
2	C19.H40	268	43	593	901	640
3	C19.H40	268	57	585	919	628
4	C12.H26.O	186	57	582	931	593
5	C12.H24	168	43	581	957	581
6	C19.H40	268	43	577	864	639
7	C16.H34	226	57	577	899	621
8	C23.H48	324	43	577	863	659
9	C19.H38.O2	298	57	576	872	583

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	- -
3	—	—	—	—	629-92-5
4	—	—	—	—	21078-65-9
5	—	—	—	—	74630-39-0
6	—	—	—	—	1560-88-9
7	—	—	—	—	544-76-3
8	—	—	—	—	638-67-5
9	—	—	—	—	15965-99-8

DATA: 30068105 #1334
CALI: 30068105 # 3

BASE M/Z: 57
RIC: 18240.

MID LIBRARY SEARCH (LIBRARYNB)

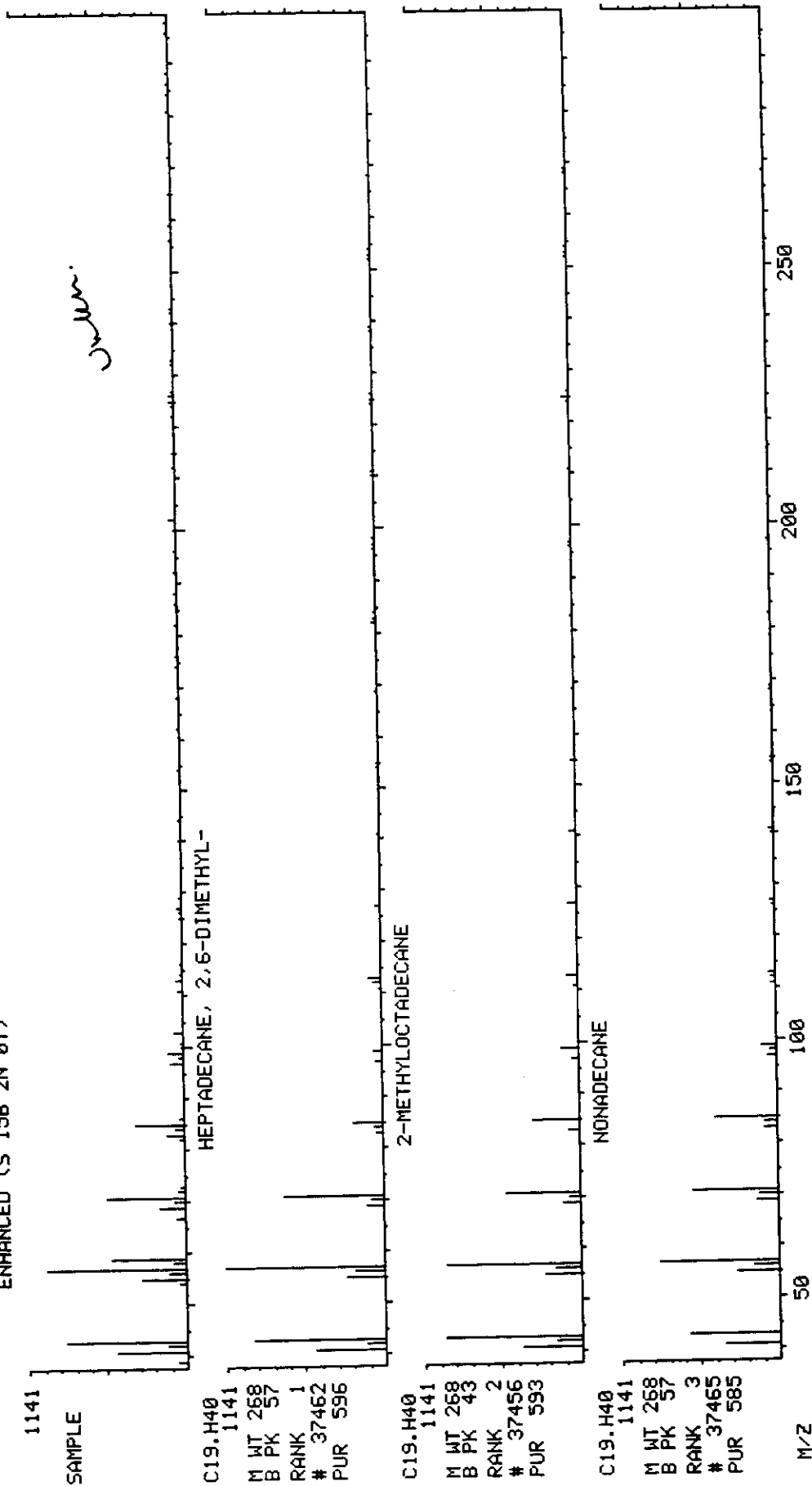
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SAMPLE: T-MM5-4 1/35A/1ML

COND5.: UG/ML *1ML *100%/100% *(NA/NA) /1/35A NA M

ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068105 #1360 Base m/z: 57
 08/31/98 21:00:00 + 15:08 Cali: 30068105 # 3 RIC: 3440.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 144 matched at least 5 of the 16 largest peaks in the unknown

Rank In. Name
 1 16268 PROPANE, 1,1'-[ETHYLIDENE BIS(OXY)] BIS[2-METHYL-
 2 3365 2-PENTANOL, 4,4-DIMETHYL-
 3 13863 PHOSPHINOUS CHLORIDE, TERT-BUTYLISOPROPYL-
 4 7691 BUTANE, 2-AZIDO-2,3,3-TRIMETHYL-
 5 38965 PHENOL, 2,6-BIS(1,1-DIMETHYLETHYL)-4-METHYL-, METHYL CARBAMATE
 6 3052 OXIRANE, (2,2-DIMETHYLPROPYL)-
 7 5791 1,3-DIOXOLANE-2-METHANOL, 2,4-DIMETHYL-
 8 5551 DI-SEC-BUTYL ETHER
 9 29529 THIOPHENE, 2,5-BIS(1,1-DIMETHYLETHOXY)-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C10.H22.O2	174	57	270	725	292
2	C7.H16.O	116	57	252	722	295
3	C7.H16.CL.P	166	57	229	707	243
4	C7.H15.N3	141	57	228	753	264
5	C17.H27.O2.N	277	205	221	521	308
6	C7.H14.O	114	57	214	741	274
7	C6.H12.O3	132	43	213	697	235
8	C8.H18.O	130	45	212	628	252
9	C12.H20.O2.S	228	41	207	700	282

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	5669-09-0
2	—	—	—	—	6144-93-0
3	—	—	—	—	29949-66-4
4	—	—	—	—	51677-41-9
5	—	—	—	—	1918-11-2
6	—	—	—	—	2245-29-6
7	—	—	—	—	53951-43-2
8	—	—	—	—	6863-58-7
9	—	—	—	—	55162-43-1

BASE M/Z: 57
RIC: 3440.

DATA: 30068105 #1360
CALI: 30068105 # 3

MID LIBRARY SEARCH (LIBRARYNB)

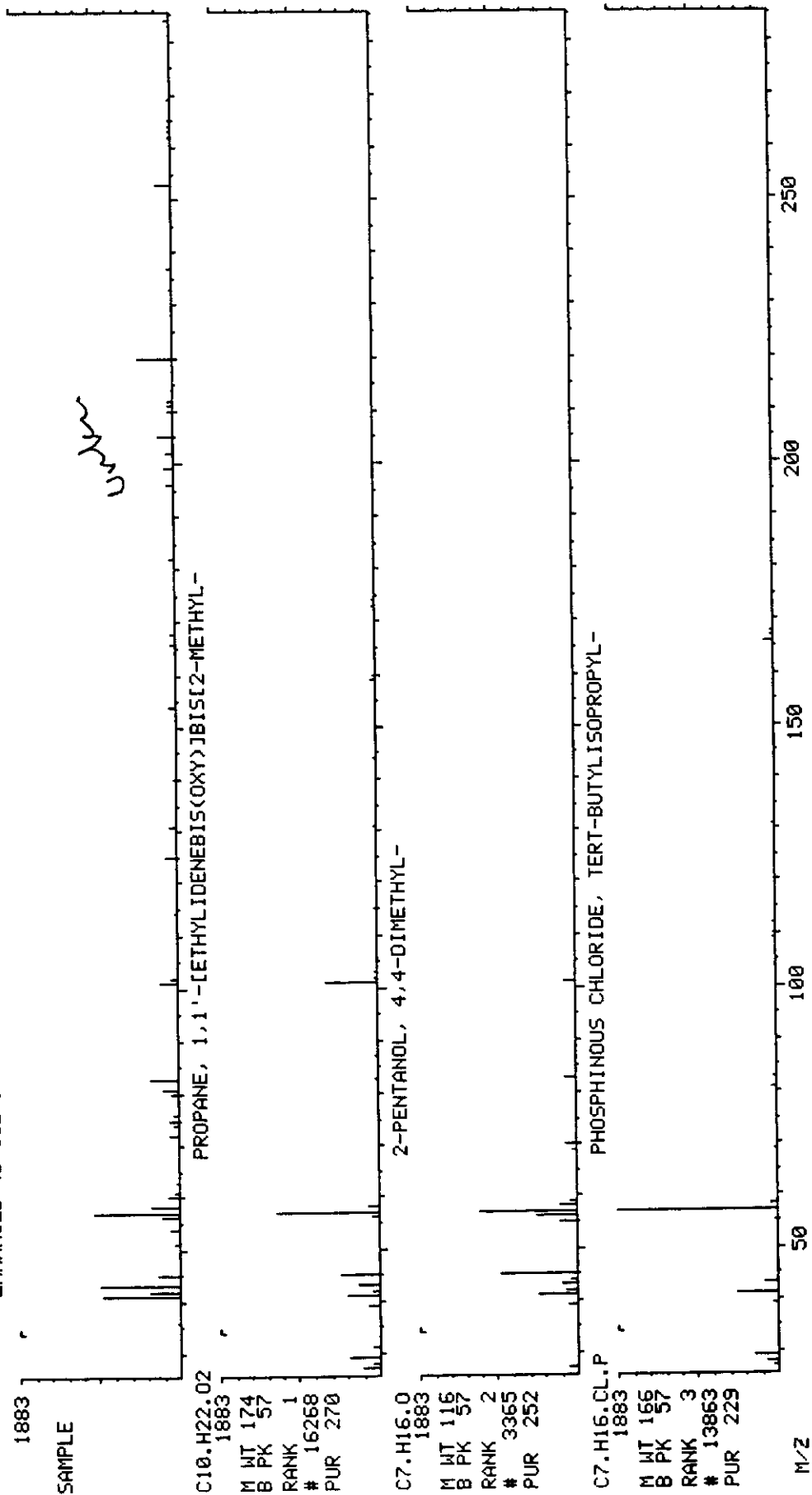
08/31/98 21:00:00 + 15:08

SAMPLE: T-MMS-4 1/35A/1ML

INST. ID: F16

CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA) >1/35A NA M

ENHANCED (S 158 2N 0T)



Library Search Data: 30068105 #1402 Base m/z: 57
 08/31/98 21:00:00 + 15:36 Cali: 30068105 # 3 RIC: 10224.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (\$ 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 488 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 46161 TRICOSANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 49555 PENTACOSANE
 4 37465 NONADECANE
 5 42197 HENEICOSANE
 6 58739 PENTATRIACONTANE
 7 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 8 42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-
 9 39858 EICOSANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C23.H48	324	43	654	890	721
2	C19.H40	268	57	653	895	688
3	C25.H52	352	43	651	870	734
4	C19.H40	268	57	635	914	684
5	C21.H44	296	57	633	879	689
6	C35.H72	492	57	627	882	698
7	C21.H44	296	57	625	877	671
8	C21.H44	296	57	612	885	661
9	C20.H42	282	57	612	852	706

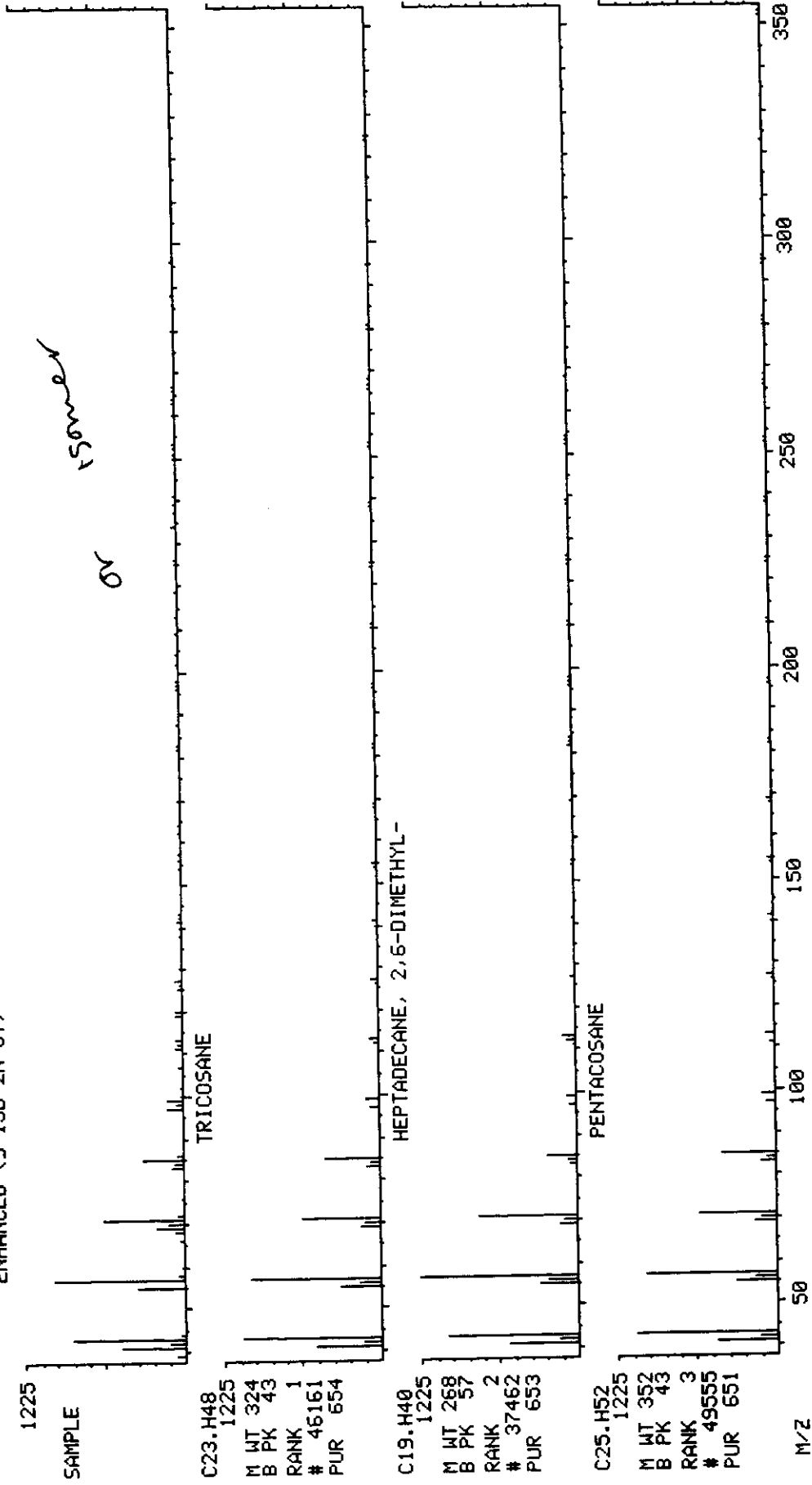
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1	—	—	—	—	638-67-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	629-99-2
4	—	—	—	—	629-92-5
5	—	—	—	—	629-94-7
6	—	—	—	—	630-07-9
7	—	—	—	—	18344-37-1
8	—	—	—	—	54833-48-6
9	—	—	—	—	112-95-8

DATA: 30068105 #1402
CALI: 30068105 # 3

BASE M/Z: 57
RIC: 10224.

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 21:00:00 + 15:36
SAMPLE: T-MM5-4 1/35A/1ML
COND.: UG/ML *1ML *100%/100% *(NA/NA) >1/35A NA M
ENHANCED (S 158 2N 0T)

INST. ID: F16



Library Search Data: 30068105 #1467 Base m/z: 57
 08/31/98 21:00:00 + 16:19 Cal: 30068105 # 3 RIC: 8352.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 751 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 46161 TRICOSANE
 2 37465 NONADECANE
 3 37462 HEPTADECANE, 2,6-DIMETHYL-
 4 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 5 49555 PENTACOSANE
 6 42190 EICOSANE, 2-METHYL-
 7 58739 PENTATRIACONTANE
 8 39858 EICOSANE
 9 25997 PENTADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C23.H48	324	43	588	891	644
2	C19.H40	268	57	572	941	600
3	C19.H40	268	57	569	900	588
4	C21.H44	296	57	567	898	587
5	C25.H52	352	43	565	862	642
6	C21.H44	296	57	549	861	608
7	C35.H72	492	57	544	864	614
8	C20.H42	282	57	538	845	622
9	C15.H32	212	57	538	949	560

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	638-67-5
2	—	—	—	—	629-92-5
3	—	—	—	—	54105-67-8
4	—	—	—	—	18344-37-1
5	—	—	—	—	629-99-2
6	—	—	—	—	1560-84-5
7	—	—	—	—	630-07-9
8	—	—	—	—	112-95-8
9	—	—	—	—	629-62-9

DATA: 30068105 #1467
CALI: 30068105 # 3

BASE M/Z: 57
RIC: 8352.

MID LIBRARY SEARCH (LIBRARYNB)

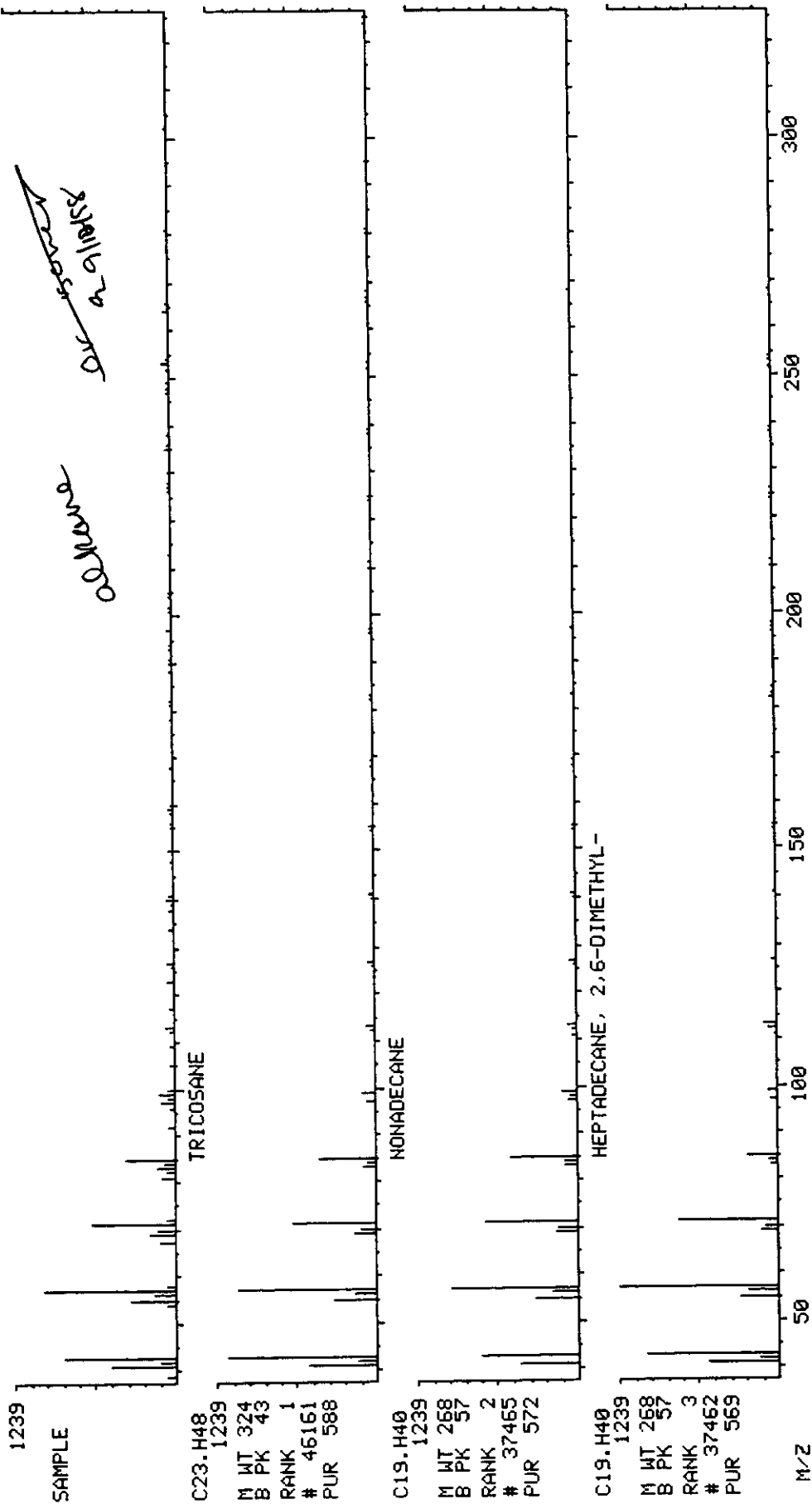
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SAMPLE: T-MM5-4 1/35A/1ML

CONDOS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M

ENHANCED (S 158 2N 0T)

INST. ID: F16



Library Search Data: 30068105 #1508 Base m/z: 59
 08/31/98 21:00:00 + 16:46 Cali: 30068105 # 3 RIC: 7352.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 179 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 24247 2-PROPANOL, 1-[2-(2-METHOXY-1-METHYLETHOXY)-1-METHYLETHOXY]-
 2 20265 3-HEXANOL, 1,5-DIMETHOXY-2,4-DIMETHYL-
 3 16192 2-PROPANOL, 1-[1-METHYL-2-(2-PROPENYLOXY)ETHOXY]-
 4 20626 2-PROPANOL, 1,1'-[(1-METHYL-1,2-ETHANEDIYL)BIS(OXY)]BIS-
 5 9228 1-PROPANOL, 2-(2-METHOXYPROPOXY)-
 6 3522 2,3-BUTANEDIOL, 2,3-DIMETHYL-
 7 8432 2-HEXANONE, 3-HYDROXY-3,5-DIMETHYL-
 8 9235 1-PROPANOL, 2-(2-METHOXY-1-METHYLETHOXY)-
 9 5832 2-PROPANOL, 1-(1-METHYLPROPOXY)-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C10.H22.O4	206	59	423	772	466
2	C10.H22.O3	190	59	423	782	466
3	C9.H18.O3	174	59	421	756	441
4	C9.H20.O4	192	59	404	794	427
5	C7.H16.O3	148	59	397	799	414
6	C6.H14.O2	118	59	384	901	391
7	C8.H16.O2	144	59	378	922	385
8	C7.H16.O3	148	59	378	740	388
9	C7.H16.O2	132	43	375	689	400

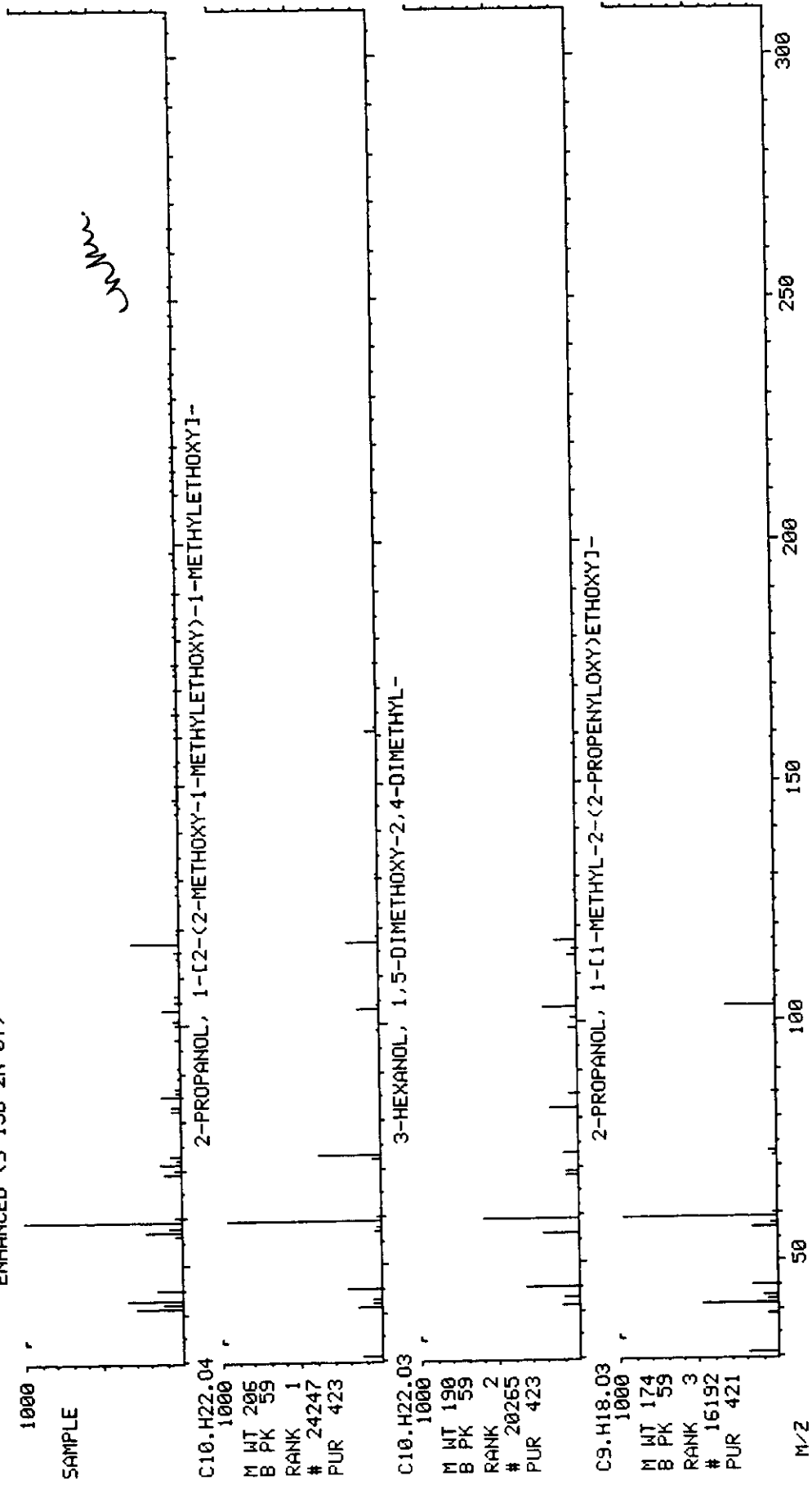
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2	---	---	---	---	13897-22-8
3	---	---	---	---	55956-25-7
4	---	---	---	---	1638-16-0
5	---	---	---	---	13588-28-8
6	---	---	---	---	76-09-5
7	---	---	---	---	6321-14-8
8	---	---	---	---	55956-21-3
9	---	---	---	---	53907-95-2

BASE M/Z: 59
RIC: 7352.

DATA: 30068105 #1508
CALL: 30068105 # 3

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 21:00:00 + 16:46
SAMPLE: T-MM5-4 1/35A/1ML
CONDOS.: UG/ML *1ML *100Z/100Z *(NA/NA) 1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



1000

Library Search Data: 30068105 #1529 Base m/z: 59
 08/31/98 21:00:00 + 17:00 Cali: 30068105 # 3 RIC: 8976.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 562 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 37465 NONADECANE
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 46161 TRICOSANE
 5 3000 OXIRANE, 2,2-DIMETHYL-3-PROPYL-
 6 19515 5,9-DIMETHYL-3-DECANOL
 7 25991 DODECANE, 2,6,10-TRIMETHYL-
 8 32059 HEPTADECANE
 9 25997 PENTADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	506	913	526
2	C19.H40	268	57	500	915	530
3	C21.H44	296	57	497	913	516
4	C23.H48	324	43	496	871	556
5	C7.H14.O	114	43	488	860	508
6	C12.H26.O	186	57	481	837	521
7	C15.H32	212	57	479	918	484
8	C17.H36	240	57	478	888	518
9	C15.H32	212	57	476	915	505

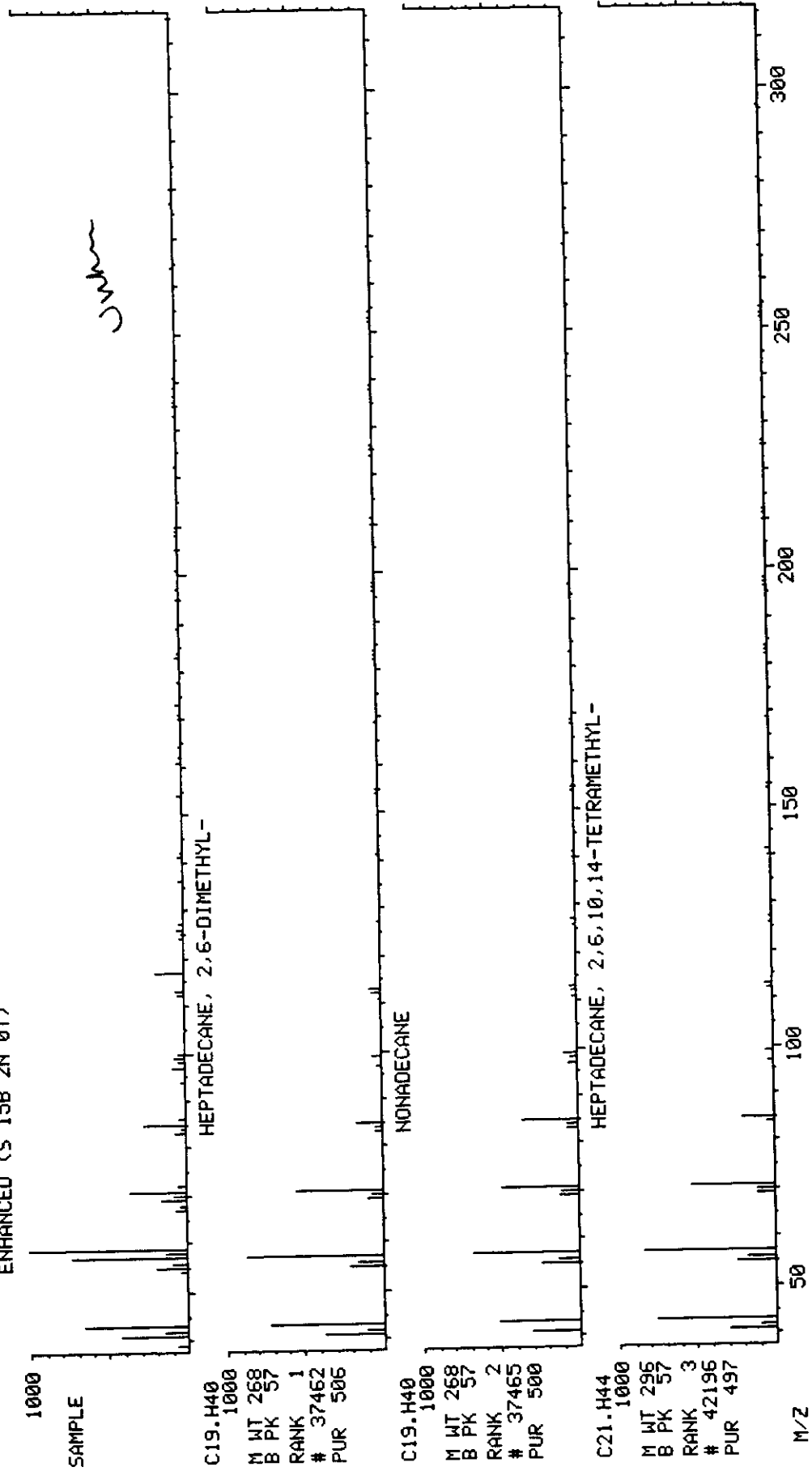
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	629-92-5
3	—	—	—	—	18344-37-1
4	—	—	—	—	638-67-5
5	—	—	—	—	17612-35-0
6	—	—	—	—	- -
7	—	—	—	—	3891-98-3
8	—	—	—	—	629-78-7
9	—	—	—	—	629-62-9

BASE M/Z: 59
RIC: 8976.

DATA: 30068105 #1529
CALI: 30068105 # 3

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 21:00:00 + 17:00
SAMPLE: T-MMS-4 1/35A/1ML
CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA) 1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068105 #1587 Base m/z: 43
 08/31/98 21:00:00 + 17:39 Cali: 30068105 # 3 RIC: 27744.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 965 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 39516 5-EICOSENE, (E)-
 2 39515 9-EICOSENE, (E)-
 3 34410 3-OCTADECENE, (E)-
 4 37062 1-NONADECENE
 5 28776 CYCLOHEXADECANE
 6 34418 9-OCTADECENE, (E)-
 7 34411 5-OCTADECENE, (E)-
 8 58701 17-PENTATRIACONTENE
 9 39517 3-EICOSENE, (E)-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C20.H40	280	55	771	974	774
2	C20.H40	280	55	768	974	771
3	C18.H36	252	69	766	957	778
4	C19.H38	266	97	763	965	771
5	C16.H32	224	55	762	970	762
6	C18.H36	252	55	759	952	771
7	C18.H36	252	55	759	952	771
8	C35.H70	490	43	759	948	793
9	C20.H40	280	57	757	977	760

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	74685-30-6
2	---	---	---	---	74685-29-3
3	---	---	---	---	7206-19-1
4	---	---	---	---	18435-45-5
5	---	---	---	---	295-65-8
6	---	---	---	---	7206-25-9
7	---	---	---	---	7206-21-5
8	---	---	---	---	6971-40-0
9	---	---	---	---	74685-33-9

DATA: 30068105 #1587
CALI: 30068105 # 3

BASE M/Z: 43
RIC: 27744.

MID LIBRARY SEARCH (LIBRARYNB)

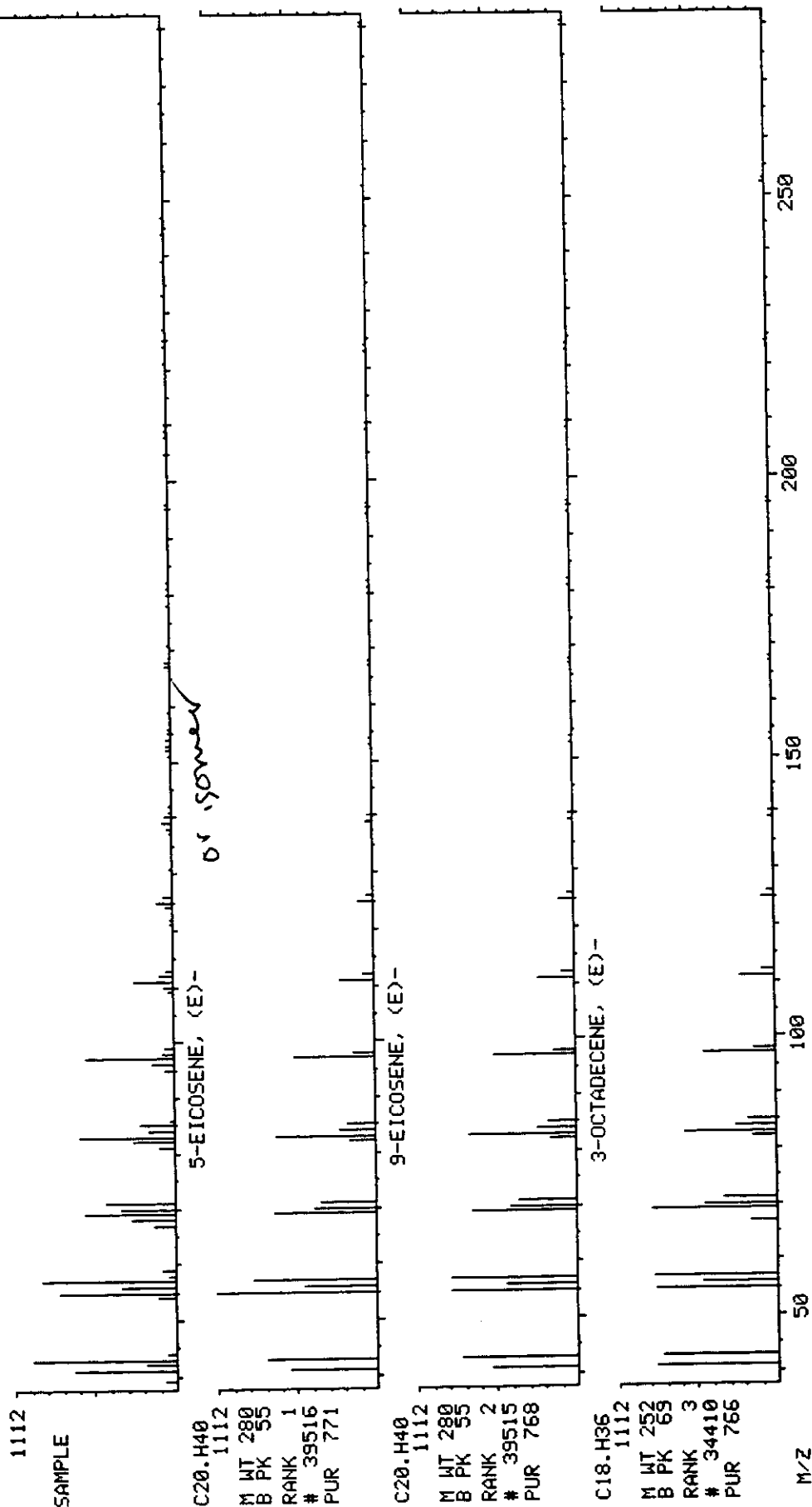
08/31/98 21:00:00 + 17:39

SAMPLE: T-MM5-4 1/35A/1ML

COND.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M

ENHANCED (S 158 2N 0T)

INST. ID: F16



Library Search Data: 30068105 #1662 Base m/z: 59
 08/31/98 21:00:00 + 18:29 Cali: 30068105 # 3 RIC: 5128.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYN8 searched for maximum PURITY
 707 matched at least 5 of the 16 largest peaks in the unknown

Rank	In.	Name
1	1909	2-PROPANOL, 1-METHOXY-2-METHYL-
2	20626	2-PROPANOL, 1,1'-[[1-METHYL-1,2-ETHANEDIYL]BIS(OXY)]BIS-
3	9235	1-PROPANOL, 2-(2-METHOXY-1-METHYLETHOXY)-
4	24247	2-PROPANOL, 1-[2-(2-METHOXY-1-METHYLETHOXY)-1-METHYLETHOXY]-
5	36537	2,5,8,11-TETRAOXATETRADECAN-13-OL, 4,7,10-TRIMETHYL-
6	6083	2-PROPANOL, 1,1'-OXYBIS-
7	9228	1-PROPANOL, 2-(2-METHOXYPROPOXY)-
8	16192	2-PROPANOL, 1-[1-METHYL-2-(2-PROPENYLOXY)ETHOXY]-
9	12856	2-PROPANOL, 1-(2-ETHOXYPROPOXY)-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C5.H12.O2	104	59	471	916	491
2	C9.H20.O4	192	59	453	812	475
3	C7.H16.O3	148	59	448	773	455
4	C10.H22.O4	206	59	447	757	513
5	C13.H28.O5	264	59	441	705	530
6	C6.H14.O3	134	59	440	886	447
7	C7.H16.O3	148	59	439	776	467
8	C9.H18.O3	174	59	438	746	466
9	C8.H18.O3	162	59	436	732	482

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	3587-64-2
2	—	—	—	—	1638-16-0
3	—	—	—	—	55956-21-3
4	—	—	—	—	20324-33-8
5	—	—	—	—	20324-34-9
6	—	—	—	—	110-98-5
7	—	—	—	—	13588-28-8
8	—	—	—	—	55956-25-7
9	—	—	—	—	10143-32-5

DATA: 30068105 #1804
CALI: 30068105 # 3

BASE M/Z: 59
RIC: 4592.

MID LIBRARY SEARCH <LIBRARYNB>

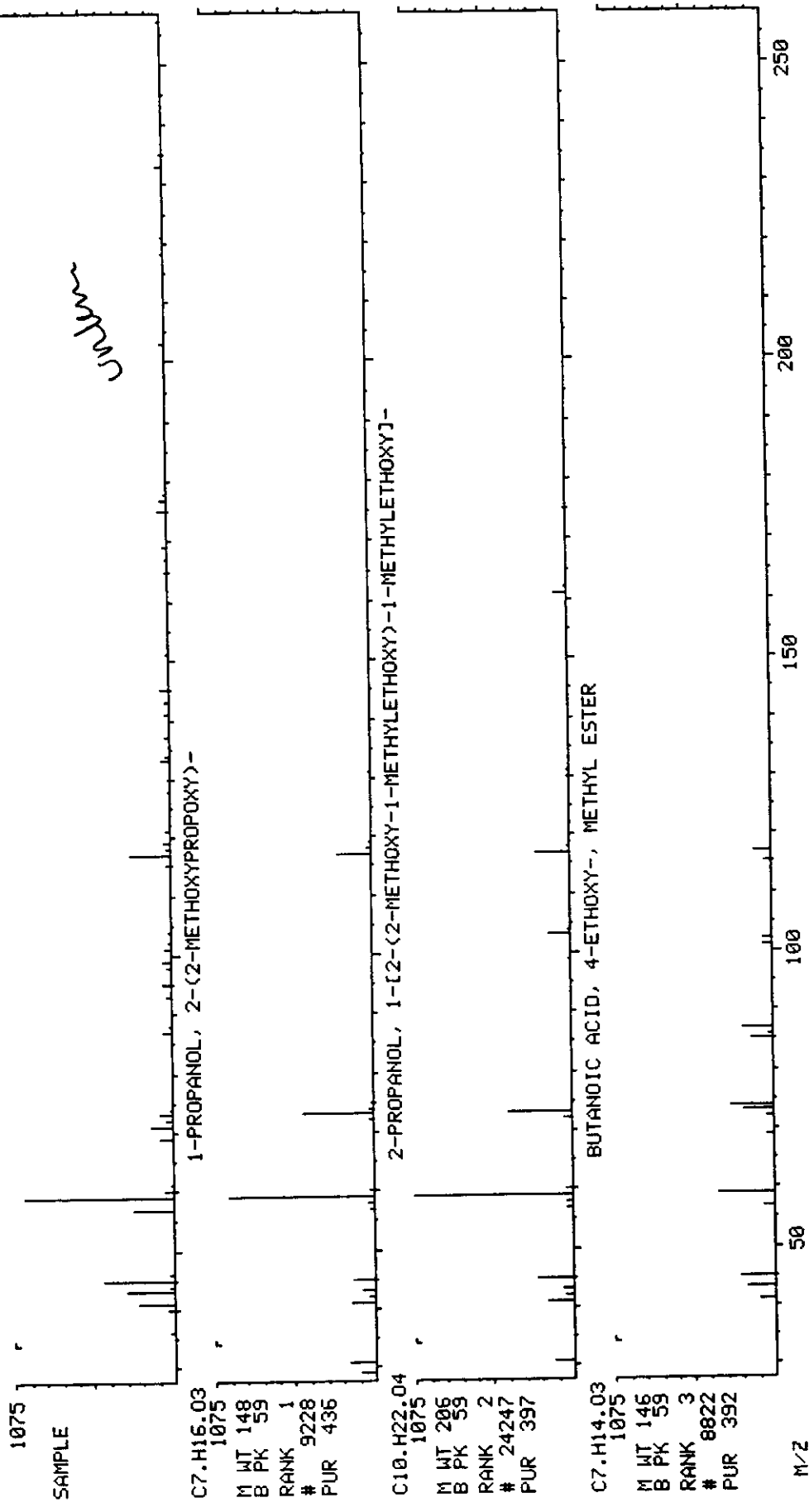
08/31/98 21:00:00 + 20:04

SAMPLE: T-MM5-4 1/35A/1ML

COND.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M

ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068105 #1916 Base m/z: 43
 08/31/98 21:00:00 + 21:19 Cal: 30068105 # 3 RIC: 8736.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

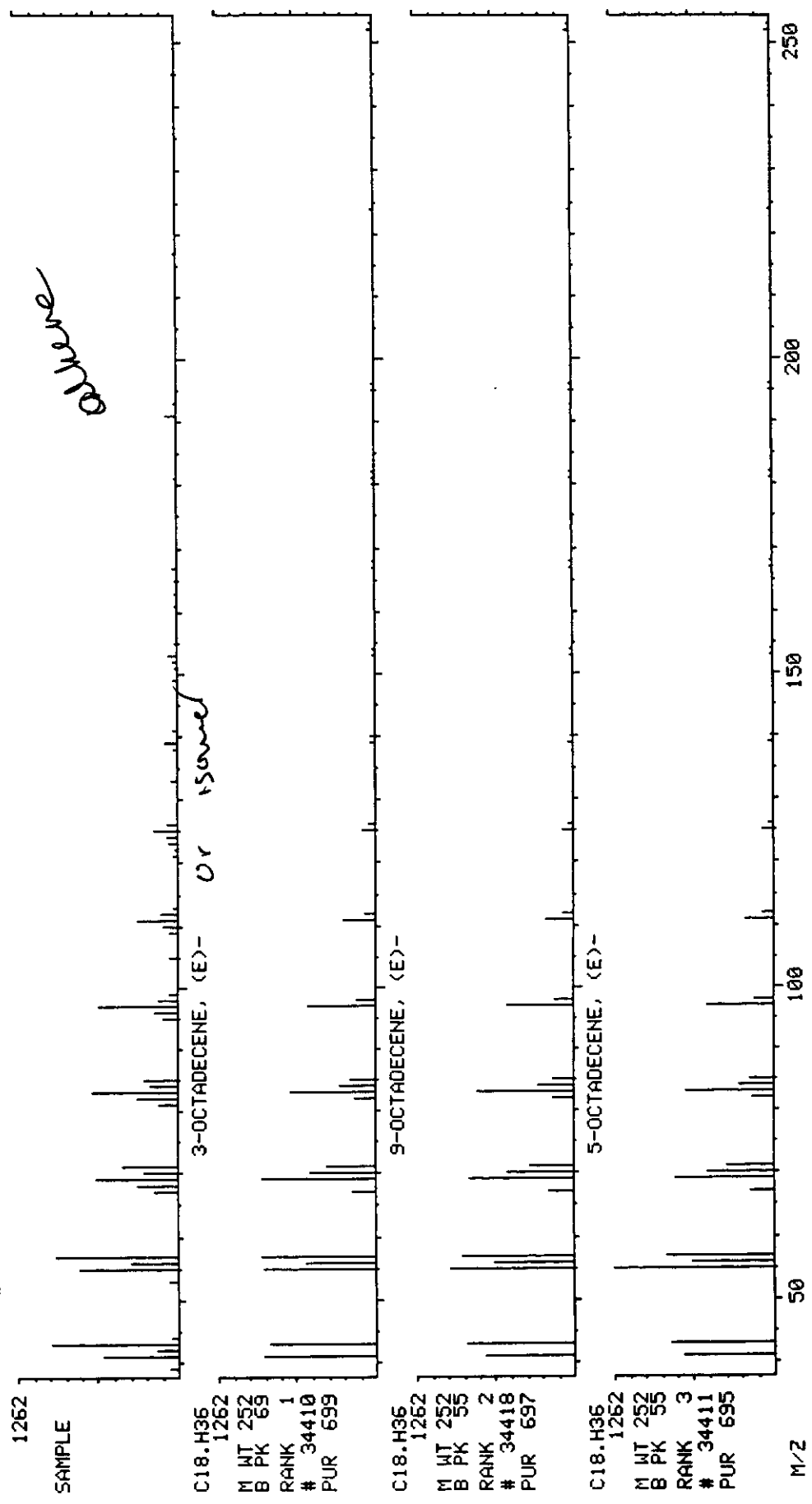
62231 spectra in LIBRARYNB searched for maximum PURITY
 950 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 34410 3-OCTADECENE, (E)-
 2 34418 9-OCTADECENE, (E)-
 3 34411 5-OCTADECENE, (E)-
 4 39515 9-EICOSENE, (E)-
 5 39516 5-EICOSENE, (E)-
 6 26416 4-TETRADECANOL
 7 31653 1-HEPTADECENE
 8 37062 1-NONADECENE
 9 40232 1-NONADECANOL

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C18.H36	252	69	699	930	723
2	C18.H36	252	55	697	927	719
3	C18.H36	252	55	695	925	718
4	C20.H40	280	55	693	942	717
5	C20.H40	280	55	691	939	716
6	C14.H30.O	214	43	689	955	711
7	C17.H34	238	55	687	933	710
8	C19.H38	266	97	687	933	722
9	C19.H40.O	284	55	684	965	703

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	7206-19-1
2	---	---	---	---	7206-25-9
3	---	---	---	---	7206-21-5
4	---	---	---	---	74685-29-3
5	---	---	---	---	74685-30-6
6	---	---	---	---	1653-33-4
7	---	---	---	---	6765-39-5
8	---	---	---	---	18435-45-5
9	---	---	---	---	1454-84-8

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:00:00 + 21:19
 SAMPLE: T-MM5-4 1/3SA/1ML
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/3SA NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068105 #1915
 CALI: 30068105 # 3
 BASE M/Z: 43
 RIC: 8736.



Library Search Data: 30068105 #1933 Base m/z: 59
 08/31/98 21:00:00 + 21:30 Cali: 30068105 # 3 RIC: 5008.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 705 matched at least 5 of the 16 largest peaks in the unknown

Rank In. Name
 1 59575 1S,20S-BICYCLO[18.16.0]-3,6,9,12,15,18,21,24,27,30,33,36-DODECAOXAH*
 2 41712 3,6,9,12,15-PENTAOXANONADECAN-1-OL
 3 33841 3,6,9,12-TETRAOXAHXADECAN-1-OL
 4 5825 2-PROPANOL, 1-(2-METHYLPROPOXY)-
 5 16648 1-BUTOXYETHOXY-2-PROPANOL
 6 5832 2-PROPANOL, 1-(1-METHYLPROPOXY)-
 7 31381 6-O-METHYL-2,4-METHYLENE-.BETA.-SEDOHEPTITOL
 8 16651 DIETHYLENE GLYCOL TERT-BUTYL ETHER METHYL ETHER
 9 1909 2-PROPANOL, 1-METHOXY-2-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C24.H46.O12	526	73	407	677	475
2	C14.H30.O6	294	45	381	704	423
3	C12.H26.O5	250	45	381	696	437
4	C7.H16.O2	132	57	380	798	410
5	C9.H20.O3	176	57	378	864	395
6	C7.H16.O2	132	43	376	795	402
7	C9.H18.O7	238	58	372	676	431
8	C9.H20.O3	176	57	372	840	389
9	C5.H12.O2	104	59	371	904	382

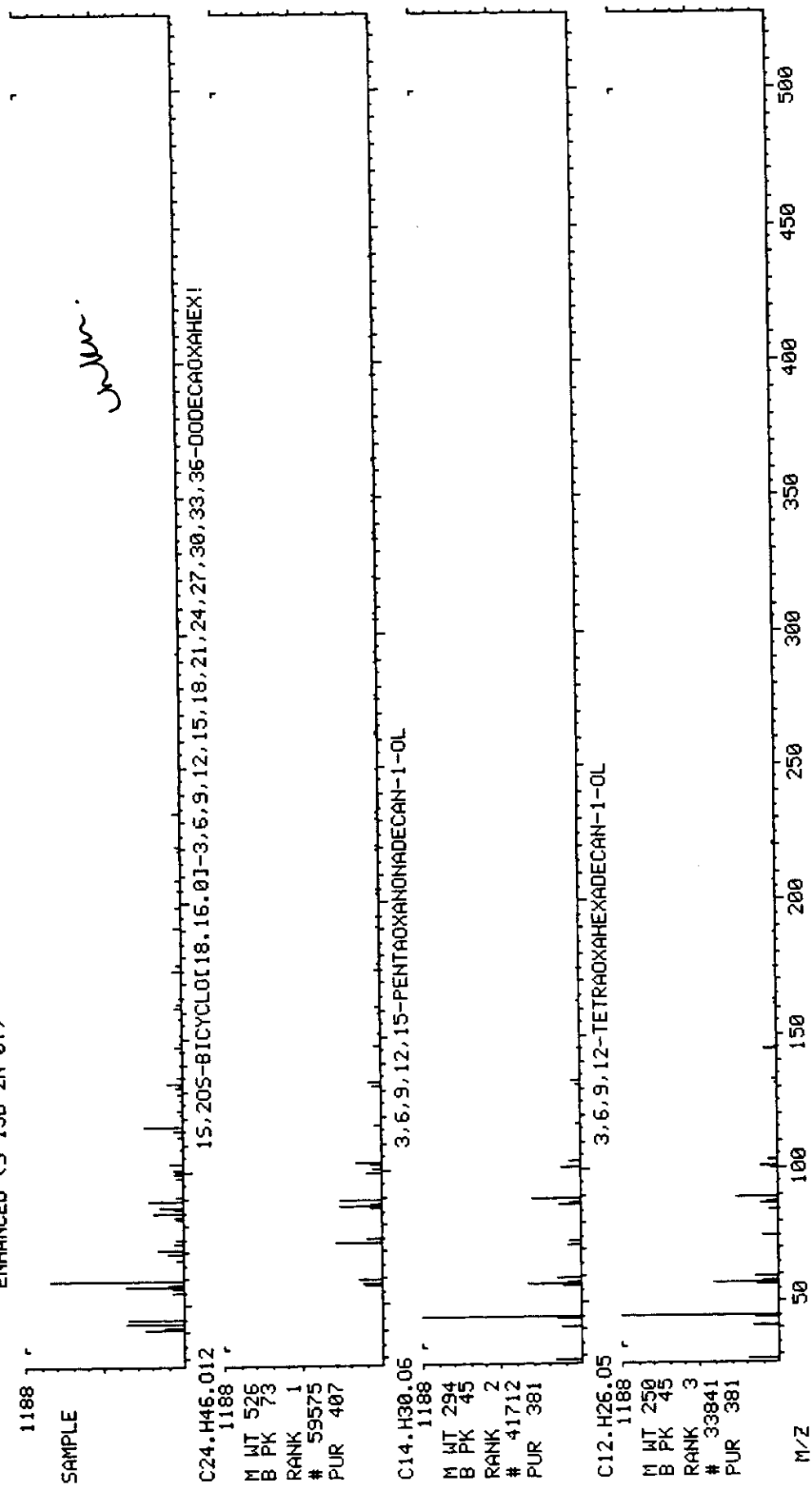
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	- -
2	---	---	---	---	1786-94-3
3	---	---	---	---	1559-34-8
4	---	---	---	---	23436-19-3
5	---	---	---	---	124-16-3
6	---	---	---	---	53907-95-2
7	---	---	---	---	- -
8	---	---	---	---	52788-79-1
9	---	---	---	---	3587-64-2

DATA: 30058105 #1933
CALI: 30058105 # 3

BASE M/Z: 59
RIC: 5008.

MID LIBRARY SEARCH <LIBRARYNB>
08/31/98 21:00:00 + 21:30
SAMPLE: T-MM5-4 1/35A/1ML
CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068105 #2057 Base m/z: 59
 08/31/98 21:00:00 + 22:53 Cali: 30068105 # 3 RIC: 3196.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 352 matched at least 5 of the 16 largest peaks in the unknown

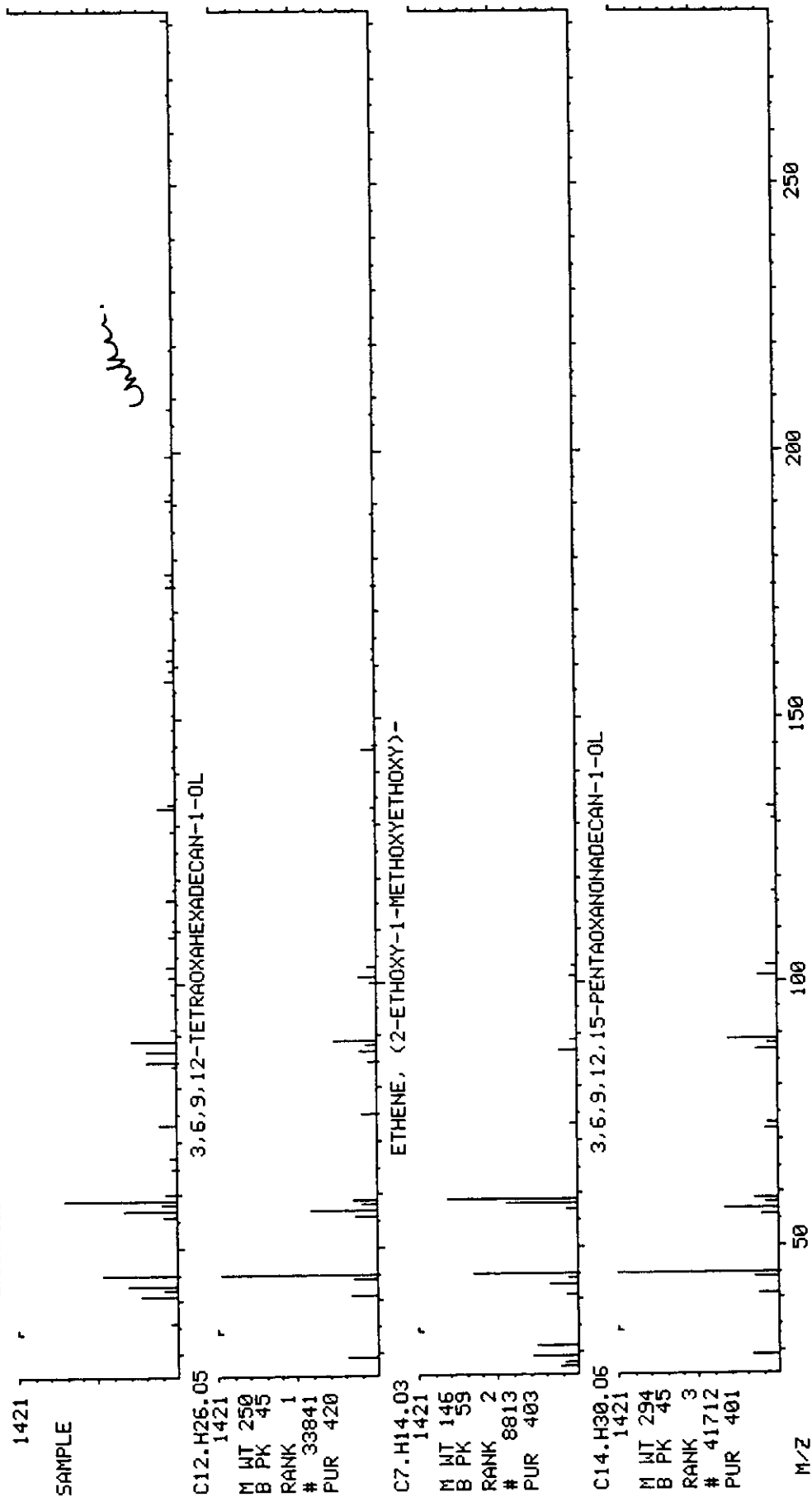
Rank In. Name
 1 33841 3,6,9,12-TETRAOXAHEXADECAN-1-OL
 2 8813 ETHENE, (2-ETHOXY-1-METHOXYETHOXY)-
 3 41712 3,6,9,12,15-PENTAOXANONADECAN-1-OL
 4 16651 DIETHYLENE GLYCOL TERT-BUTYL ETHER METHYL ETHER
 5 34224 2,5,8,11,14-PENTAOXAHEXADECAN-16-OL
 6 5832 2-PROPANOL, 1-(1-METHYLPROPOXY)-
 7 8871 METHANE, TERT-BUTOXYISOPROPOXY-
 8 16648 1-BUTOXYETHOXY-2-PROPANOL
 9 1909 2-PROPANOL, 1-METHOXY-2-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H26.O5	250	45	420	704	492
2	C7.H14.O3	146	59	403	823	409
3	C14.H30.O6	294	45	401	735	469
4	C9.H20.O3	176	57	397	782	436
5	C11.H24.O6	252	45	394	809	438
6	C7.H16.O2	132	43	383	863	419
7	C8.H18.O2	146	57	381	748	410
8	C9.H20.O3	176	57	380	780	444
9	C5.H12.O2	104	59	375	832	405

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	1559-34-8
2	—	—	—	—	54063-18-2
3	—	—	—	—	1786-94-3
4	—	—	—	—	52788-79-1
5	—	—	—	—	23778-52-1
6	—	—	—	—	53907-95-2
7	—	—	—	—	4346-01-4
8	—	—	—	—	124-16-3
9	—	—	—	—	3587-64-2

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 21:00:00 + 22:53

08/31/98 21:00:00 + 22:53 INST. ID: F16
SAMPLE: T-MM5-4 1/35A/ML
CONDS.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)



1017

TIC SELECTION REPORT

DATA FILE: 30068105

THE FOLLOWING PEAKS WERE REJECTED BECAUSE
AT LEAST 40 % OF THEIR SIZE WAS ACCOUNTED FOR BY
TARGET COMPOUNDS ELUTING WITHIN 4 SCANS OF THE
PEAK TOP.

SCAN	SIZE	AMOUNT
198	221024.	26.074
308	308392.	36.381
340	226080.	26.670
421	158208.	18.664
529	339072.	40.000
722	329502.	38.833
828	339403.	40.000
969	255936.	26.223
1086	390400.	40.000
1384	295584.	66.337
1384	285536.	64.082
1551	178232.	40.000
1782	139640.	40.000

TOTAL NUMBER OF UNIDENTIFIED PEAKS WITH SIZE
GREATER THAN 10 % OF THE CLOSEST INTERNAL STANDARD
THAT DOES NOT HAVE INTERFERENCES = 41

INTERNAL STANDARDS THAT HAVE RIC SIZE LESS THAN
50 % OR GREATER THAN 200 % OF THE ESTIMATED RIC SIZE
ARE CONSIDERED TO HAVE INTERFERENCES AND WILL NOT BE USED
FOR QUANTITATION.

#	INTERNAL STANDARD	PERCENT OF ESTIMATED		SCAN
		RIC SIZE	RIC SIZE	
1	C140 NAPHTHALENE-D8	339072.	157	529
2	C150 ACENAPHTHENE-D10	339403.	130	828
3	C160 PHENANTHRENE-D10	390400.	146	1086
4	C170 CHRYSENE-D12	178232.	130	1551
5	C175 PERYLENE-D12	139640.	121	1782

* INDICATES INTERFERENCE

SIZE = AREA

Semivolatile Organics
Method 0010/8270

Client Name: Pacific Environmental Services
Client ID: T-MM5-3-F, FH, XAD, COND, BH
LAB ID: 300681-0006-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 27 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
Phenol	53	ug/Sample	30	R
bis(2-Chloroethyl)ether	ND	ug/Sample	30	
2-Chlorophenol	ND	ug/Sample	30	
1,3-Dichlorobenzene	ND	ug/Sample	30	
1,4-Dichlorobenzene	ND	ug/Sample	30	
Benzyl alcohol	ND	ug/Sample	30	
1,2-Dichlorobenzene	ND	ug/Sample	30	
2-Methylphenol	ND	ug/Sample	30	
2,2'-Oxybis(1-chloropropane)	ND	ug/Sample	30	
3/4-Methylphenol	ND	ug/Sample	30	
N-Nitroso-di-n-propylamine	ND	ug/Sample	30	
Hexachloroethane	ND	ug/Sample	30	
Nitrobenzene	ND	ug/Sample	30	
Isophorone	ND	ug/Sample	30	
2-Nitrophenol	ND	ug/Sample	30	
2,4-Dimethylphenol	ND	ug/Sample	30	
Benzoic acid	ND	ug/Sample	150	
bis(2-Chloroethoxy)-methane	ND	ug/Sample	30	
2,4-Dichlorophenol	ND	ug/Sample	30	
1,2,4-Trichlorobenzene	ND	ug/Sample	30	
Naphthalene	39	ug/Sample	30	
4-Chloroaniline	ND	ug/Sample	30	
Hexachlorobutadiene	ND	ug/Sample	30	
4-Chloro-3-methylphenol	ND	ug/Sample	30	
2-Methylnaphthalene	74	ug/Sample	30	
Hexachlorocyclopentadiene	ND	ug/Sample	30	
2,4,6-Trichlorophenol	ND	ug/Sample	30	
2,4,5-Trichlorophenol	ND	ug/Sample	150	
2-Chloronaphthalene	ND	ug/Sample	30	
2-Nitroaniline	ND	ug/Sample	30	
Dimethyl phthalate	ND	ug/Sample	30	
Acenaphthylene	ND	ug/Sample	30	
3-Nitroaniline	ND	ug/Sample	150	
Acenaphthene	ND	ug/Sample	30	
2,4-Dinitrophenol	ND	ug/Sample	150	
4-Nitrophenol	ND	ug/Sample	150	
Dibenzofuran	ND	ug/Sample	30	

Note R = Reporting limit(s) raised due to sample volume limitations.
ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

1013

The cover letter is an integral part of this report.
Rev 230787

Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: T-MM5-3-F, FH, XAD, COND, BH
LAB ID: 300681-0006-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 27 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
2,4-Dinitrotoluene	ND	ug/Sample	30	
2,6-Dinitrotoluene	ND	ug/Sample	30	
Diethyl phthalate	ND	ug/Sample	30	
4-Chlorophenyl phenyl ether	ND	ug/Sample	30	
Fluorene	ND	ug/Sample	30	
4-Nitroaniline	ND	ug/Sample	150	
4,6-Dinitro-2-methylphenol	ND	ug/Sample	150	
N-Nitrosodiphenylamine	ND	ug/Sample	30	
4-Bromophenyl phenyl ether	ND	ug/Sample	30	
Hexachlorobenzene	ND	ug/Sample	30	
Pentachlorophenol	ND	ug/Sample	150	
Phenanthrene	ND	ug/Sample	30	
Anthracene	ND	ug/Sample	30	
Di-n-butyl phthalate	ND	ug/Sample	30	
Fluoranthene	ND	ug/Sample	30	
Pyrene	ND	ug/Sample	30	
Butyl benzyl phthalate	ND	ug/Sample	30	
3,3'-Dichlorobenzidine	ND	ug/Sample	60	
Benzo(a)anthracene	ND	ug/Sample	30	
bis(2-Ethylhexyl)-phthalate	ND	ug/Sample	30	
Chrysene	ND	ug/Sample	30	
Di-n-octyl phthalate	ND	ug/Sample	30	
Benzo(b)fluoranthene	ND	ug/Sample	30	
Benzo(k)fluoranthene	ND	ug/Sample	30	
Benzo(a)pyrene	ND	ug/Sample	30	
Indeno(1,2,3-cd)pyrene	ND	ug/Sample	30	
Dibenz(a,h)anthracene	ND	ug/Sample	30	
Benzo(g,h,i)perylene	ND	ug/Sample	30	
Acetophenone	ND	ug/Sample	30	
4-Aminobiphenyl	ND	ug/Sample	150	
Aniline	ND	ug/Sample	30	
Benzidine	ND	ug/Sample	300	
3,3'-Dimethylbenzidine	ND	ug/Sample	60	
N-Nitrosodimethylamine	ND	ug/Sample	30	
N-Nitrosomorpholine	ND	ug/Sample	30	
Pentachloronitrobenzene (PCNB)	ND	ug/Sample	150	
o-Toluidine	ND	ug/Sample	60	

ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.
Rev 230787

1020

Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: T-MM5-3-F, FH, XAD, COND, BH
LAB ID: 300681-0006-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 27 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
2-Methoxybenzenamine	ND	ug/Sample	--	K
Biphenyl	ND	ug/Sample	--	K
Chloroacetophenone	ND	ug/Sample	--	K
Cumene	ND	ug/Sample	--	K
DBCP (1,2-Dibromo-3-chloropropane)	ND	ug/Sample	--	K
Benzo(e)pyrene	ND	ug/Sample	--	K
N-N-Diethylaniline	ND	ug/Sample	--	K
Dimethylaniline	ND	ug/Sample	--	K
3,3'-Dimethoxybenzidine	ND	ug/Sample	--	K
Hydroquinone	ND	ug/Sample	--	K
4,4'-Methyl-bis(2-chloroaniline)	ND	ug/Sample	--	K
4-Nitrodiphenyl	ND	ug/Sample	--	K
Trifluralin	ND	ug/Sample	--	K

Surrogate	Recovery	Acceptable Range	
Nitrobenzene-d5	78 %	45 - 107	
2-Fluorobiphenyl	106 %	62 - 110	
Terphenyl-d14	80 %	58 - 135	
Phenol-d5	69 %	43 - 130	
2-Fluorophenol	71 %	36 - 111	
2,4,6-Tribromophenol	52 %	58 - 131	i

Note i = Surrogate recovery is outside of control limits.

Note K = Identified by mass spectrum only; quantitation based on 1:1 response with internal standard.

ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.
Rev 230787

1021

Semivolatiles Library Search (20 Compound TID)
Method 8270

Client Name: Pacific Environmental Services
Client ID: T-MM5-3-F, FH, XAD, COND, BH
LAB ID: 300681-0006-SA

Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 27 JUL 98
Prepared: NA

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Result	Units	Reporting Limit	Qualifier
n-Nonane	390	ug/Sample	--	0
Unknown	270	ug/Sample	--	
Undecane	180	ug/Sample	--	0
Dodecane	160	ug/Sample	--	0
Unknown alkane	240	ug/Sample	--	
1H-Indene, 1-ethylidene-	170	ug/Sample	--	0
Undecane, 4,7-dimethyl-	240	ug/Sample	--	0
4-Hydroxy-4-methyl-2-pentanone	350	ug/Sample	--	0
Heptadecane, 2,6,10,14 -tetramethyl-	250	ug/Sample	--	0
Tetradecane	430	ug/Sample	--	0
Unknown aromatic hydrocarbon	230	ug/Sample	--	
Heptadecane, 2,6,10,14 -tetramethyl-	300	ug/Sample	--	0
Pentadecane	440	ug/Sample	--	
Nonadecane	400	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	230	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	520	ug/Sample	--	0
Nonadecane	170	ug/Sample	--	0
Nonadecane	180	ug/Sample	--	0
Nonadecane	150	ug/Sample	--	0
9-Eicosene (E)-	240	ug/Sample	--	0

Note 0 = Or structurally similar compound (isomer).
NA = Not Applicable

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.
Rev 230787

1022

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068106
Std Id: ST16980831

Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 21:30
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Surrogate Spike Recoveries
8270-G, LIMS 10/10/96

Surrogate	Surrogate Amount (ug)		% Recovery	
	Spiked	Measured	Measured	QC limits
CS20 NITROBENZENE-D5	50.00	39.20	78.4	45 107
CS25 2-FLUOROBIPHENY	50.00	52.90	106.	62 110
CS30 TERPHENYL-D14	50.00	39.90	79.8	58 135
CS45 PHENOL-D5	100.0	68.90	68.9	43 130
CS50 2-FLUOROPHENOL	100.0	71.40	71.4	36 111
CS55 2,4,6-TRIBROMOP	100.0	52.50	52.5	* 58 131

RI confirms

Target Compounds: SAP9

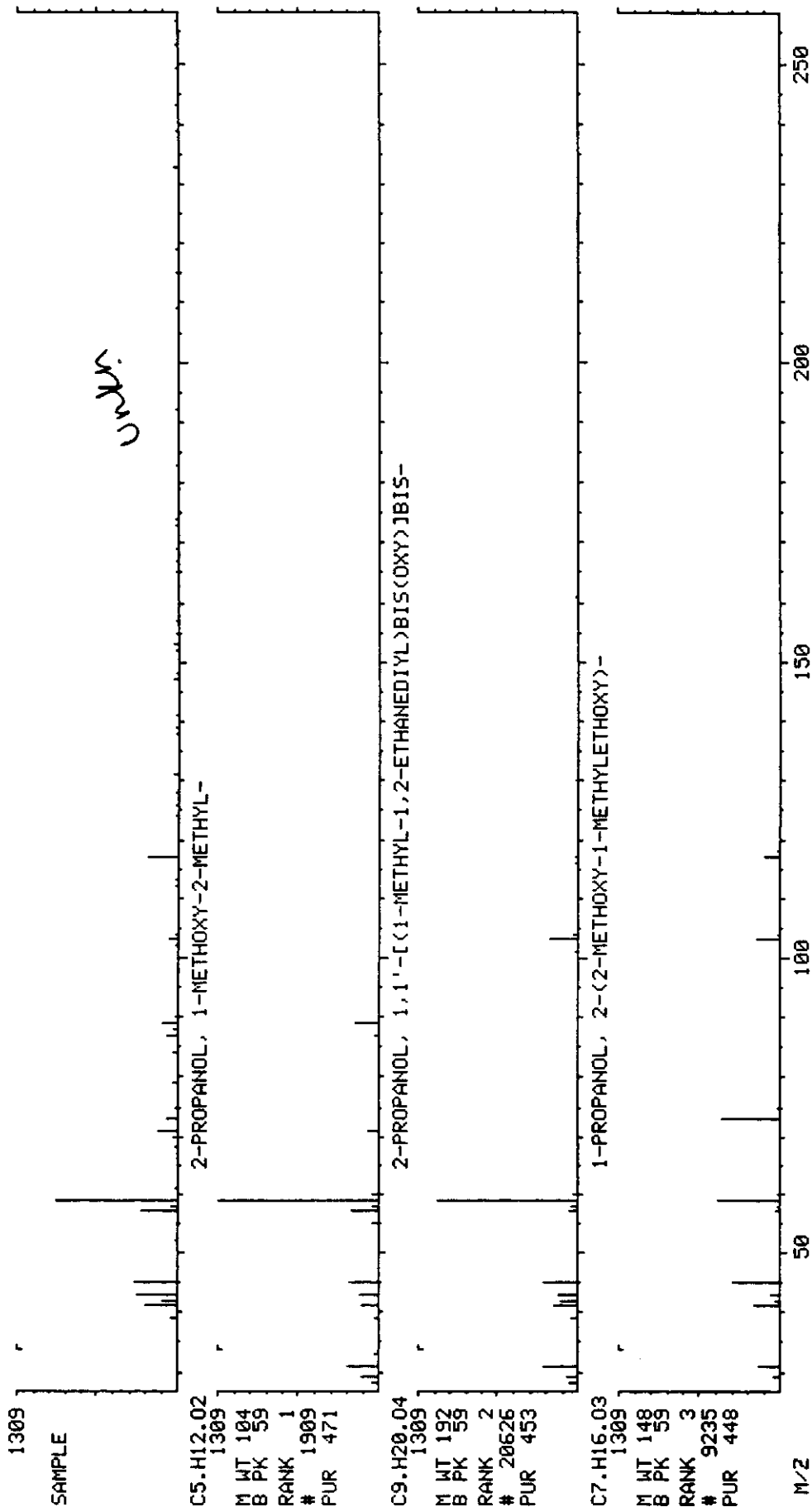
29/16/98

Parameter	Scan	Result	Units	Reporting Limit
HEXACHLOROBENZENE-C13		ND	UG/A	30.0
C310 N-NITROSODIMETHYLAMINE		ND	UG/A	30.0
PYRIDINE		ND	UG/A	60.0
2-PICOLINE		ND	UG/A	30.0
N-NITROSOMETHYLETHYLAMINE	157	58 ND	UG/A	30.0
METHYLMETHANESULFONATE		ND	UG/A	30.0
N-NITROSODIETHYLAMINE		ND	UG/A	30.0
ETHYLMETHANE SULFONATE		ND	UG/A	30.0
PENTACHLOROETHANE		ND	UG/A	30.0
C320 ANILINE		ND	UG/A	30.0
C315 PHENOL	311	53	UG/A	30.0
C325 BIS(2-CHLOROETHYL)ETHER		ND	UG/A	30.0
C330 2-CHLOROPHENOL		ND	UG/A	30.0
C335 1,3-DICHLOROBENZENE		ND	UG/A	30.0
C340 1,4-DICHLOROBENZENE		ND	UG/A	30.0
C345 BENZYL ALCOHOL		ND	UG/A	30.0
C350 1,2-DICHLOROBENZENE		ND	UG/A	30.0
C355 2-METHYLPHENOL		ND	UG/A	30.0
C360 2,2'-OXYBIS(1-CLPROPAN)		ND	UG/A	30.0
C361 ACETOPHENONE	401	31 NA	UG/A	30.0
N-NITROSOPYRROLIDINE		ND	UG/A	30.0
N-NITROSOMORPHOLINE	410	31 ND	UG/A	30.0
3-METHYL PHENOL		ND	UG/A	30.0
C365 4-METHYLPHENOL		ND	UG/A	30.0
C370 N-NITROSO-DI-N-PROPYLAM		ND	UG/A	30.0
O-TOLUIDINE		ND	UG/A	30.0
C375 HEXACHLOROETHANE		ND	UG/A	30.0
C410 NITROBENZENE		ND	UG/A	30.0
N-NITROSOPIPERIDINE		ND	UG/A	30.0

Reviewed by:

29/17/98

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:00:00 + 18:29
 SAMPLE: T-MM5-4 1/35A/1ML
 COND5.: UG/ML *1ML *100%/100% *(NA/NA) >1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30068105 #1652
 CALI: 30068105 # 3
 BASE M/Z: 59
 RIC: 5128.



Library Search Data: 30068105 #1688 Base m/z: 41
 08/31/98 21:00:00 + 18:47 Cali: 30068105 # 3 RIC: 2780.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 246 matched at least 5 of the 16 largest peaks in the unknown

Rank In. Name
 1 15753 2,2-DIMETHYL-3-HEXANOL ACETATE
 2 43380 BUTANOIC ACID, 4-(2,4-DICHLOROPHENOXY)-, 2-METHYLPROPYL ESTER
 3 43381 2,4-DB BUTYL ESTER
 4 11858 1,3,2-DIOXABOROLANE, 4-METHYL-2-(2-METHYLPROPOXY)-
 5 11857 1,3,2-DIOXABORINANE, 2-(2-METHYLPROPOXY)-
 6 3130 2-BUTANONE, 3-METHYL-, O-METHYLOXIME
 7 12049 4-OCTANOL, 4,7-DIMETHYL-
 8 3139 2-PENTANONE, O-METHYLOXIME
 9 52020 D-MANNITOL, 2,4-DI-O-METHYL-, TETRAACETATE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C10.H20.O2	172	43	240	646	291
2	C14.H18.O3.CL2	304	87	224	660	270
3	C14.H18.O3.CL2	304	87	220	664	247
4	C7.H15.O3.B	158	115	208	609	273
5	C7.H15.O3.B	158	115	199	557	237
6	C6.H13.O.N	115	42	198	758	211
7	C10.H22.O	158	87	187	660	234
8	C6.H13.O.N	115	42	185	707	196
9	C16.H26.O10	378	43	175	622	217

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	- -
2	—	—	—	—	51550-64-2
3	—	—	—	—	6753-24-8
4	—	—	—	—	52910-21-1
5	—	—	—	—	55162-67-9
6	—	—	—	—	27685-13-8
7	—	—	—	—	19781-13-6
8	—	—	—	—	31377-00-1
9	—	—	—	—	24406-90-4

Library Search Data: 30068105 #1804 Base m/z: 59
 08/31/98 21:00:00 + 20:04 Cali: 30068105 # 3 RIC: 4392.
 Sample: T-MM5-4 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

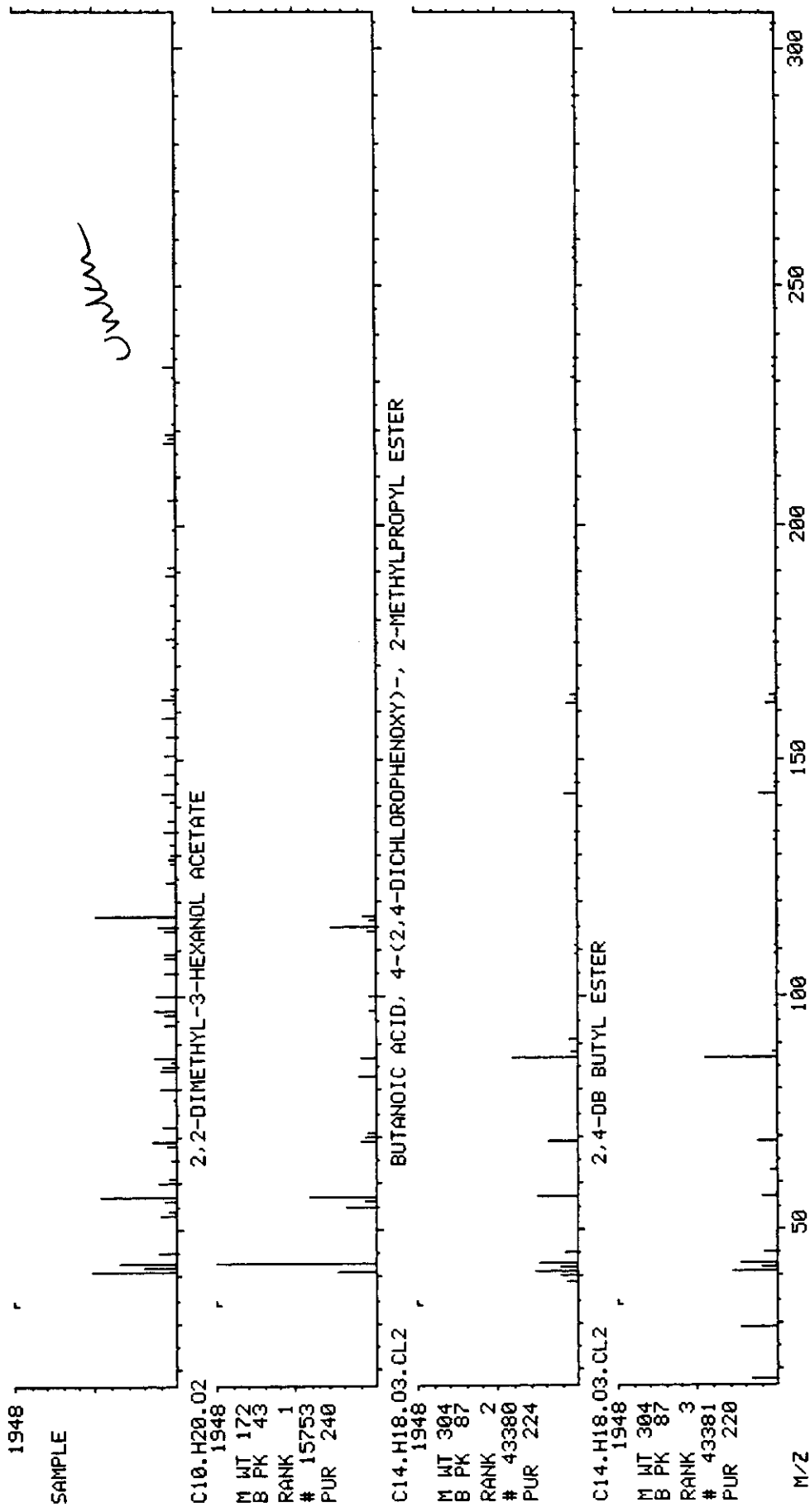
62231 spectra in LIBRARYNB searched for maximum PURITY
 465 matched at least 5 of the 16 largest peaks in the unknown

Rank In.	Name
1	9228 1-PROPANOL, 2-(2-METHOXYPROPOXY)-
2	24247 2-PROPANOL, 1-[2-(2-METHOXY-1-METHYLETHOXY)-1-METHYLETHOXY]-
3	8822 BUTANOIC ACID, 4-ETHOXY-, METHYL ESTER
4	8432 2-HEXANONE, 3-HYDROXY-3,5-DIMETHYL-
5	1760 2-PENTANOL, 2-METHYL-
6	1905 2-BUTANOL, 3-METHOXY-
7	16649 1-PROPANOL, 3-[3-(1-METHYLETHOXY)PROPOXY]-
8	5832 2-PROPANOL, 1-(1-METHYLPROPOXY)-
9	9235 1-PROPANOL, 2-(2-METHOXY-1-METHYLETHOXY)-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C7.H16.O3	148	59	436	799	454
2	C10.H22.O4	206	59	397	681	490
3	C7.H14.O3	146	59	392	687	433
4	C8.H16.O2	144	59	369	851	391
5	C6.H14.O	102	59	365	867	385
6	C5.H12.O2	104	59	362	825	392
7	C9.H20.O3	176	59	356	768	413
8	C7.H16.O2	132	43	354	624	433
9	C7.H16.O3	148	59	349	668	423

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	13588-28-8
2	—	—	—	—	20324-33-8
3	—	—	—	—	29006-04-0
4	—	—	—	—	6321-14-8
5	—	—	—	—	590-36-3
6	—	—	—	—	53778-72-6
7	—	—	—	—	54518-03-5
8	—	—	—	—	53907-95-2
9	—	—	—	—	55956-21-3

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:00:00 + 18:47
 SAMPLE: T-NMS-4 1/35A/1ML
 COND.: UG/ML *1ML *1002/1002 *(NA/NA >1/35A NA M
 ENHANCED (5 15B 2N 0T)
 DATA: 30068105 #1688
 CALI: 30068105 # 3
 BASE M/Z: 41
 RIC: 2780.



QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068106
Std Id: ST16980831

Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 21:30
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C415 ISOPHORONE		ND	UG/A	30.0
C420 2-NITROPHENOL		ND	UG/A	30.0
C425 2,4-DIMETHYLPHENOL		ND	UG/A	30.0
C435 BIS(2-CHLOROETHOXY)METH		ND	UG/A	30.0
C440 2,4-DICHLOROPHENOL		ND	UG/A	30.0
C430 BENZOIC ACID		ND	UG/A	150.0
AA-DIMETHYLPHENETHYLAMINE		ND	UG/A	30.0
C445 1,2,4-TRICHLOROBENZENE		ND	UG/A	30.0
C450 NAPHTHALENE	534	39	UG/A	30.0
C455 4-CHLOROANALINE		ND	UG/A	30.0
2,6-DICHLOROPHENOL		ND	UG/A	30.0
HEXACHLOROPROPENE		ND	UG/A	30.0
C460 HEXACHLOROBUTADIENE		ND	UG/A	30.0
P-PHENYLENE DIAMINE		ND	UG/A	30.0
N-NITROSODI-N-BUTYLAMINE		ND	UG/A	30.0
C465 4-CHLORO-3-METHYLPHENO		ND	UG/A	30.0
SAFROLE		ND	UG/A	30.0
C470 2-METHYLNAPHTHALENE	653	74	UG/A	30.0
1,2,4,5-TETRACHLOROBENZENE		ND	UG/A	30.0
ISOSAFROLE (#1)		ND	UG/A	60.0
C510 HEXACHLOROCYCLOPENTADI		ND	UG/A	30.0
C515 2,4,6-TRICHLOROPHENOL		ND	UG/A	30.0
C520 2,4,5-TRICHLOROPHENOL		ND	UG/A	30.0
ISOSAFROLE (#2)		ND	UG/A	60.0
C525 2-CHLORONAPHTHALENE		ND	UG/A	30.0
1-CHLORONAPHTHALENE		ND	UG/A	15.0
C530 2-NITROANALINE		ND	UG/A	150.0
1,4-NAPHTHOQUINONE		ND	UG/A	30.0
C535 DIMETHYLPHTHALATE		ND	UG/A	30.0
1,3-DINITROBENZENE		ND	UG/A	30.0
C540 ACENAPHTHYLENE		ND	UG/A	30.0
C543 2,6-DINITROTOLUENE		ND	UG/A	30.0
C545 3-NITROANILINE		ND	UG/A	150.0
C550 ACENAPHTHENE		ND	UG/A	30.0
C555 2,4-DINITROPHENOL		ND	UG/A	150.0
C565 DIBENZOFURAN		ND	UG/A	30.0
C560 4-NITROPHENOL		ND	UG/A	150.0
PENTACHLOROBENZENE		ND	UG/A	30.0
C570 2,4-DINITROTOLUENE		ND	UG/A	30.0
1-NAPHTHYLAMINE		ND	UG/A	30.0
2-NAPHTHYLAMINE		ND	UG/A	30.0
2,3,4,6-TETRACHLOROPHENOL		ND	UG/A	60.0
C580 DIETHYLPHTHALATE		ND	UG/A	30.0
C590 FLUORENE		ND	UG/A	30.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068106
Std Id: ST16980831

Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 21:30
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
=====				
C585 4-CHLOROPHENYL-PHENYLE		ND	UG/A	30.0
5-NITRO-O-TOLUIDINE		ND	UG/A	30.0
C595 4-NITROANALINE		ND	UG/A	150.0
C610 4,6-DINITRO-2-METHYLPH		ND	UG/A	150.0
C615 N-NITROSODIPHENYLAMINE		ND	UG/A	30.0
C620 AZOBENZENE		ND	UG/A	30.0
SYM-TRINITROBENZENE		ND	UG/A	30.0
C625 4-BROMOPHENYL-PHENYLET		ND	UG/A	30.0
PHENACETIN		ND	UG/A	30.0
DIALATE	234	ND	UG/A	30.0
C630 HEXACHLOROBENZENE		ND	UG/A	30.0
4-AMINOBIHENYL		ND	UG/A	30.0
C635 PENTACHLOROPHENOL		ND	UG/A	150.0
PRONAMIDE		ND	UG/A	30.0
PENTACHLORONITROBENZENE		ND	UG/A	150.0
C640 PHENANTHRENE		ND	UG/A	30.0
C645 ANTHRACENE		ND	UG/A	30.0
2SECBUTYL-4,6-DINITROPHENOL		ND	UG/A	30.0
C647 CARBAZOLE		ND	UG/A	30.0
C650 DI-N-BUTYLPHTHALATE		ND	UG/A	30.0
4-NITROQUINOLINE-1-OXIDE		ND	UG/A	30.0
METHAPYRILENE		ND	UG/A	30.0
ISODRIN		ND	UG/A	30.0
C655 FLUORANTHENE		ND	UG/A	30.0
CHLOROBENZILATE		ND	UG/A	30.0
C710 BENZIDINE		ND	UG/A	300.0
C715 PYRENE		ND	UG/A	30.0
ARAMITE (#1)		ND	UG/A	30.0
ARAMITE (#2)		ND	UG/A	30.0
P-DIMETHYLAMINOAZOBENZENE		ND	UG/A	30.0
3,3'-DIMETHYLBENZIDINE		ND	UG/A	30.0
KEPONE		ND	UG/A	150.0
C720 BUTYLBENZYLPHTHALATE		ND	UG/A	30.0
2-ACETYLAMINOFLUORENE		ND	UG/A	30.0
C730 BENZO(A)ANTHRACENE		ND	UG/A	30.0
C725 3,3'-DICHLOROBENZIDINE		ND	UG/A	60.0
C740 CHRYSENE		ND	UG/A	30.0
C745 BIS(2-ETHYLHEXYL)PHTHA		ND	UG/A	30.0
3-METHYLCHOLANTHRENE		ND	UG/A	30.0
C760 DI-N-OCTYL PHTHALATE		ND	UG/A	30.0
C765 BENZO(B)FLUORANTHENE		ND	UG/A	30.0
7,12-DIMETHYLBENZANTHRACENE		ND	UG/A	30.0
C770 BENZO(K)FLUORANTHENE		ND	UG/A	30.0
HEXACHLOROPHENE		ND	UG/A	30.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

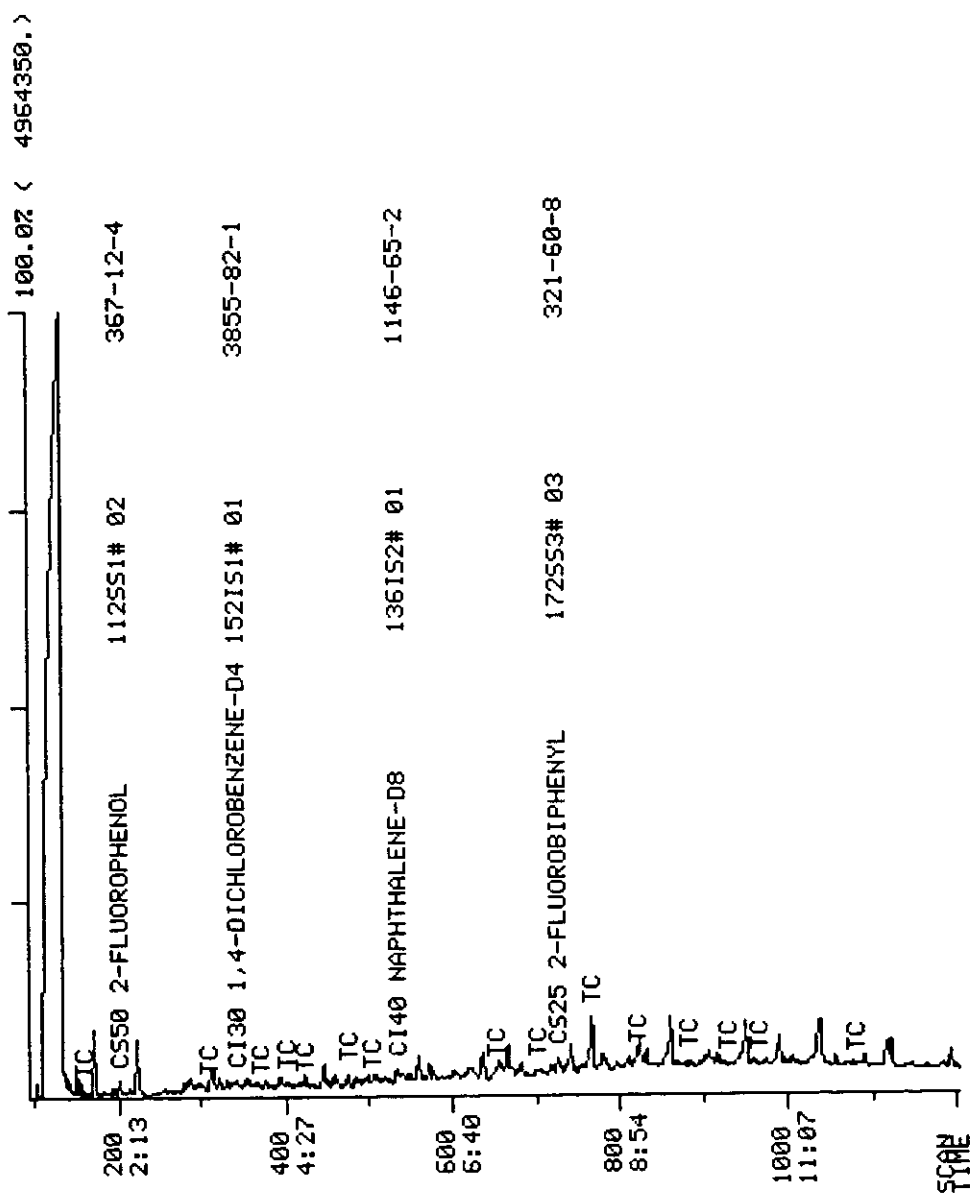
Data File: 30068106
Std Id: ST16980831

Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 21:30
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

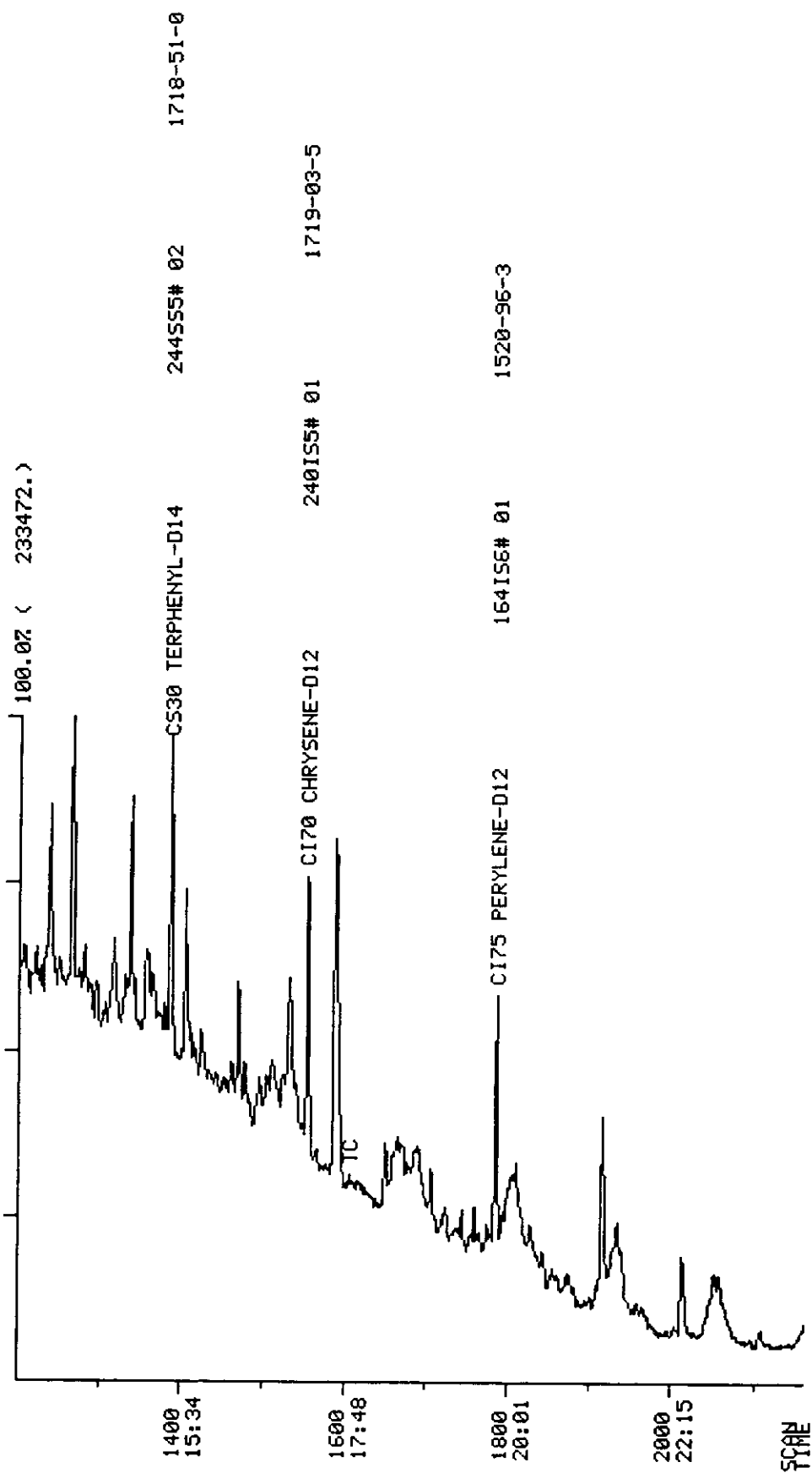
Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C775 BENZO(A)PYRENE		ND	UG/A	30.0
C780 INDENO(1,2,3-CD)PYRENE		ND	UG/A	30.0
C785 DIBENZ(A,H)ANTHRACENE		ND	UG/A	30.0
C790 BENZO(G,H,I)PERYLENE		ND	UG/A	30.0

DATA FROM FILE: 30068106 SCANS 93 TO 1203 ACQUIRED: 08/31/98 21:30:00
 CALI: 30068106 #3
 SAMPLE: T-MM5-3 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA) /1/35A NA M



DATA FROM FILE: 30068106
SCANS 1203 TO 2158 ACQUIRED: 08/31/98 21:30:00
CALI: 30068106 #3
SAMPLE: T-HMS-3 1/35A/1ML INST. ID: F16
CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA) /1/35A NA M



ADDITIONAL STANDARD USED FOR THIS REPORT IS: ST16980831A

QUANTERRA QUANTITATION SUMMARY

File: 30068106

Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Analyst: DAT Instrument Id: F16 Analyzed: 08/31/98 21:30

Compounds with amounts less than 0.20 reported as NOT FOUND

Library						Units: UG/ML	
No Entry	Name	Mass Meth	Scan Ref	Area	RRF(L)	Amount	
1 S1#	1 C130 1,4-DICHLOROBENZENE	152 A BB	341 1	33301.	1.000	40.000	
2 S2#	1 C140 NAPHTHALENE-D8	136 A BB	531 2	120808.	1.000	40.000	
3 S3#	1 C150 ACENAPHTHENE-D10	164 A BB	831 3	61921.	1.000	40.000	
4 S4#	1 C160 PHENANTHRENE-D10	188 A BB	1088 4	101964.	1.000	40.000	
5 S5#	1 C170 CHRYSENE-D12	240 A BB	1553 5	101687.	1.000	40.000	
6 S6#	1 C175 PERYLENE-D12	264 A VB	1784 6	95704.	1.000	40.000	
7 S2#	2 CS20 NITROBENZENE-D5	82 A BV	422 2	55580.	0.470	39.192	
8 S3#	3 CS25 2-FLUOROBIPHENYL	172 A VB	724 3	101124.	1.234	52.925	
9 S5#	2 CS30 TERPHENYL-D14	244 A BB	1385 5	100901.	0.995	39.882	
10 S1#	3 CS45 PHENOL-D5	99 A BB	310 1	116232.	2.025	68.929	
11 S1#	2 CS50 2-FLUOROPHENOL	112 A BB	198 1	81364.	1.368	71.429	
12 S3#	2 CS55 2,4,6-TRIBROMOPHENO	330 A BB	972 3	19151.	0.236	52.497	
13 S4#	4 HEXACHLOROBENZENE-C13	294	4 NOT FOUND				
14 S1#	15 C310 N-NITROSODIMETHYLAM	74	1 NOT FOUND				
15 S1#	20 PYRIDINE	79	1 NOT FOUND				
16 S1#	30 2-PICOLINE	93	1 NOT FOUND				
17 S1#	40 N-NITROSOMETHYLETHYLAMIN	42 A VV	157 1	8518.	0.811	12.614	
18 S1#	60 METHYLMETHANESULFONATE	80 A BB	183 1	468.	0.636	0.884	
19 S1#	70 N-NITROSODIETHYLAMINE	102	1 NOT FOUND				
20 S1#	85 ETHYLMETHANE SULFONATE	79	1 NOT FOUND				
21 S1#	95 PENTACHLOROETHANE	117 A BB	307 1	1144.	0.602	2.285	
22 S1#	100 C320 ANILINE	93	1 NOT FOUND				
23 S1#	105 C315 PHENOL	94 A BB	311 1	28734.	1.968	17.540	
24 S1#	110 C325 BIS(2-CHLOROETHYL)E	93 A BB	320 1	3814.	1.368	3.349	
25 S1#	115 C330 2-CHLOROPHENOL	128 A BV	315 1	330.	1.497	0.265	
26 S1#	125 C335 1,3-DICHLOROBENZENE	146	1 NOT FOUND				
27 S1#	130 C340 1,4-DICHLOROBENZENE	146	1 NOT FOUND				
28 S1#	145 C345 BENZYL ALCOHOL	108 A VB	370 1	6573.	0.854	9.241	
29 S1#	150 C350 1,2-DICHLOROBENZENE	146	1 NOT FOUND				
30 S1#	160 C355 2-METHYLPHENOL	108	1 NOT FOUND				
31 S1#	165 C360 2,2'-OXYBIS(1-CLPRO	45 A BV	386 1	640.	2.539	0.303	
32 S1#	170 C361 ACETOPHENONE	105 A BB	401 1	16818.	1.934	10.446	
33 S1#	175 N-NITROSOPYRROLIDINE	100 A BB	404 1	250.	0.553	0.543	
34 S1#	180 N-NITROSOMORPHOLINE	56 A BB	410 1	7023.	0.821	10.279	
35 S1#	182 3-METHYL PHENOL	108 A BB	420 1	4135.	2.252	2.206	
36 S1#	185 C365 4-METHYLPHENOL	108	1 NOT FOUND				
37 S1#	190 C370 N-NITROSO-DI-N-PROP	70 A VB	404 1	1969.	0.906	2.611	
38 S1#	195 O-TOLUIDINE	106	1 NOT FOUND				
39 S1#	200 C375 HEXACHLOROETHANE	117	1 NOT FOUND				
40 S2#	10 C410 NITROBENZENE	77 A VB	423 2	2891.	0.445	2.151	
41 S2#	15 N-NITROSOPIPERIDINE	42	2 NOT FOUND				
42 S2#	20 C415 ISOPHORONE	82 A BB	473 2	10146.	0.741	4.532	
43 S2#	25 C420 2-NITROPHENOL	139	2 NOT FOUND				
44 S2#	30 C425 2,4-DIMETHYLPHENOL	107 A BB	501 2	1308.	0.330	1.313	
45 S2#	35 C435 BIS(2-CHLOROETHOXY)	93 A BB	516 2	744.	0.511	0.483	

46	S2#	40	C440	2,4-DICHLOROPHENOL	162	A	BB	520	2	548.	0.308	0.590
47	S2#	45	C430	BENZOIC ACID	122			2	NOT FOUND			
48	S2#	50	AA	DIMETHYLPHENETHYLAMIN	58			2	NOT FOUND			
49	S2#	55	C445	1,2,4-TRICHLOROBENZ	180			2	NOT FOUND			

50	S2# 60	C450 NAPHTHALENE	128 A BV	534 2	40335. 1.022	13.065
51	S2# 80	C455 4-CHLOROANALINE	127 A BB	557 2	1230. 0.454	0.898
52	S2# 85	2,6-DICHLOROPHENOL	162	2 NOT FOUND		
53	S2# 90	HEXACHLOROPROPENE	213	2 NOT FOUND		
54	S2# 95	C460 HEXACHLOROBUTADIENE	225	2 NOT FOUND		
55	S2#115	P-PHENYLENE DIAMINE	108	2 NOT FOUND		
56	S2#120	N-NITROSODI-N-BUTYLAMINE	84	2 NOT FOUND		
57	S2#130	C465 4-CHLORO-3-METHYLPH	107	2 NOT FOUND		
58	S2#140	SAFROLE	162 A BB	649 2	1554. 0.248	2.077
59	S2#145	C470 2-METHYLNAPHTHALENE	142 A BB	653 2	56007. 0.751	24.684
60	S3# 10	1,2,4,5-TETRACHLOROBENZE	214	3 NOT FOUND		
61	S3# 15	ISOSAFROLE (#1)	162 A BB	699 3	1059. 0.044	15.673
62	S3# 20	C510 HEXACHLOROCYCLOPENT	237	3 NOT FOUND		
63	S3# 25	C515 2,4,6-TRICHLOROPHEN	196	3 NOT FOUND		
64	S3# 30	C520 2,4,5-TRICHLOROPHEN	196	3 NOT FOUND		
65	S3# 35	ISOSAFROLE (#2)	104	3 NOT FOUND		
66	S3# 40	C525 2-CHLORONAPHTHALENE	162	3 NOT FOUND		
67	S3# 42	1-CHLORONAPHTHALENE	162	3 NOT FOUND		
68	S3# 45	C530 2-NITROANALINE	65 A BB	765 3	1407. 0.502	1.812
69	S3# 50	1,4-NAPHTHOQUINONE	158	3 NOT FOUND		
70	S3# 55	C535 DIMETHYLPHTHALATE	163	3 NOT FOUND		
71	S3# 60	1,3-DINITROBENZENE	168 A BB	808 3	278. 0.197	0.909
72	S3# 65	C540 ACENAPHTHYLENE	152 A BB	809 3	1258. 1.838	0.442
73	S3# 70	C543 2,6-DINITROTOLUENE	165 A BB	818 3	645. 0.324	1.287
74	S3# 75	C545 3-NITROANILINE	138	3 NOT FOUND		
75	S3# 80	C550 ACENAPHTHENE	153 A VB	836 3	2658. 1.159	1.482
76	S3# 85	C555 2,4-DINITROPHENOL	184 A BV	859 3	224. 0.189	0.764
77	S3# 90	C565 DIBENZOFURAN	168 A BB	868 3	1768. 1.669	0.684
78	S3# 95	C560 4-NITROPHENOL	109	3 NOT FOUND		
79	S3#100	PENTACHLOROBENZENE	250	3 NOT FOUND		
80	S3#105	C570 2,4-DINITROTOLUENE	165 A BB	882 3	682. 0.397	1.109
81	S3#110	1-NAPHTHYLAMINE	143 A BV	883 3	602. 0.837	0.464
82	S3#115	2-NAPHTHYLAMINE	143	3 NOT FOUND		
83	S3#120	2,3,4,6-TETRACHLOROPHENO	232	3 NOT FOUND		
84	S3#130	C580 DIETHYLPHTHALATE	149 A VB	934 3	1179. 1.369	0.556
85	S3#135	C590 FLUORENE	166 A BB	927 3	5987. 1.298	2.980
86	S3#140	C585 4-CHLOROPHENYL-PHEN	204	3 NOT FOUND		
87	S3#145	5-NITRO-O-TOLUIDINE	152 A VB	943 3	2476. 0.356	4.499
88	S3#150	C595 4-NITROANALINE	138	3 NOT FOUND		
89	S4# 10	C610 4,6-DINITRO-2-METHY	198	4 NOT FOUND		
90	S4# 15	C615 N-NITROSODIPHENYLAM	169 A VB	965 4	5818. 0.551	4.145
91	S4# 20	C620 AZOBENZENE	77 A VB	966 4	643. 1.171	0.216
92	S4# 25	SYM-TRINITROBENZENE	75 A BV	1026 4	206. 0.162	0.501
93	S4# 30	C625 4-BROMOPHENYL-PHENY	248	4 NOT FOUND		
94	S4# 35	PHENACETIN	108	4 NOT FOUND		
95	S4# 37	DIALATE	234 A VB	1021 4	156. 0.090	0.678
96	S4# 40	C630 HEXACHLOROBENZENE	284	4 NOT FOUND		
97	S4# 45	4-AMINOBIIPHENYL	169 A BB	1077 4	3167. 0.609	2.040
98	S4# 50	C635 PENTACHLOROPHENOL	266 A VB	1064 4	261. 0.155	0.660
99	S4# 55	PRONAMIDE	173	4 NOT FOUND		
100	S4# 60	PENTACHLORONITROBENZENE	237	4 NOT FOUND		
101	S4# 65	C640 PHENANTHRENE	178 A BB	1092 4	18337. 1.033	6.966
102	S4# 70	C645 ANTHRACENE	178	4 NOT FOUND		
103	S4# 75	2SECBUTYL-4,6-DINITROPHE	211	4 NOT FOUND		
104	S4# 80	C647 CARBAZOLE	167	4 NOT FOUND		
105	S4# 85	C650 DI-N-BUTYLPHTHALATE	149	4 NOT FOUND		
106	S4#100	4-NITROQUINOLINE-1-OXIDE	190	4 NOT FOUND		
107	S4#105	METHAPYRILENE	58	4 NOT FOUND		
108	S4#106	ISODRIN	193	4 NOT FOUND		
109	S4#110	C655 FLUORANTHENE	202 A BB	1300 4	602. 0.900	0.263
110	S4#120	CHLOROBENZILATE	139	4 NOT FOUND		

111	S5# 10	C710 BENZIDINE	184	5	NOT FOUND		
112	S5# 15	C715 PYRENE	202	A BB	1336 5	2881.	1.323 0.857
113	S5# 20	ARAMITE (#1)	185	5	NOT FOUND		
114	S5# 25	ARAMITE (#2)	185	5	NOT FOUND		
115	S5# 30	P-DIMETHYLAMINOAZOBENZEN	120	5	NOT FOUND		
116	S5# 35	3,3'-DIMETHYLBENZIDINE	212	5	NOT FOUND		
117	S5# 37	KEPONE	272	5	NOT FOUND		
118	S5# 40	C720 BUTYLBENZYLPHTHALAT	149	5	NOT FOUND		
119	S5# 45	2-ACETYLAMINOFLUORENE	181	5	NOT FOUND		
120	S5# 50	C730 BENZO(A)ANTHRACENE	228	A BB	1556 5	2250.	1.125 0.787
121	S5# 55	C725 3,3'-DICHLOROBENZID	252	5	NOT FOUND		
122	S5# 60	C740 CHRYSENE	228	A BB	1556 5	2250.	1.014 0.873
123	S5# 65	C745 BIS(2-ETHYLHEXYL)PH	149	A BB	1604 5	6255.	0.980 2.511
124	S5# 85	3-METHYLCHOLANTHRENE	268	5	NOT FOUND		
125	S6# 10	C760 DI-N-OCTYL PHTHALAT	149	A VB	1717 6	1093.	1.925 0.237
126	S6# 15	C765 BENZO(B)FLUORANTHEN	252	A BV	1728 6	1158.	1.443 0.335
127	S6# 20	7,12-DIMETHYLBENZANTHRAC	256	6	NOT FOUND		
128	S6# 25	C770 BENZO(K)FLUORANTHEN	252	6	NOT FOUND		
129	S6# 30	HEXACHLOROPHENE	196	6	NOT FOUND		
130	S6# 35	C775 BENZO(A)PYRENE	252	A BB	1787 6	643.	1.182 0.227
131	S6# 55	C780 INDENO(1,2,3-CD)PYR	276	6	NOT FOUND		
132	S6# 60	C785 DIBENZ(A,H)ANTHRACE	278	6	NOT FOUND		
133	S6# 65	C790 BENZO(G,H,I)PERYLEN	276	6	NOT FOUND		

QUANTERRA Internal Standard Check

Standard Filename: ST16980831

Analyzed: 08/31/98 12:31

Standard	IS# 1			IS# 2			IS# 3		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	28649.		3.78	100432.		5.90	54347.		9.21
Upper Limit	57297.=+200		4.28	200864.=+200		6.40	108695.=+200		9.71
Lower Limit	14324.=+ 50		3.28	50216.=+ 50		5.40	27174.=+ 50		8.71
Filename									
1 30068106	33301.= 116		3.79	120808.= 120		5.91	61921.= 113		9.24

IS# 1 = C130 1,4-DICHLOROBENZENE-D4

IS# 2 = C140 NAPHTHALENE-D8

IS# 3 = C150 ACENAPHTHENE-D10

* - indicates an I.S. is outside QC limit(s)

Standard	IS# 4			IS# 5			IS# 6		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	98520.		12.08	62289.		17.26	50750.		19.83
Upper Limit	197039.=+200		12.58	124578.=+200		17.76	101500.=+200		20.33
Lower Limit	49260.=+ 50		11.58	31145.=+ 50		16.76	25375.=+ 50		19.33
Filename									
1 30068106	101964.= 103		12.10	101687.= 163		17.27	95704.= 188		19.84

IS# 4 = C160 PHENANTHRENE-D10

IS# 5 = C170 CHRYSENE-D12

IS# 6 = C175 PERYLENE-D12

* - indicates an I.S. is outside QC limit(s)

TARGET COMPOUND COMPARISON

COMPOUND: C315 PHENOL

94 S1#105CC 108-95-2

RAW DATA: 30068106 #311

08/31/98 21:30

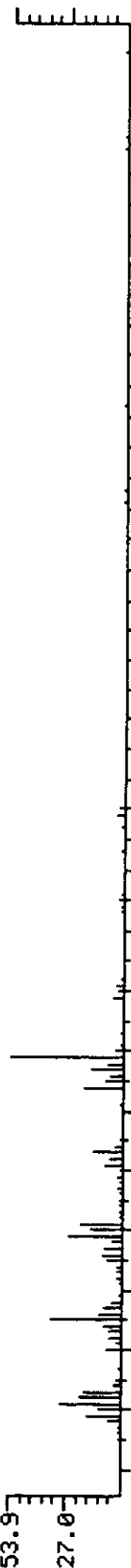
SAMPLE: T-MM5-3 1/35A/IML

COND.: UG/ML *1ML *100%/100% *(NA/NA) /1/35A NA M

BASE M/Z: 99

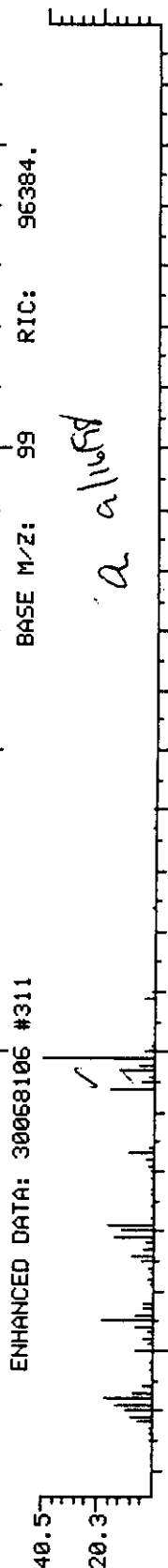
RIC: 170752.

17568.



STANDARD FILE: ST16980831 #309

13200.

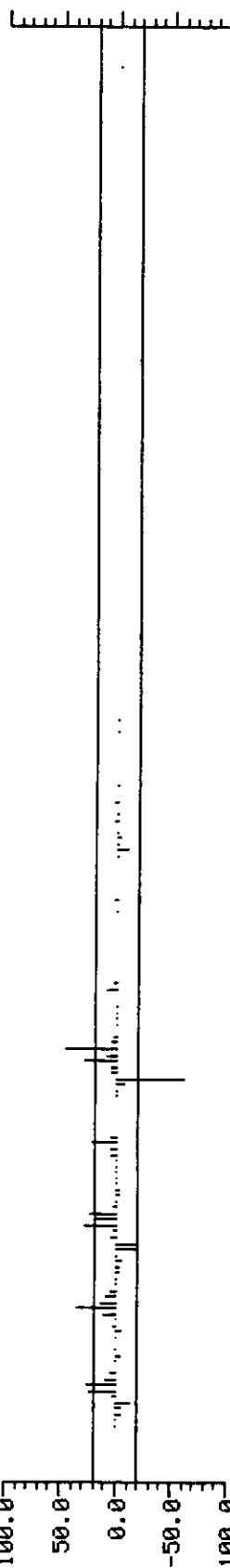
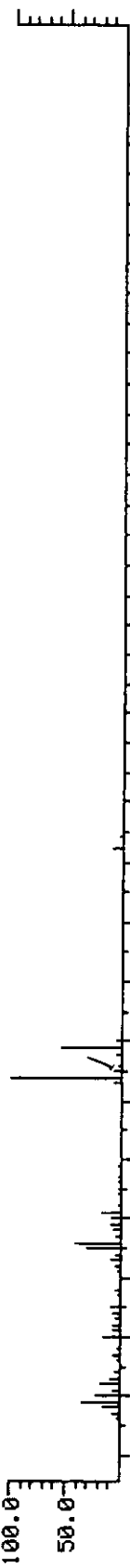


08/31/98 12:31

BASE M/Z: 94

RIC: 170752.

32576.



OUT

TARGET COMPOUND COMPARISON

COMPOUND: C450 NAPHTHALENE 128 S2# 60 91-20-3
 BASE M/Z: 128 RIC: 139264.

RAW DATA: 30068106 #534

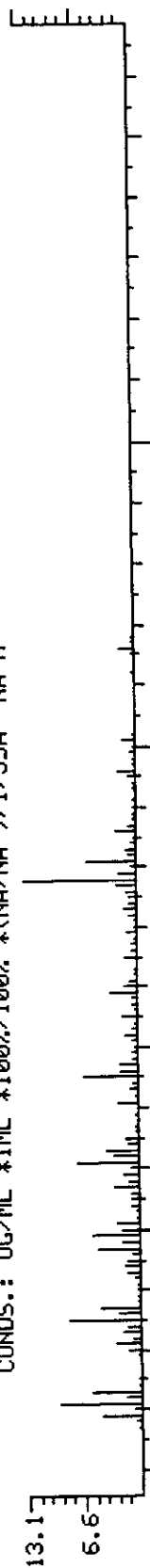
08/31/98 21:30

SAMPLE: T-MM5-3 1/35A/1ML

INST. ID: F16

CONDS.: UG/ML *1ML *100%/100% *(NA/NA) /1/35A NA M

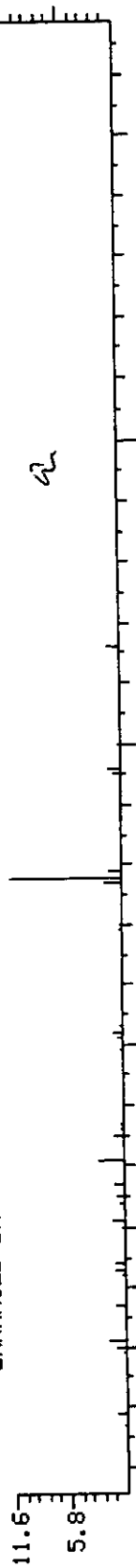
10384.



ENHANCED DATA: 30068106 #534

BASE M/Z: 128 RIC: 28736.

9152.

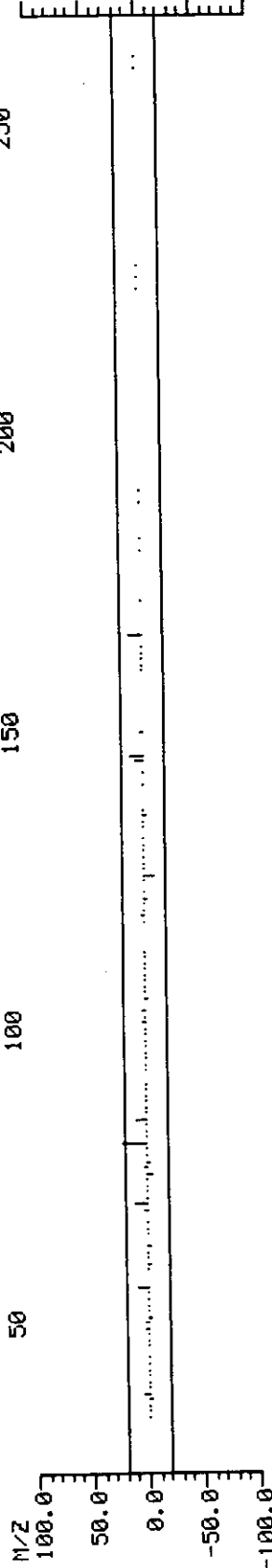
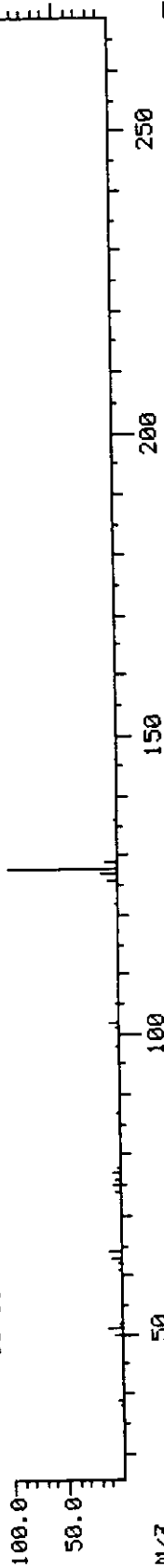


STANDARD FILE: ST16980831 #533

08/31/98 12:31

BASE M/Z: 128 RIC: 195840.

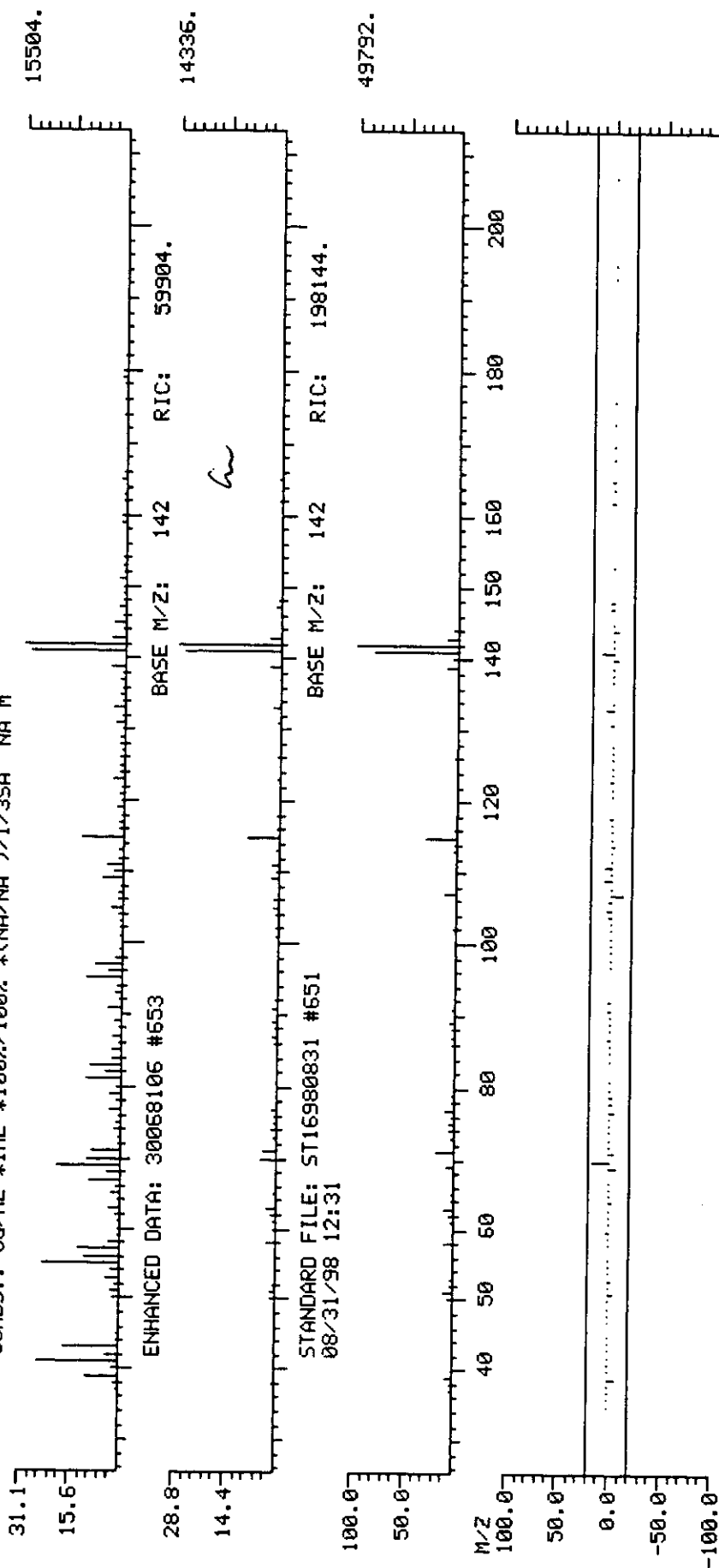
79232.



TARGET COMPOUND COMPARISON

COMPOUND: C470 2-METHYLNAPHTHALENE 142 52#145 91-57-6

RAW DATA: 30068106 #653 BASE M/Z: 142 RIC: 218368.
 08/31/98 21:30
 SAMPLE: T-MM5-3 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA) >1/35A NA M



Data Reduced by: *9/17/98* Date: *9/17/98*
Data Reviewed by: *9/17/98* Date: *9/17/98*

Data File: 30068106

QUANTERRA GC/MS TIC REPORT (Part 1)

Sample: T-MM5-3 1/3SA/1ML
Analyst: DAT

INST. ID: F16

Date Analyzed: 08/31/98 21:30

Run Factor: 3.00

Concentration
in Sample

SCAN (UG/A) CAS

100 170 350. 123-42-2

2-PENTANONE, 4-HYDROXY-4-METHYL-

or isomer

2 221 390. 111-84-2

NONANE

or isomer

3 311 270. 00-00-0

UNKNOWN

5700

4 445 180. 1120-21-4

UNDECANE

or isomer

5 557 160. 112-40-3

DODECANE

or isomer

6 634 240. ~~54105-67-8~~

~~HEPTADECANE, 2,6-DIMETHYL-~~

alkane

5746400

7 653 170. 2471-83-2

1H-INDENE, 1-ETHYLIDENE-

or isomer

8 664 240. 17301-32-5

UNDECANE, 4,7-DIMETHYL-

or isomer

9 724 150. 321-60-8

1,1'-BIPHENYL, 2-FLUORO-

or isomer

surrogate

10 739 250. 18344-37-1

HEPTADECANE, 2,6,10,14-TETRAMETHYL-

or isomer

11 765 430. 629-59-4

TETRADECANE *isomer*

12 778 230. ~~00-00-0~~

5740201000

~~UNKNOWN~~

PAH

13 822 300. 18344-37-1

HEPTADECANE, 2,6,10,14-TETRAMETHYL- *or isomer*

14 860 440. 629-62-9

PENTADECANE *or isomer*

~~15~~ 905 120. 52783-43-4

~~NONADECANOL~~ *or isomer*

16 949 400. 629-92-5

NONADECANE

17 989 230. 54105-67-8

HEPTADECANE, 2,6-DIMETHYL-

18 1037 520. 54105-67-8

HEPTADECANE, 2,6-DIMETHYL-

19 1115 170. 629-92-5

NONADECANE *or isomer*

20 1120 180. 629-92-5

NONADECANE

~~21~~ 1192 140. 629-92-5

~~NONADECANE~~

~~22~~ 1336 130.

~~2-METHYLOCTADECANE~~

23 1531 150. 629-92-5

NONADECANE

or isomer

24 1587 240.

74685-29-3

9-EICOSENE, (E)-

~~25~~ 1809 120.

00-00-0

UNKNOWN

DATA FILE: 30068106

QUANTERRA GC/MS TIC REPORT (Part 2)

CONCENTRATION = AREA(TIC)*CONC(IS)/AREA(IS)

				INT.		AMOUNT						
						AS ANALYZED						
#	FIT	PURITY	STD.	RT	RRT	AREA	HEIGHT	(UG/ML)	LIB	LIB #		
1	948	906	2	1:53	0.205	1066940.	387584.	117.319	NB	3241.		
2	986	942	2	2:27	0.266	1174820.	346624.	129.180	NB	5159.		
3	829	529	2	3:27	0.375	827993.	123904.	91.044	UK	1.		
4	975	722	2	4:56	0.536	541092.	131657.	59.497	NB	11607.		
5	970	895	2	6:11	0.671	494208.	143104.	54.342	NB	15343.		
6	938	687	2	7:03	0.764	724224.	159488.	79.634	NB	37462.		
7	956	760	2	7:15	0.787	513280.	72192.	56.439	NB	8109.		
8	939	829	2	7:23	0.800	715264.	174336.	78.649	NB	19016.		
9	994	843	2	8:03	0.872	444928.	87552.	48.923	NB	15856.		
10	930	812	2	8:13	0.890	768256.	166144.	84.476	NB	42196.		
11	960	744	2	8:30	0.922	1312510.	320000.	144.321	NB	22530.		
12	910	451	2	8:39	0.937	701952.	90624.	77.185	UK	1.		
13	959	792	2	9:08	0.990	914944.	165632.	100.605	NB	42196.		
14	963	816	2	9:33	1.036	1321220.	301056.	145.278	NB	25997.		
15	934	653	2	10:03	1.090	375040.	81152.	41.239	NB	40233.		
16	952	842	2	10:33	1.143	1198340.	266752.	131.766	NB	37465.		
17	955	834	2	11:00	1.192	708891.	168771.	77.948	NB	37462.		
18	974	872	2	11:32	1.249	1565700.	264192.	172.160	NB	37462.		
19	947	806	2	12:24	1.343	528896.	153600.	58.156	NB	37465.		
20	952	822	2	12:27	1.349	551424.	169472.	60.633	NB	37465.		
21	950	774	4	13:15	0.768	357120.	115968.	45.964	NB	37465.		
22	908	662	4	14:51	0.860	335488.	76717.	43.180	NB	37456.		
23	935	663	4	17:01	0.986	392960.	48384.	50.577	NB	37465.		
24	984	795	4	17:39	1.022	617728.	115968.	79.506	NB	39515.		
25	706	329	5	20:07	1.014	281792.	23360.	41.621	UK	1.		

Library Search Data: 30068106 # 170 Base m/z: 43
 08/31/98 21:30:00 + 1:53 Cali: 30068106 # 3 RIC: 273920.
 Sample: T-MMS-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 495 matched at least 6 of the 16 largest peaks in the unknown

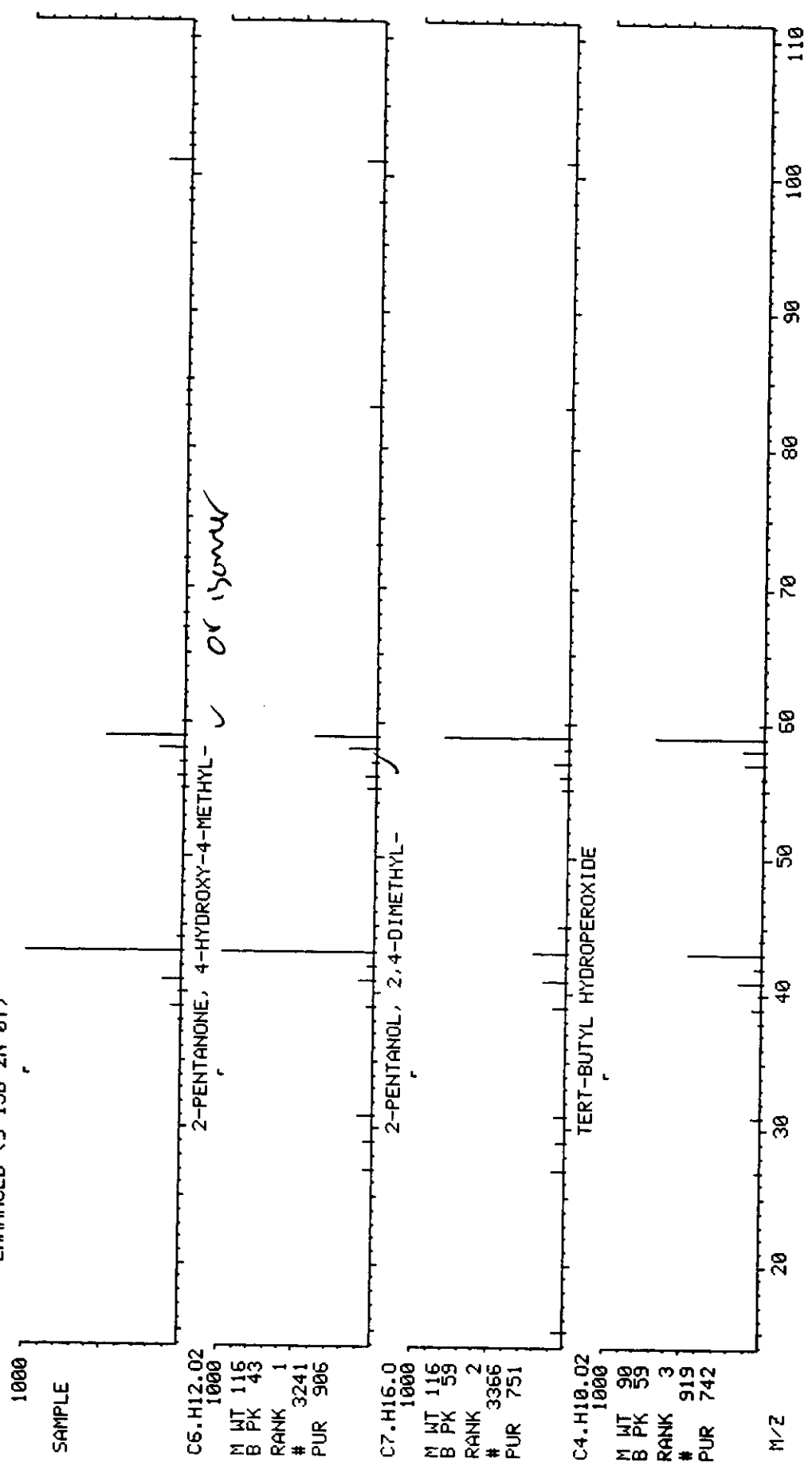
Rank In. Name
 1 3241 2-PENTANONE, 4-HYDROXY-4-METHYL-
 2 3366 2-PENTANOL, 2,4-DIMETHYL-
 3 919 TERT-BUTYL HYDROPEROXIDE
 4 5791 1,3-DIOXOLANE-2-METHANOL, 2,4-DIMETHYL-
 5 3288 ACETIC ACID, 1,1-DIMETHYLETHYL ESTER
 6 5594 2-PROPANOL, 2-NITROSO-, ACETATE (ESTER)
 7 3349 2-HEXANOL, 2-METHYL-
 8 1576 OXIRANE, TETRAMETHYL-
 9 3353 2-PENTANOL, 2,3-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C6.H12.O2	116	43	906	948	928
2	C7.H16.O	116	59	751	778	769
3	C4.H10.O2	90	59	742	894	761
4	C6.H12.O3	132	43	734	805	800
5	C6.H12.O2	116	43	729	779	742
6	C5.H9.O3.N	131	43	713	830	760
7	C7.H16.O	116	59	698	736	750
8	C6.H12.O	100	59	653	805	705
9	C7.H16.O	116	59	643	689	706

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	123-42-2
2	—	—	—	—	625-06-9
3	—	—	—	—	75-91-2
4	—	—	—	—	53951-43-2
5	—	—	—	—	540-88-5
6	—	—	—	—	6931-04-0
7	—	—	—	—	625-23-0
8	—	—	—	—	5076-20-0
9	—	—	—	—	4911-70-0

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:30:00 + 1:53
 SAMPLE: T-MM5-3 1/35A/1ML
 CONDS.: UG/ML *1ML *1002/1002 *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068106 # 170
 CALI: 30068106 # 3
 BASE M/Z: 43
 RIC: 273920.



Library Search Data: 30068106 # 221 Base m/z: 43
 08/31/98 21:30:00 + 2:27 Cali: 30068106 # 3 RIC: 275968.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (\$ 158 2N 0T)

62231 spectra in LIBRARYN8 searched for maximum PURITY
 921 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 5159 NONANE
 2 5151 HEPTANE, 3,4-DIMETHYL-
 3 3086 HEXANE, 2,4-DIMETHYL-
 4 5141 HEPTANE, 2,4-DIMETHYL-
 5 19015 DECANE, 2,5,6-TRIMETHYL-
 6 5144 HEPTANE, 4-ETHYL-
 7 8085 HEPTANE, 2,4,6-TRIMETHYL-
 8 11607 UNDECANE
 9 5154 HEXANE, 4-ETHYL-2-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C9.H20	128	43	942	986	942
2	C9.H20	128	43	905	928	905
3	C8.H18	114	43	897	962	908
4	C9.H20	128	43	877	926	890
5	C13.H28	184	57	871	907	891
6	C9.H20	128	43	869	912	869
7	C10.H22	142	43	865	903	881
8	C11.H24	156	43	863	918	919
9	C9.H20	128	57	862	909	862

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	111-84-2
2	—	—	—	—	922-28-1
3	—	—	—	—	589-43-5
4	—	—	—	—	2213-23-2
5	—	—	—	—	62108-23-0
6	—	—	—	—	2216-32-2
7	—	—	—	—	2613-61-8
8	—	—	—	—	1120-21-4
9	—	—	—	—	3074-75-7

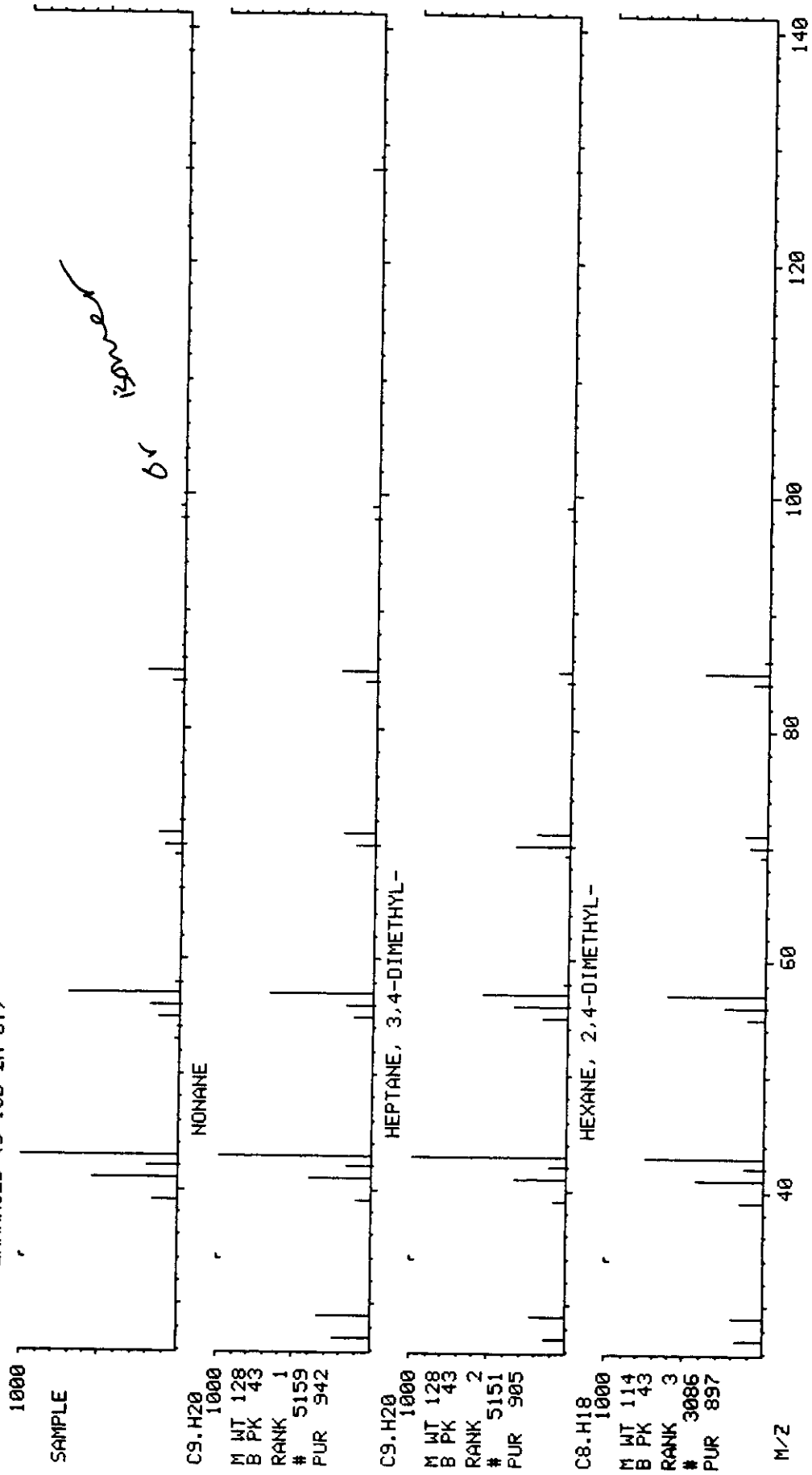
MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 21:30:00 + 2:27

SAMPLE: T-MM5-3 1/35A/1ML

CONDS.: UG/ML *1ML *1002/1002 *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)

DATA: 30058106 # 221
CALI: 30058106 # 3
BASE M/Z: 43
RIC: 275968.



Library Search Data: 30068106 # 311 Base m/z: 99
 08/31/98 21:30:00 + 3:28 Cali: 30068106 # 3 RIC: 84864.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 346 matched at least 6 of the 16 largest peaks in the unknown

Rank	In.	Name
1	15124	2,5-FURANDIONE, 3-(1,1-DIMETHYLPROPYL)DIHYDRO-
2	1394	2-PIPERIDINONE
3	18863	CIS-4-HYDROXY-3-METHYLDECAHOIC ACID LACTONE
4	18864	TRANS-4-HYDROXY-3-METHYLDECAHOIC ACID LACTONE
5	1501	PHENOL-D6-
6	11410	5-HYDROXYNONANOIC ACID LACTONE
7	22383	CIS-4-HYDROXY-3-METHYLUNDECAHOIC ACID LACTONE
8	1399	1-PYRROLIDINECARBOXALDEHYDE
9	4906	2H-PYRAN-2-ONE, 6-ETHYLTETRAHYDRO-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C9.H14.O3	170	99	529	829	553
2	C5.H9.O.N	99	30	527	870	544
3	C11.H20.O2	184	99	517	819	587
4	C11.H20.O2	184	99	514	814	598
5	C6.O.D6	100	99	512	879	549
6	C9.H16.O2	156	99	501	856	555
7	C12.H22.O2	198	99	495	856	548
8	C5.H9.O.N	99	99	494	868	514
9	C7.H12.O2	128	42	478	871	525

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	56666-76-3
2	—	—	—	—	675-20-7
3	—	—	—	—	- -
4	—	—	—	—	- -
5	—	—	—	—	13127-88-3
6	—	—	—	—	- -
7	—	—	—	—	- -
8	—	—	—	—	3760-54-1
9	—	—	—	—	3301-90-4

MID LIBRARY SEARCH <LIBRARYNB>

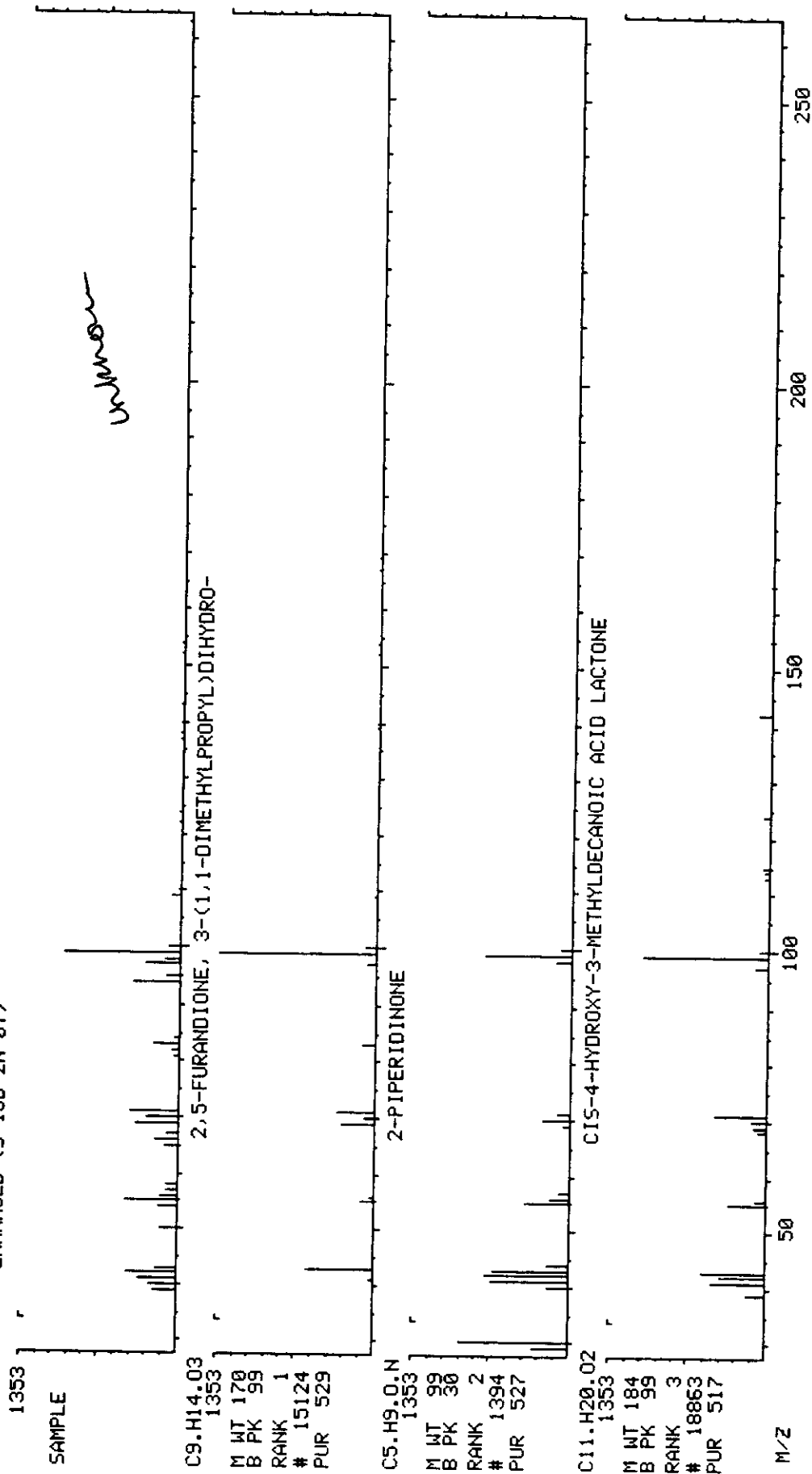
08/31/98 21:30:00 + 3:28

SAMPLE: T-MM5-3 1/35A/1ML

COND.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)

DATA: 30068105 # 311
CALI: 30068105 # 3

BASE M/Z: 99
RIC: 84864.



Library Search Data: 30068106 # 445 Base m/z: 43
 08/31/98 21:30:00 + 4:57 Cali: 30068106 # 3 RIC: 110592.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

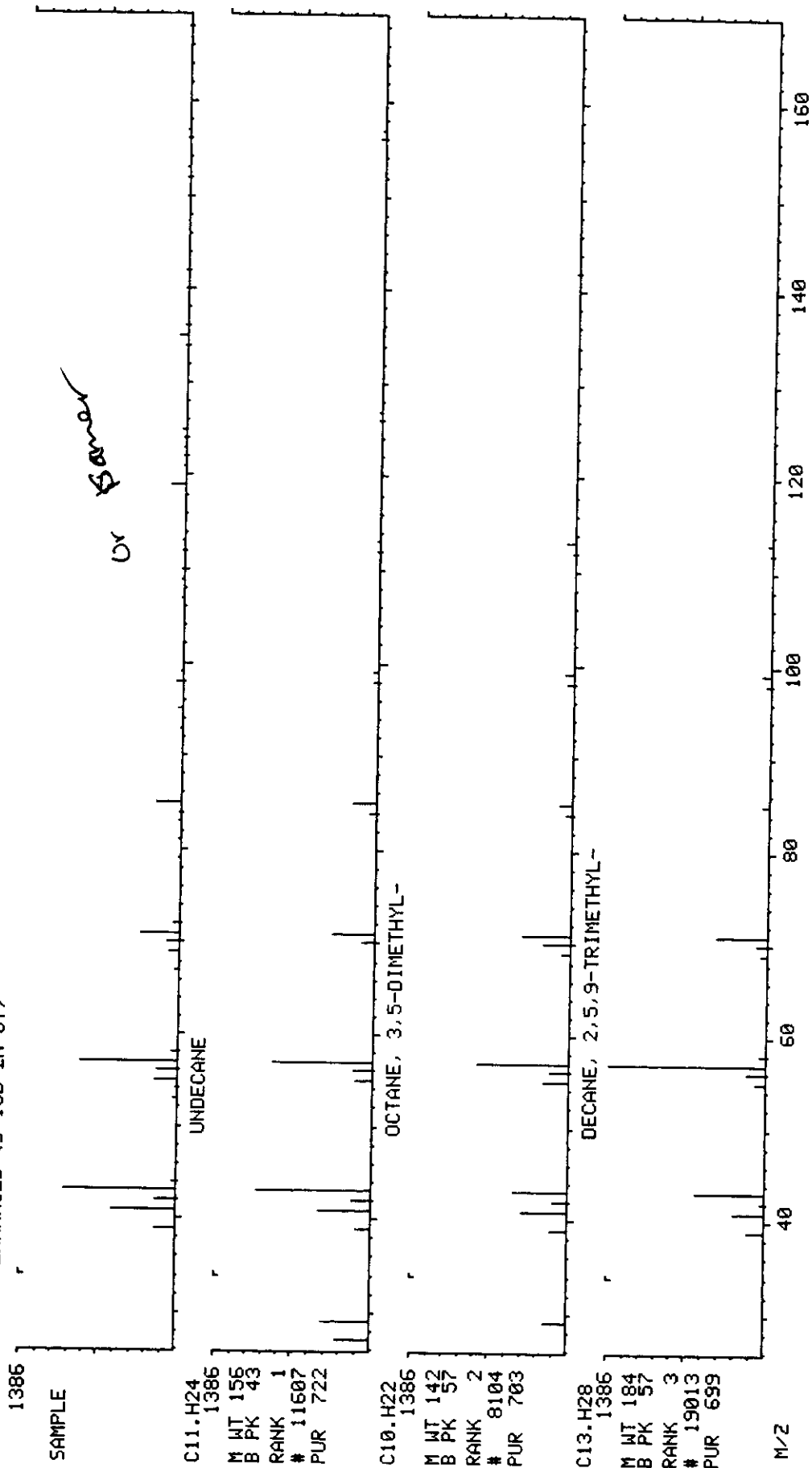
62231 spectra in LIBRARYNB searched for maximum PURITY
 578 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 11607 UNDECANE
 2 8104 OCTANE, 3,5-DIMETHYL-
 3 19013 DECANE, 2,5,9-TRIMETHYL-
 4 19015 DECANE, 2,5,6-TRIMETHYL-
 5 5151 HEPTANE, 3,4-DIMETHYL-
 6 5154 HEXANE, 4-ETHYL-2-METHYL-
 7 8089 NONANE, 2-METHYL-
 8 11602 OCTANE, 2,4,6-TRIMETHYL-
 9 5160 PENTANE, 2,2,3,4-TETRAMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C11.H24	156	43	722	975	730
2	C10.H22	142	57	703	936	703
3	C13.H28	184	57	699	923	699
4	C13.H28	184	57	695	913	702
5	C9.H20	128	43	688	916	690
6	C9.H20	128	57	686	920	690
7	C10.H22	142	43	680	924	710
8	C11.H24	156	57	680	939	691
9	C9.H20	128	57	677	919	677

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	1120-21-4
2	—	—	—	—	15869-93-9
3	—	—	—	—	62108-22-9
4	—	—	—	—	62108-23-0
5	—	—	—	—	922-28-1
6	—	—	—	—	3074-75-7
7	—	—	—	—	871-83-0
8	—	—	—	—	62016-37-9
9	—	—	—	—	1186-53-4

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:30:00 + 4:57
 SAMPLE: T-MM5-3 1/35A/1ML
 CONDS.: UG/ML *1ML *1002/1002 *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30058106 # 445
 CALI: 30058106 # 3
 BASE M/Z: 43
 RIC: 110592.



Library Search Data: 30068106 # 557 Base m/z: 43
08/31/98 21:30:00 + 6:12 Cali: 30068106 # 3 RIC: 132352.
Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
990 matched at least 7 of the 16 largest peaks in the unknown

Rank	In.	Name
1	15343	DODECANE
2	11607	UNDECANE
3	19016	UNDECANE, 4,7-DIMETHYL-
4	15353	2,6-DIMETHYLDECANE
5	8104	OCTANE, 3,5-DIMETHYL-
6	19015	DECANE, 2,5,6-TRIMETHYL-
7	22530	TETRADECANE
8	19028	DECANE, 2,4,6-TRIMETHYL-
9	11602	OCTANE, 2,4,6-TRIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H26	170	57	895	970	895
2	C11.H24	156	43	874	966	901
3	C13.H28	184	43	871	962	872
4	C12.H26	170	43	862	923	869
5	C10.H22	142	57	860	946	860
6	C13.H28	184	57	860	933	861
7	C14.H30	198	43	855	924	912
8	C13.H28	184	43	855	941	861
9	C11.H24	156	57	855	965	855

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	112-40-3
2	—	—	—	—	1120-21-4
3	—	—	—	—	17301-32-5
4	—	—	—	—	13150-81-7
5	—	—	—	—	15869-93-9
6	—	—	—	—	62108-23-0
7	—	—	—	—	629-59-4
8	—	—	—	—	62108-27-4
9	—	—	—	—	62016-37-9

MID LIBRARY SEARCH (LIBRARYNB)

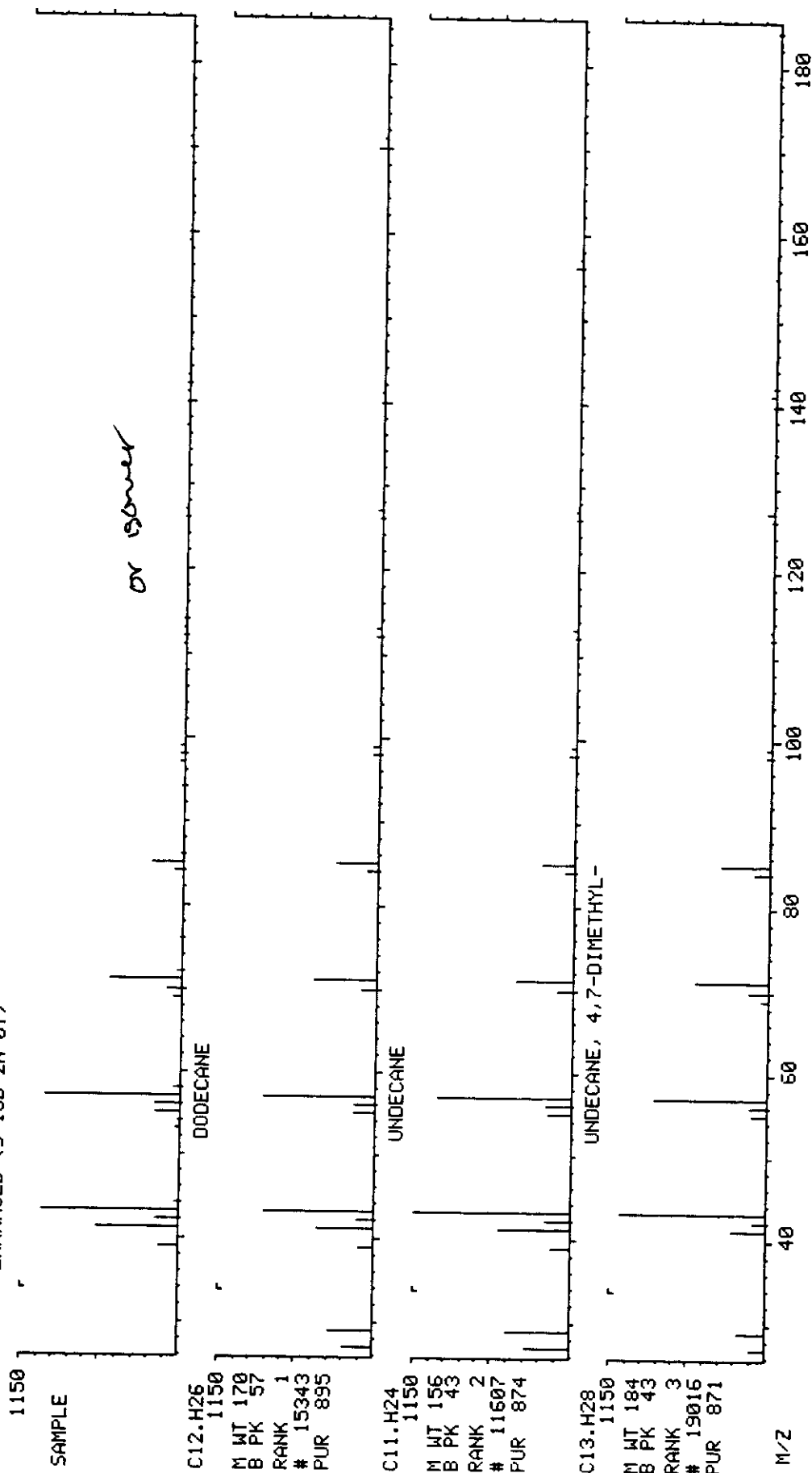
08/31/98 21:30:00 + 6:12

SAMPLE: T-MMS-3 1/35A/1ML

CONDS.: UG/ML *1ML *1002/1002 *(NA/NA) /1/35A NA M
ENHANCED (S 15B 2N 0T)

DATA: 30068106 # 557
CALI: 30068106 # 3

BASE M/Z: 43
RIC: 132352.



Library Search Data: 30068106 # 634 Base m/z: 57
 08/31/98 21:30:00 + 7:03 Cali: 30068106 # 3 RIC: 136960.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 748 matched at least 7 of the 16 largest peaks in the unknown

Rank	In.	Name
1	37462	HEPTADECANE, 2,6-DIMETHYL-
2	14799	1-DECENE, 3,4-DIMETHYL-
3	19523	1-DECANOL, 2-ETHYL-
4	22534	TRIDECANE, 7-METHYL-
5	15353	2,6-DIMETHYLDECANE
6	26408	2-ETHYL-1-DODECANOL
7	42196	HEPTADECANE, 2,6,10,14-TETRAMETHYL-
8	26409	2-HEXYL-1-OCTANOL
9	32418	2-HEXYL-1-DECANOL

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	687	938	702
2	C12.H24	168	57	686	940	686
3	C12.H26.O	186	57	679	928	692
4	C14.H30	198	57	667	913	681
5	C12.H26	170	43	659	918	664
6	C14.H30.O	214	57	656	899	675
7	C21.H44	296	57	654	910	675
8	C14.H30.O	214	57	653	899	683
9	C16.H34.O	242	57	652	897	680

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	50871-03-9
3	—	—	—	—	21078-65-9
4	—	—	—	—	26730-14-3
5	—	—	—	—	13150-81-7
6	—	—	—	—	- -
7	—	—	—	—	18344-37-1
8	—	—	—	—	- -
9	—	—	—	—	- -

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 21:30:00 + 7:03

SAMPLE: T-MM5-3 1/35A/1ML

COND.: UG/ML *1ML *100%/100% *(NA/NA >1/35A

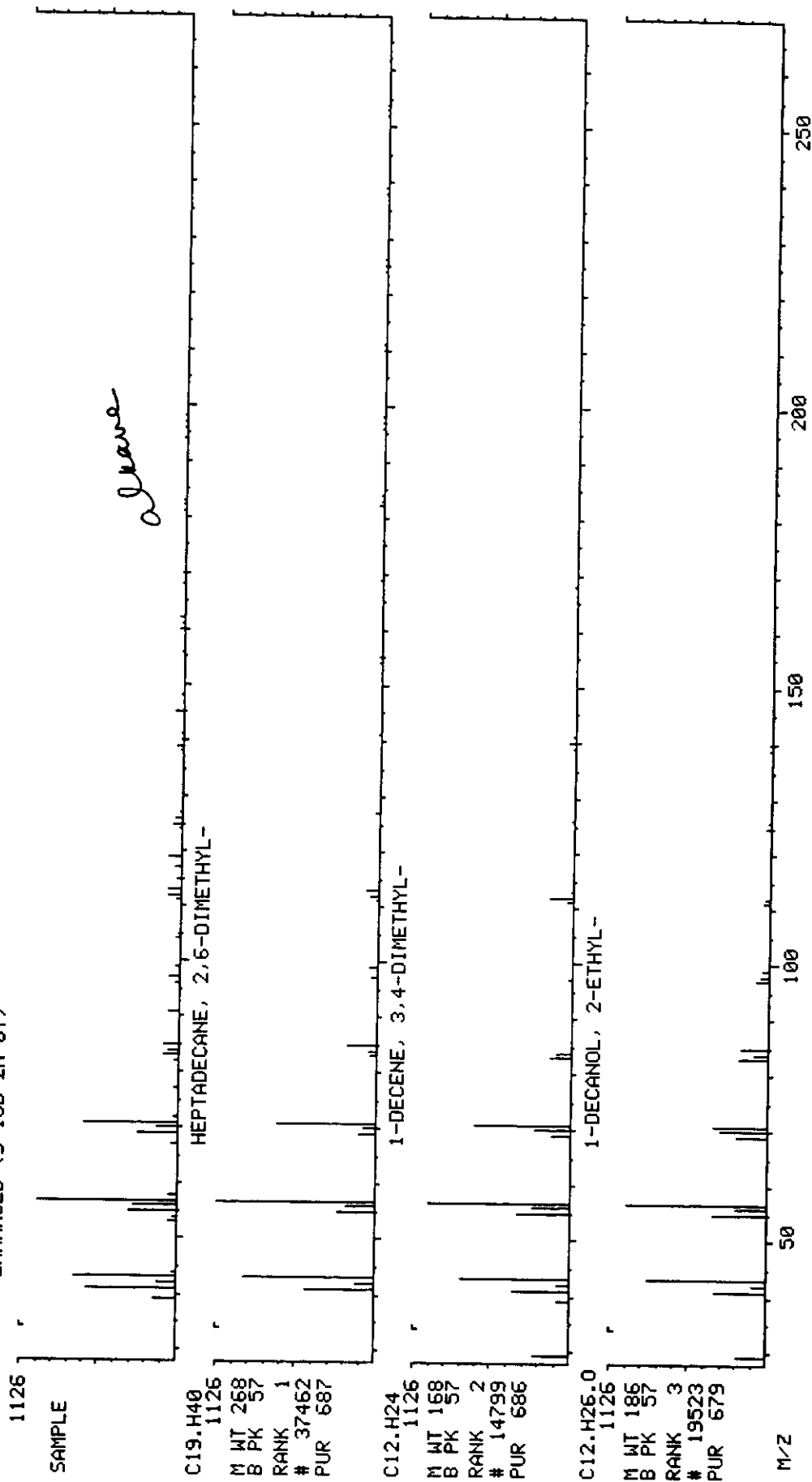
ENHANCED (S 15B 2N ØT)

DATA: 30068106 # 534
CALI: 30068106 # 3

CALI: 30068106

BASE M/Z: 57
RIC: 136960.

PHF WZ: 57
RIC: 136960.



1058

Library Search Data: 30068106 # 653 Base m/z: 142
 08/31/98 21:30:00 + 7:16 Cali: 30068106 # 3 RIC: 57920.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 298 matched at least 4 of the 16 largest peaks in the unknown

Rank In. Name
 1 8109 1H-INDENE, 1-ETHYLIDENE-
 2 8107 NAPHTHALENE, 1-METHYL-
 3 8108 1,4-METHANONAPHTHALENE, 1,4-DIHYDRO-
 4 8111 NAPHTHALENE, 2-METHYL-
 5 8110 BENZOCYCLOHEPTATRIENE
 6 12521 2-CYCLOPENTEN-1-OL, 1-PHENYL-
 7 7992 BENZENEACETONITRILE, 4-CYANO-
 8 15865 1-NAPHTHALENEETHANOL
 9 15368 BICYCLO[2.2.1]HEPT-2-ENE, 2-PHENYL-

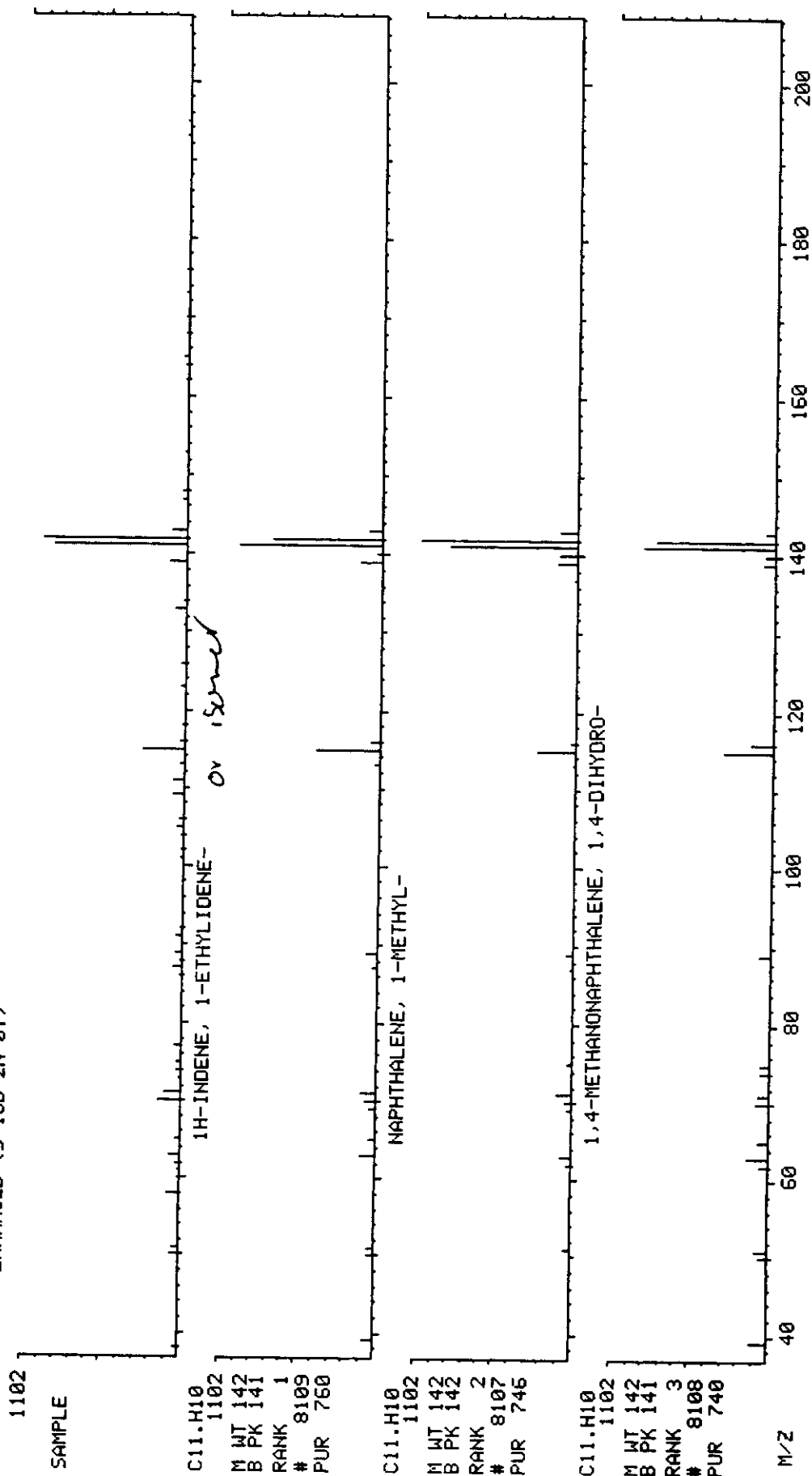
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C11.H10	142	141	760	956	779
2	C11.H10	142	142	746	938	787
3	C11.H10	142	141	740	932	767
4	C11.H10	142	142	728	926	779
5	C11.H10	142	141	706	931	721
6	C11.H12.O	160	142	577	758	743
7	C9.H6.N2	142	142	574	790	622
8	C12.H12.O	172	141	563	733	677
9	C13.H14	170	142	562	738	622

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	2471-83-2
2	—	—	—	—	90-12-0
3	—	—	—	—	4453-90-1
4	—	—	—	—	91-57-6
5	—	—	—	—	264-09-5
6	—	—	—	—	56667-10-8
7	—	—	—	—	876-31-3
8	—	—	—	—	773-99-9
9	—	—	—	—	4237-08-5

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:30:00 + 7:16
 SAMPLE: T-MM5-3 1/35A/1ML
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068106 # 653
 CALI: 30068106 # 3

BASE M/Z: 142
 RIC: 57920.



1056

Library Search Data: 30068106 # 664 Base m/z: 57
 08/31/98 21:30:00 + 7:23 Cali: 30068106 # 3 RIC: 165888.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 324 matched at least 8 of the 16 largest peaks in the unknown

Rank	In.	Name
1	19016	UNDECANE, 4,7-DIMETHYL-
2	8104	OCTANE, 3,5-DIMETHYL-
3	19015	DECANE, 2,5,6-TRIMETHYL-
4	19026	DECANE, 2,6,8-TRIMETHYL-
5	11602	OCTANE, 2,4,6-TRIMETHYL-
6	19013	DECANE, 2,5,9-TRIMETHYL-
7	18985	TRIDECANE
8	8089	NONANE, 2-METHYL-
9	8102	HEXANE, 2,2,3,3-TETRAMETHYL-

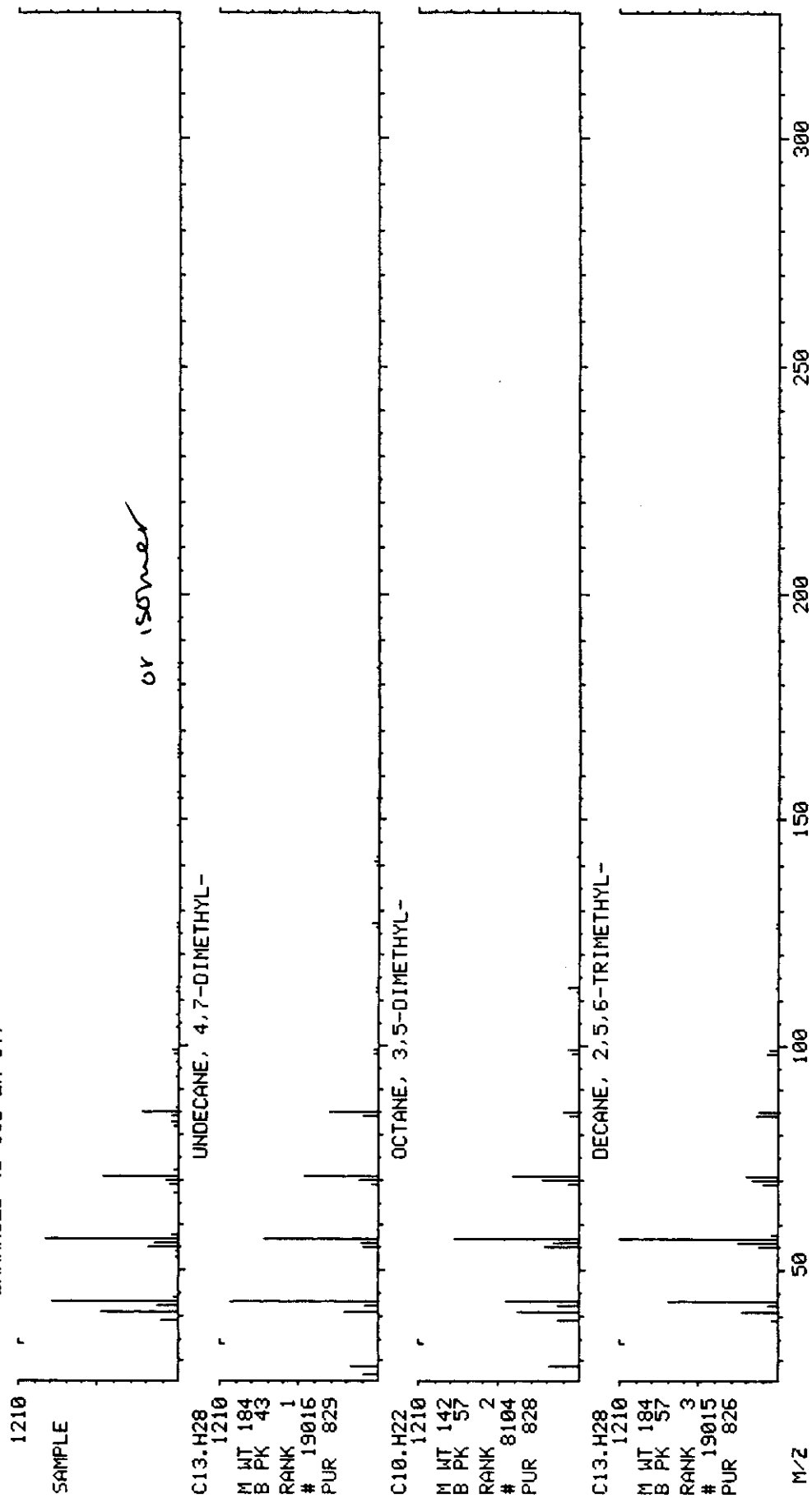
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H28	184	43	829	939	845
2	C10.H22	142	57	828	941	836
3	C13.H28	184	57	826	931	826
4	C13.H28	184	57	823	940	827
5	C11.H24	156	57	823	965	825
6	C13.H28	184	57	816	935	816
7	C13.H28	184	57	815	948	831
8	C10.H22	142	43	798	934	808
9	C10.H22	142	57	795	918	795

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	17301-32-5
2	—	—	—	—	15869-93-9
3	—	—	—	—	62108-23-0
4	—	—	—	—	62108-26-3
5	—	—	—	—	62016-37-9
6	—	—	—	—	62108-22-9
7	—	—	—	—	629-50-5
8	—	—	—	—	871-83-0
9	—	—	—	—	13475-81-5

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:30:00 + 7:23
 SAMPLE: T-MM5-3 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *1ML *1002/1002 *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30058106 # 664
 CALI: 30058106 # 3

BASE M/Z: 57
 RIC: 165888.



Library Search Data: 30068106 # 724 Base m/z: 172
 08/31/98 21:30:00 + 8:03 Cali: 30068106 # 3 RIC: 82816.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 224 matched at least 4 of the 16 largest peaks in the unknown

Rank In. Name
 1 15856 1,1'-BIPHENYL, 2-FLUORO-
 2 15857 1,1'-BIPHENYL, 4-FLUORO-
 3 15721 4-(2-HYDROXYPHENYL)PYRIMIDINE
 4 15828 1,8-NAPHTHYRIDINE, 2,4,7-TRIMETHYL-
 5 15722 2-(2-HYDROXYPHENYL)PYRIMIDINE
 6 15004 BENZENE, 1-BROMO-4-METHYL-
 7 15719 5-(4-HYDROXYPHENYL)PYRIMIDINE
 8 39206 1-(2-(2-METHYL-5-PYRIDYL)ETHYL)-2,3,5-TRIMETHYLINDOLE
 9 14964 IMIDAZOLE, 4-CHLORO-2-TRIFLUOROMETHYL-

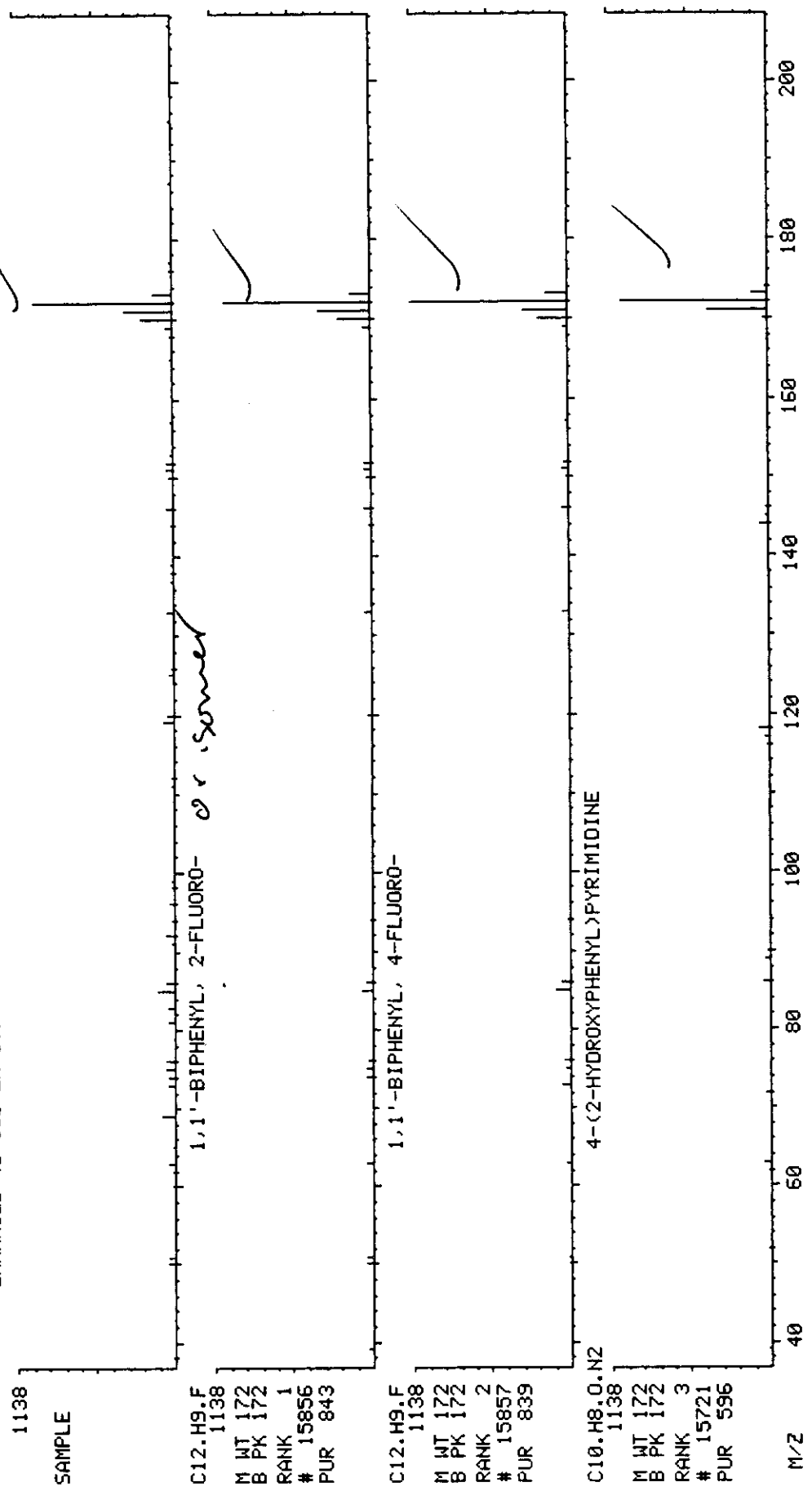
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H9.F	172	172	843	994	845
2	C12.H9.F	172	172	839	989	839
3	C10.H8.O.N2	172	172	596	854	652
4	C11.H12.N2	172	172	507	885	547
5	C10.H8.O.N2	172	172	474	755	529
6	C7.H7.BR	170	91	431	596	474
7	C10.H8.O.N2	172	172	426	684	555
8	C19.H22.N2	278	172	379	571	535
9	C4.H2.N2.CL.F3	170	170	378	542	451

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	321-60-8
2	—	—	—	—	324-74-3
3	—	—	—	—	68535-55-7
4	—	—	—	—	14757-44-9
5	—	—	—	—	64435-20-7
6	—	—	—	—	106-38-7
7	—	—	—	—	- -
8	—	—	—	—	- -
9	—	—	—	—	- -

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:30:00 + 8:03
 SAMPLE: T-MM5-3 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (5 158 2N 0T)

DATA: 30068105 # 724
 CALL: 30068105 # 3

BASE M/Z: 172
 RIC: 82816.



Library Search Data: 30068106 # 739 Base m/z: 57
 08/31/98 21:30:00 + 8:13 Cali: 30068106 # 3 RIC: 158720.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 712 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 2 25991 DODECANE, 2,6,10-TRIMETHYL-
 3 25997 PENTADECANE
 4 22535 DODECANE, 4,6-DIMETHYL-
 5 37462 HEPTADECANE, 2,6-DIMETHYL-
 6 19007 DODECANE, 3-METHYL-
 7 26001 DODECANE, 2,7,10-TRIMETHYL-
 8 22530 TETRADECANE
 9 18998 UNDECANE, 3,7-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C21.H44	296	57	812	930	835
2	C15.H32	212	57	805	959	810
3	C15.H32	212	57	800	936	815
4	C14.H30	198	57	796	940	800
5	C19.H40	268	57	791	924	824
6	C13.H28	184	57	787	931	788
7	C15.H32	212	57	785	973	795
8	C14.H30	198	43	782	912	799
9	C13.H28	184	43	777	924	778

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	18344-37-1
2	—	—	—	—	3891-98-3
3	—	—	—	—	629-62-9
4	—	—	—	—	61141-72-8
5	—	—	—	—	54105-67-8
6	—	—	—	—	17312-57-1
7	—	—	—	—	74645-98-0
8	—	—	—	—	629-59-4
9	—	—	—	—	17301-29-0

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 21:30:00 + 8:13

SAMPLE: T-MM5-3 1/35A/1ML

CONDOS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M

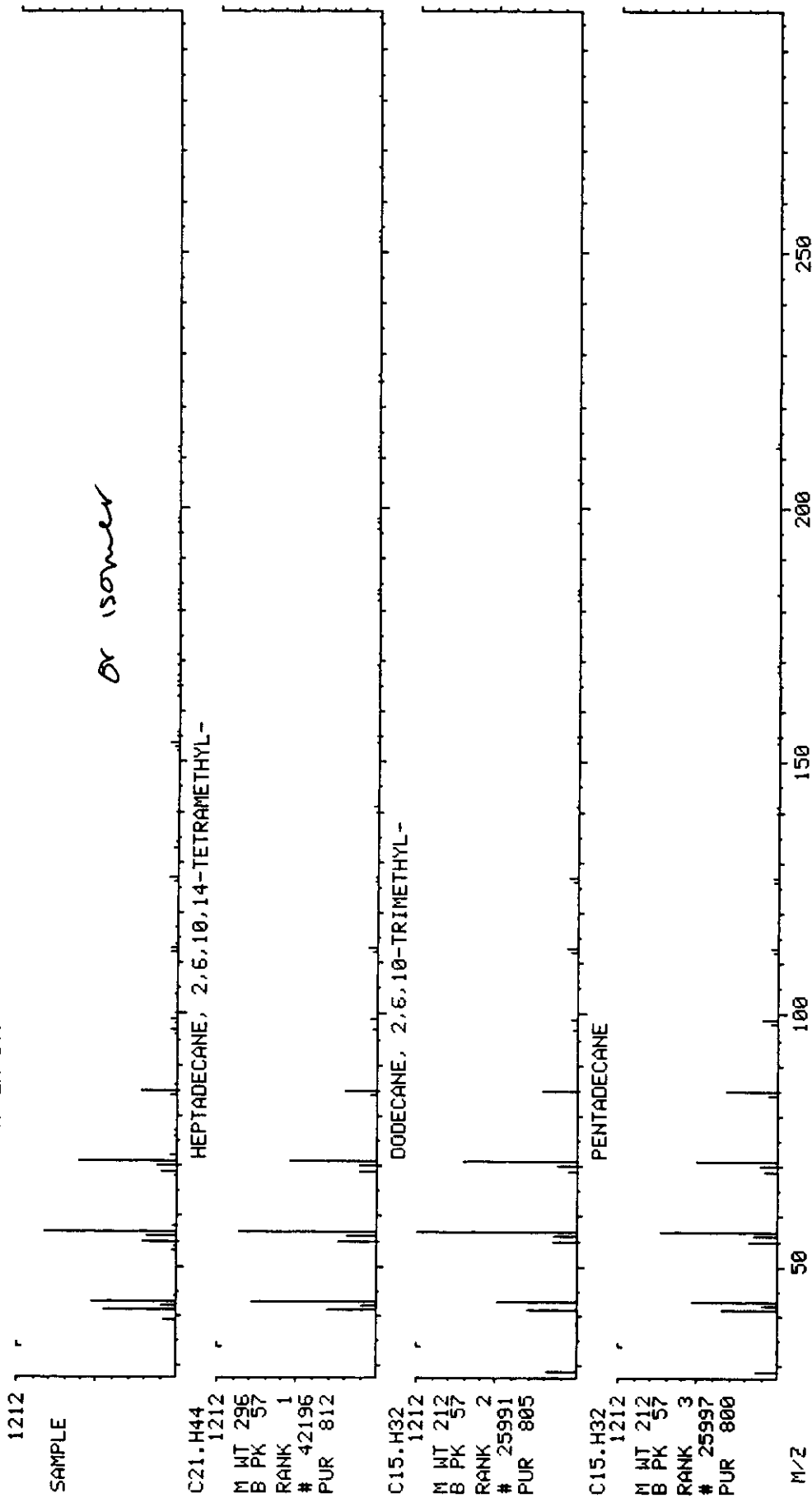
ENHANCED (S 15B 2N ØT)

DATA:	30068106	#	739
CALI:	30068106	#	3

CALI: 30068106 #

BASE M/2: 57
RIC: 158720.

RIC: 158720.



Library Search Data: 30068106 # 765 Base m/z: 43
08/31/98 21:30:00 + 8:31 Cali: 30068106 # 3 RIC: 285184.
Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
Enhanced (S 15B 2N 0T)

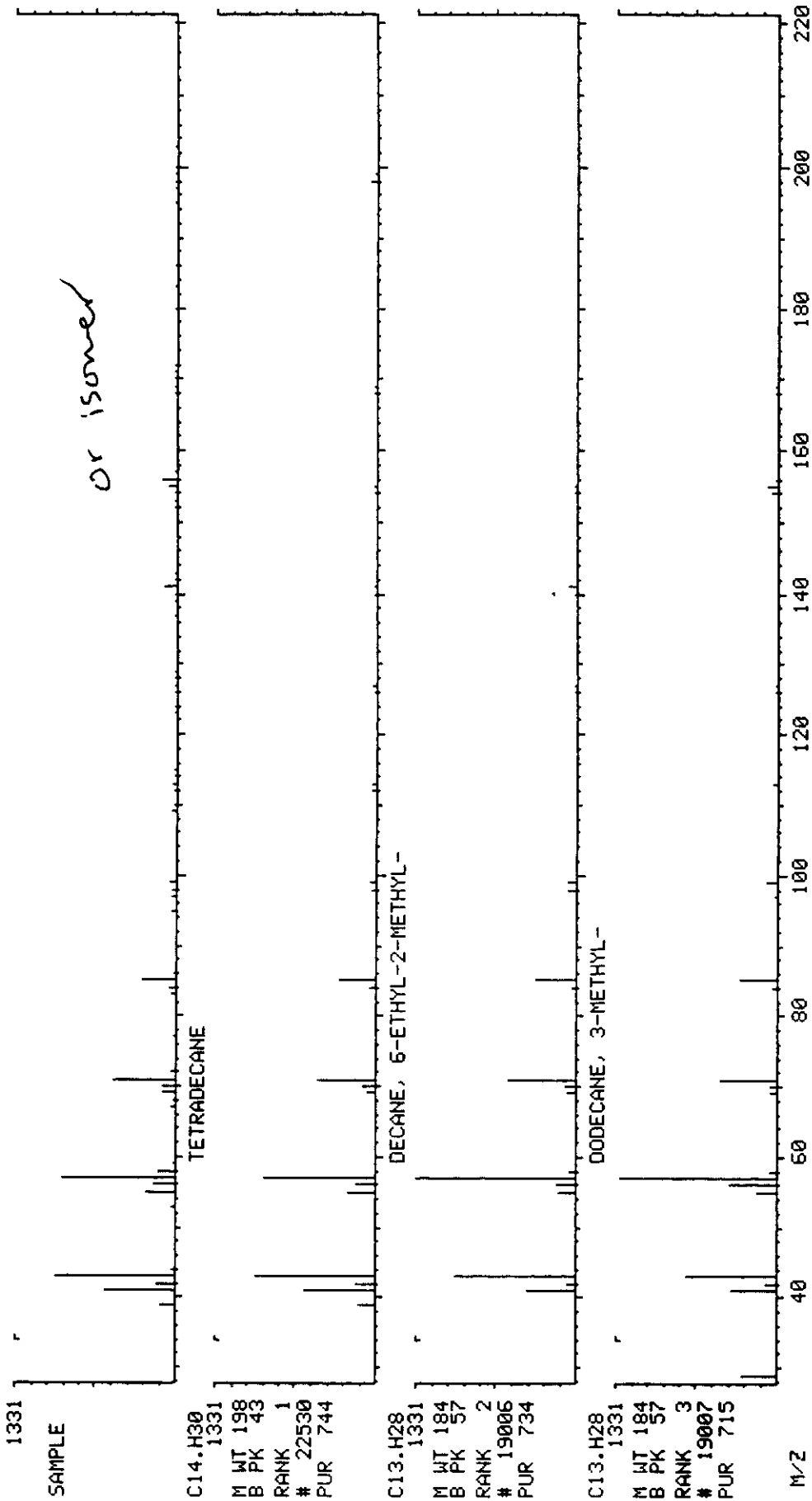
62231 spectra in LIBRARYNB searched for maximum PURITY
634 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
1 22530 TETRADECANE
2 19006 DECANE, 6-ETHYL-2-METHYL-
3 19007 DODECANE, 3-METHYL-
4 11607 UNDECANE
5 11612 NONANE, 2,5-DIMETHYL-
6 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
7 25996 TRIDECANE, 4,8-DIMETHYL-
8 39681 1-iodoundecane
9 25997 PENTADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C14.H30	198	43	744	960	744
2	C13.H28	184	57	734	963	734
3	C13.H28	184	57	715	917	715
4	C11.H24	156	43	711	977	711
5	C11.H24	156	57	708	877	708
6	C21.H44	296	57	706	925	726
7	C15.H32	212	57	705	929	710
8	C11.H23.I	282	57	703	888	707
9	C15.H32	212	57	703	928	722

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-59-4
2	—	—	—	—	62108-21-8
3	—	—	—	—	17312-57-1
4	—	—	—	—	1120-21-4
5	—	—	—	—	17302-27-1
6	—	—	—	—	18344-37-1
7	—	—	—	—	55030-62-1
8	—	—	—	—	4282-44-4
9	—	—	—	—	629-62-9

MID LIBRARY SEARCH <LIBRARYNB>
 08/31/98 21:30:00 + 8:31
 SAMPLE: T-MM5-3 1/35A/1ML
 COND.: UG/ML *1ML *1002/1002 *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30058106 # 765
 CALI: 30058106 # 3
 BASE M/Z: 43
 RIC: 285184.



Library Search Data: 30068106 # 778 Base m/z: 141
 08/31/98 21:30:00 + 8:39 Cali: 30068106 # 3 RIC: 79232.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 421 matched at least 5 of the 16 largest peaks in the unknown

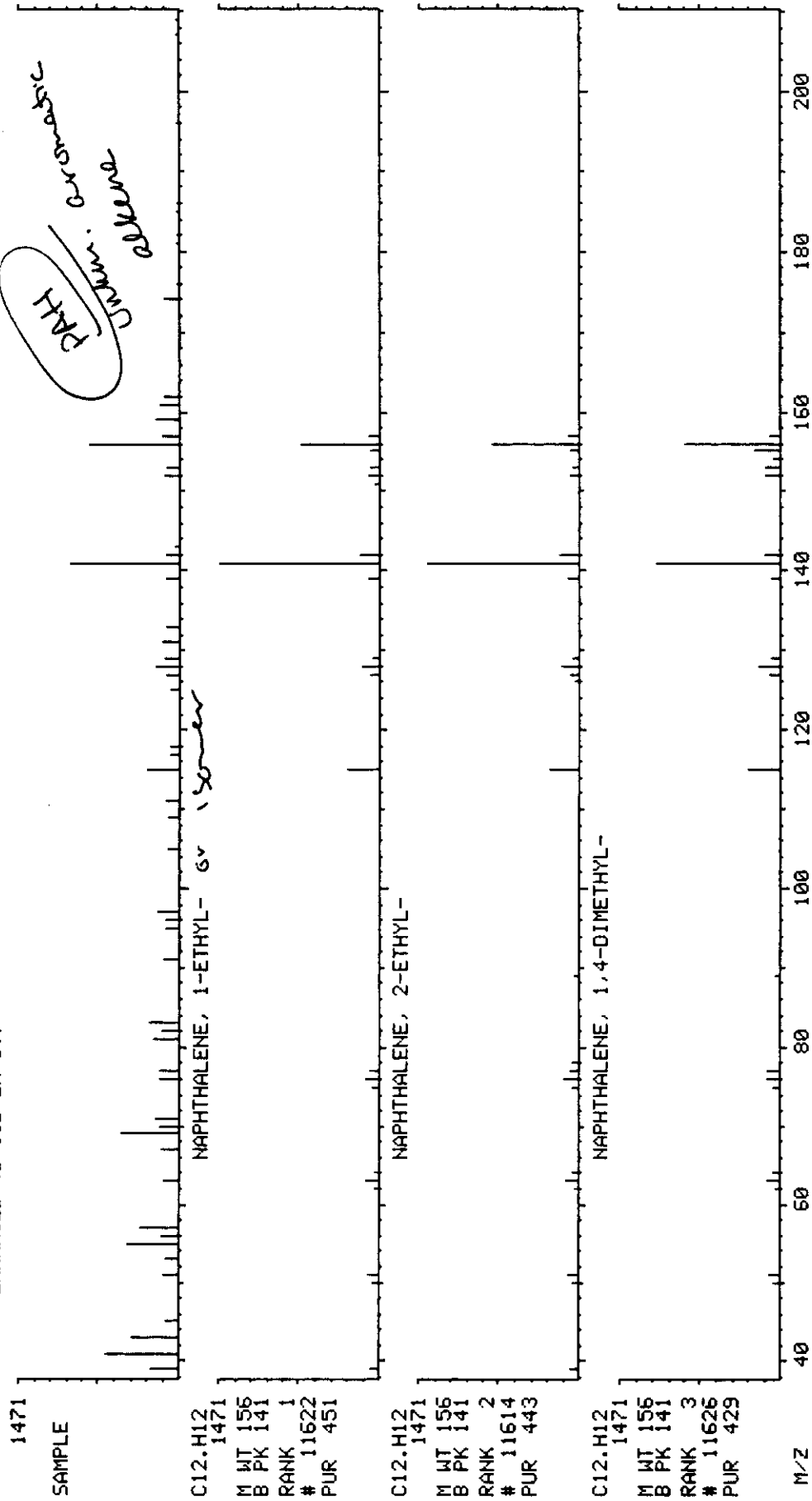
Rank In.	Name
1	11622 NAPHTHALENE, 1-ETHYL-
2	11614 NAPHTHALENE, 2-ETHYL-
3	11626 NAPHTHALENE, 1,4-DIMETHYL-
4	11619 NAPHTHALENE, 1,2-DIMETHYL-
5	11625 NAPHTHALENE, 1,3-DIMETHYL-
6	11618 NAPHTHALENE, 2,3-DIMETHYL-
7	11621 NAPHTHALENE, 1,8-DIMETHYL-
8	11624 NAPHTHALENE, 1,7-DIMETHYL-
9	11620 NAPHTHALENE, 2,6-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H12	156	141	451	910	477
2	C12.H12	156	141	443	892	480
3	C12.H12	156	141	429	882	481
4	C12.H12	156	156	429	866	477
5	C12.H12	156	156	421	850	483
6	C12.H12	156	156	417	841	474
7	C12.H12	156	156	408	840	471
8	C12.H12	156	156	408	823	484
9	C12.H12	156	156	407	838	479

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	1127-76-0
2	—	—	—	—	939-27-5
3	—	—	—	—	571-58-4
4	—	—	—	—	573-98-8
5	—	—	—	—	575-41-7
6	—	—	—	—	581-40-8
7	—	—	—	—	569-41-5
8	—	—	—	—	575-37-1
9	—	—	—	—	581-42-0

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:30:00 + 8:39
 SAMPLE: T-MM5-3 1/35A/1ML
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA) 1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068106 # 778
 CALI: 30068106 # 3
 BASE M/Z: 141
 RIC: 79232.



Library Search Data: 30068106 # 822 Base m/z: 57
 08/31/98 21:30:00 + 9:09 Cali: 30068106 # 3 RIC: 145152.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 855 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 2 19016 UNDECANE, 4,7-DIMETHYL-
 3 37462 HEPTADECANE, 2,6-DIMETHYL-
 4 15353 2,6-DIMETHYLDECANE
 5 18987 UNDECANE, 2,8-DIMETHYL-
 6 19023 DECANE, 2,6,7-TRIMETHYL-
 7 8104 OCTANE, 3,5-DIMETHYL-
 8 19035 DODECANE, 2-METHYL-
 9 19054 UNDECANE, 2,6-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C21.H44	296	57	792	959	816
2	C13.H28	184	43	758	947	765
3	C19.H40	268	57	757	923	803
4	C12.H26	170	43	754	916	774
5	C13.H28	184	43	753	959	758
6	C13.H28	184	57	749	948	749
7	C10.H22	142	57	747	961	747
8	C13.H28	184	43	745	950	755
9	C13.H28	184	57	742	897	757

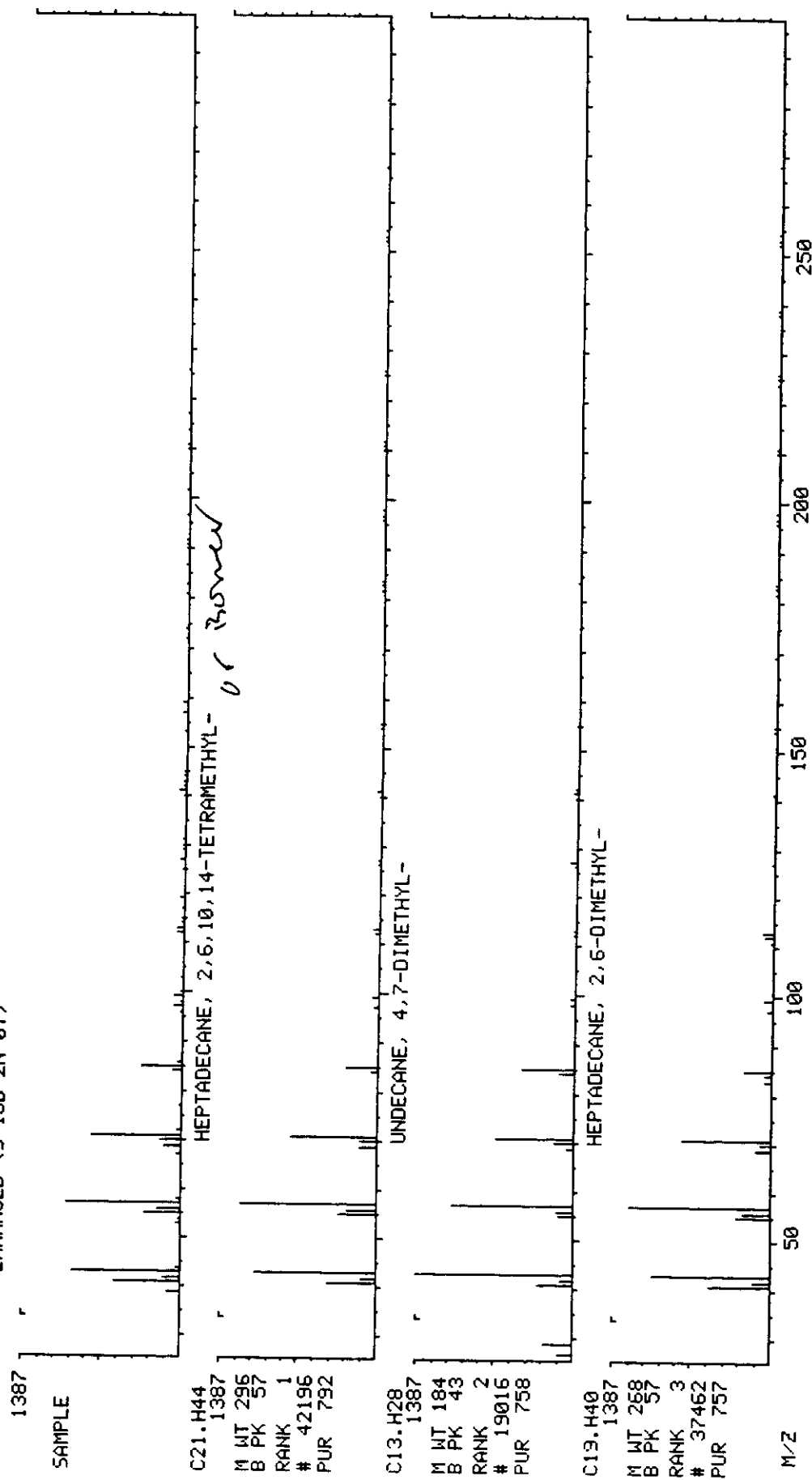
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	18344-37-1
2	—	—	—	—	17301-32-5
3	—	—	—	—	54105-67-8
4	—	—	—	—	13150-81-7
5	—	—	—	—	17301-25-6
6	—	—	—	—	62108-25-2
7	—	—	—	—	15869-93-9
8	—	—	—	—	1560-97-0
9	—	—	—	—	17301-23-4

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:30:00 + 9:09
 SAMPLE: T-MM5-3 1/35A/1ML
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA) >1/35A NA M
 ENHANCED (S 158 2N 0T)

BASE M/Z: 57
 RIC: 145152.

DATA: 30068106 # 822
 CALI: 30068106 # 3

INST. ID: F16
 NA M



Library Search Data: 30068106 # 860 Base m/z: 57
 08/31/98 21:30:00 + 9:34 Cali: 30068106 # 3 RIC: 246528.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

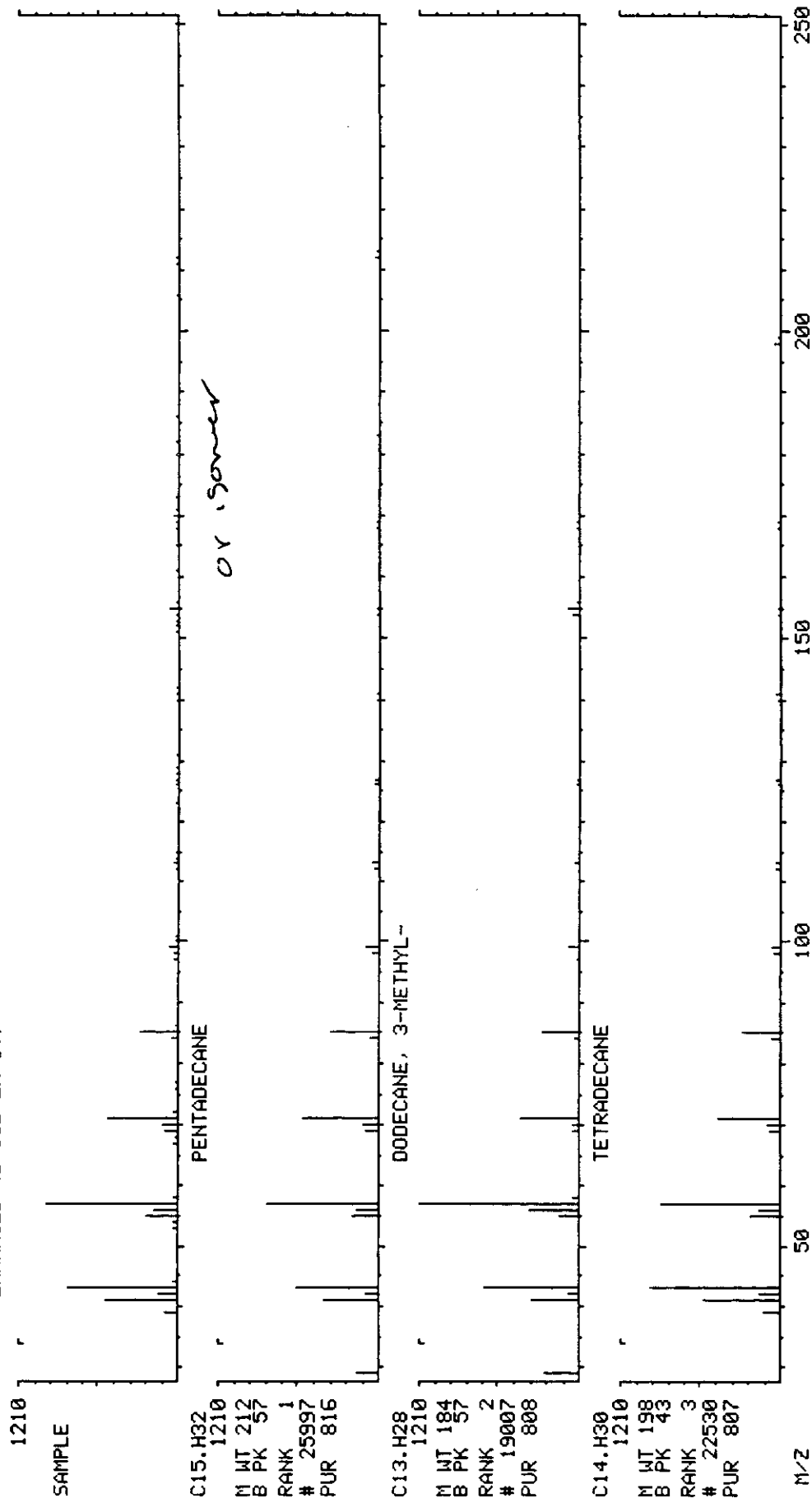
62231 spectra in LIBRARYNB searched for maximum PURITY
 792 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 25997 PENTADECANE
 2 19007 DODECANE, 3-METHYL-
 3 22530 TETRADECANE
 4 15352 UNDECANE, 2-METHYL-
 5 15343 DODECANE
 6 19006 DECAN, 6-ETHYL-2-METHYL-
 7 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 8 15353 2,6-DIMETHYLDECANE
 9 18998 UNDECANE, 3,7-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C15.H32	212	57	816	963	823
2	C13.H28	184	57	808	965	808
3	C14.H30	198	43	807	945	841
4	C12.H26	170	43	807	952	807
5	C12.H26	170	57	800	985	800
6	C13.H28	184	57	799	967	799
7	C21.H44	296	57	797	939	819
8	C12.H26	170	43	796	923	796
9	C13.H28	184	43	790	949	791

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-62-9
2	—	—	—	—	17312-57-1
3	—	—	—	—	629-59-4
4	—	—	—	—	7045-71-8
5	—	—	—	—	112-40-3
6	—	—	—	—	62108-21-8
7	—	—	—	—	18344-37-1
8	—	—	—	—	13150-81-7
9	—	—	—	—	17301-29-0

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:30:00 + 9:34
 SAMPLE: T-MM5-3 1/35A/1ML
 COND5.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30068106 # 860
 CALI: 30068106 # 3
 BASE M/Z: 57
 RIC: 246528.



Library Search Data: 30068106 # 905 Base m/z: 41
 08/31/98 21:30:00 + 10:04 Cali: 30068106 # 3 RIC: 89088.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 751 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 40233 NONADECANOL
 2 42521 1-EICOSANOL
 3 37444 OXIRANE, HEXADECYL-
 4 40193 CIS-9,10-EPOXYOCTADECAN-1-OL
 5 37449 OCTADECANAL
 6 42180 OCTADECANE, 1-(ETHENYLOXY)-
 7 25971 OXIRANE, DODECYL-
 8 22482 2-TRIDECEN-1-OL, (E)-
 9 32052 OXIRANE, TETRADECYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40.O	284	43	653	934	682
2	C20.H42.O	298	43	647	928	680
3	C18.H36.O	268	57	638	853	674
4	C18.H36.O2	284	55	638	904	668
5	C18.H36.O	268	43	628	851	687
6	C20.H40.O	296	43	625	939	650
7	C14.H28.O	212	41	625	891	643
8	C13.H26.O	198	57	621	898	644
9	C16.H32.O	240	41	619	935	619

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	52783-43-4
2	—	—	—	—	629-96-9
3	—	—	—	—	7390-81-0
4	—	—	—	—	13980-12-6
5	—	—	—	—	638-66-4
6	—	—	—	—	930-02-9
7	—	—	—	—	3234-28-4
8	—	—	—	—	74962-98-4
9	—	—	—	—	7320-37-8

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 21:30:00 + 10:04

SAMPLE: T-MM5-3 1/35A/1ML

COND.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M

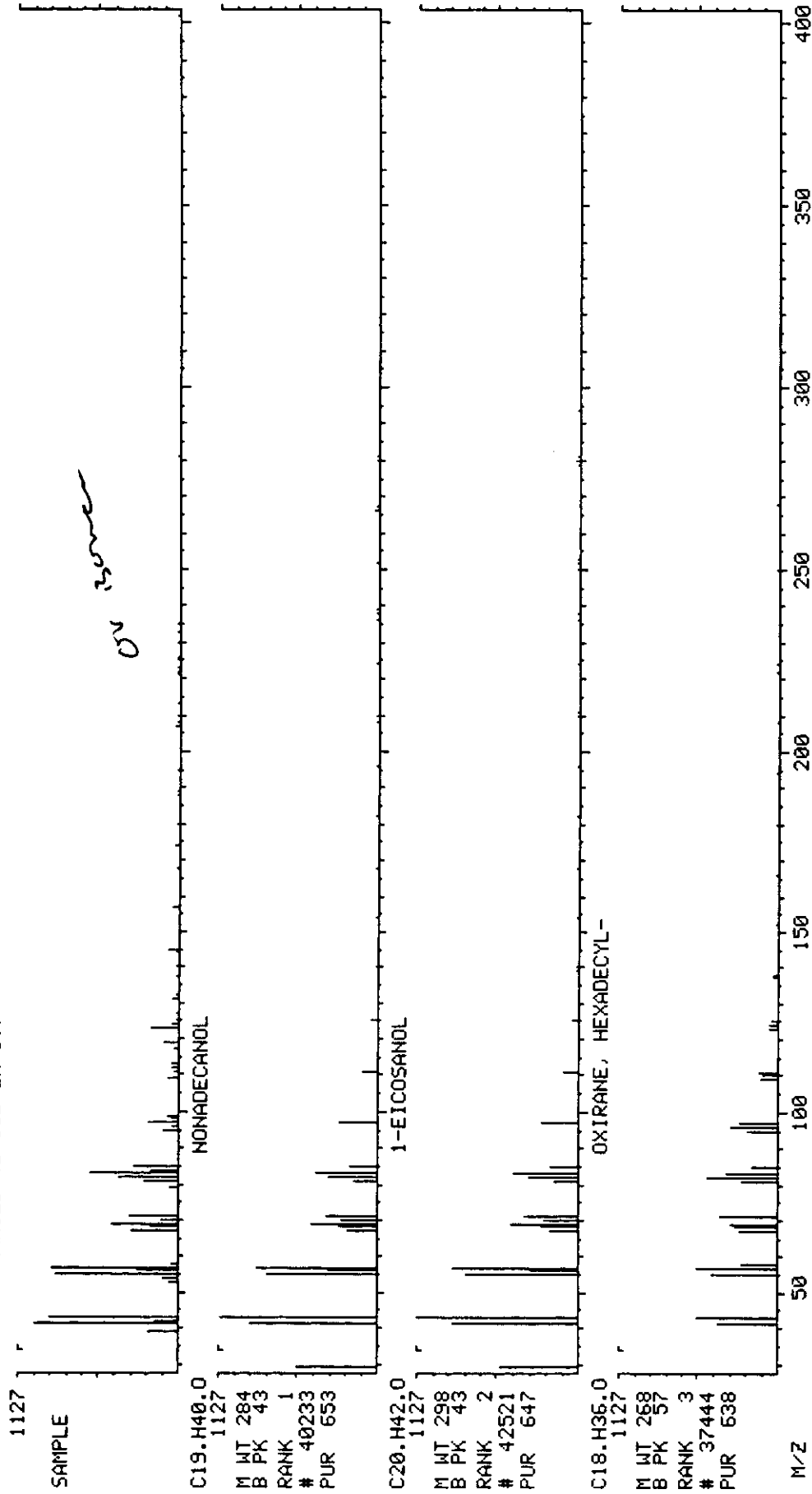
ENHANCED (S 15B 2N 0T)

DATA: 30068106 # 905
CALI: 30068106 # 3

CALI: 30068106 # 3

BASE M/Z: 41
RIC: 89088.

RIC: 89088.



Library Search Data: 30068106 # 949 Base m/z: 57
08/31/98 21:30:00 + 10:33 Cali: 30068106 # 3 RIC: 260096.
Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
360 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
1 37465 NONADECANE
2 37462 HEPTADECANE, 2,6-DIMETHYL-
3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
4 15969 HYDROXYLAMINE, O-DECYL-
5 19016 UNDECANE, 4,7-DIMETHYL-
6 25997 PENTADECANE
7 19054 UNDECANE, 2,6-DIMETHYL-
8 18985 TRIDECANE
9 19523 1-DECANOL, 2-ETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	842	952	867
2	C19.H40	268	57	840	944	865
3	C21.H44	296	57	831	951	849
4	C10.H23.O.N	173	43	823	962	826
5	C13.H28	184	43	819	962	819
6	C15.H32	212	57	819	953	844
7	C13.H28	184	57	805	919	805
8	C13.H28	184	57	804	964	807
9	C12.H26.O	186	57	800	927	813

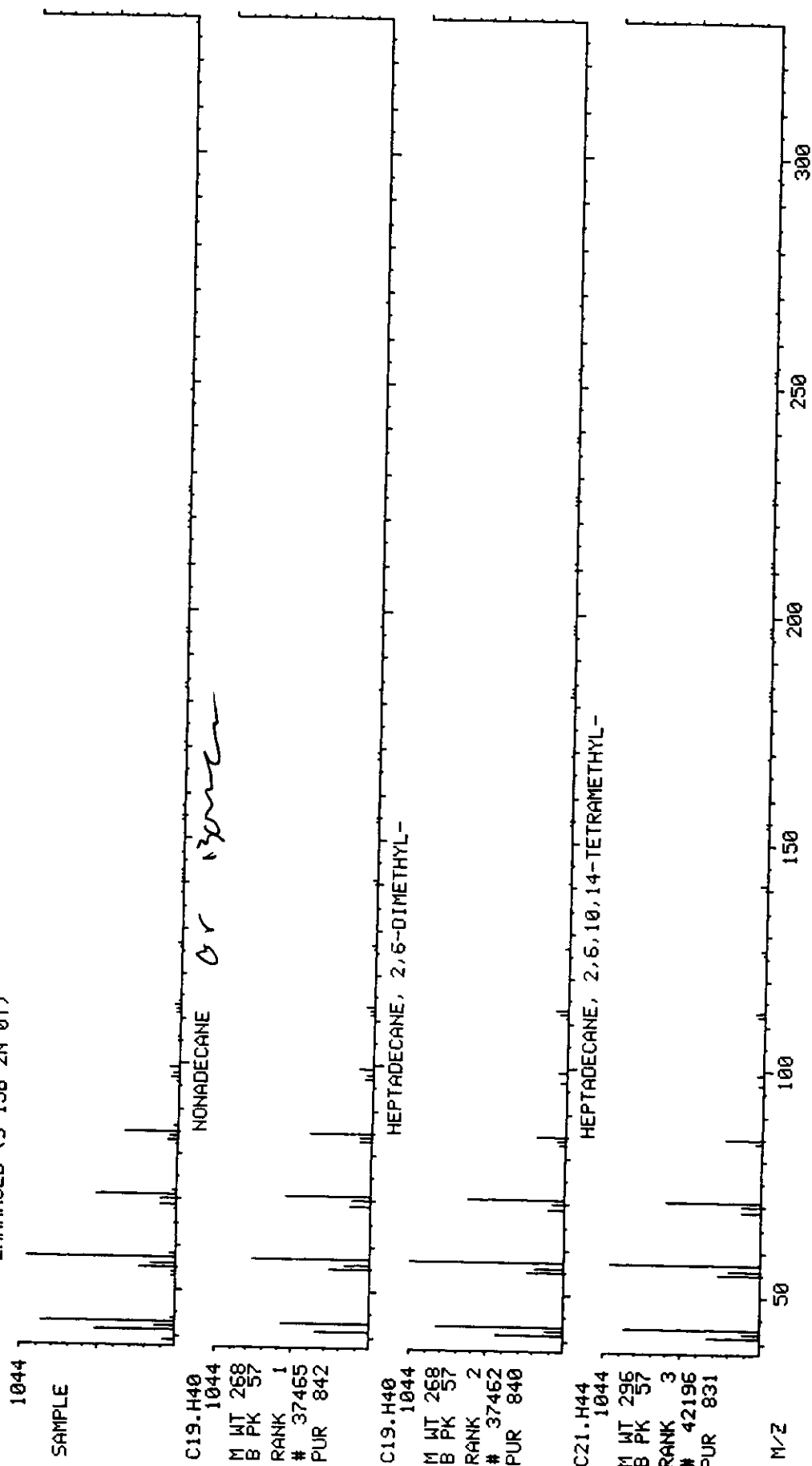
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	18344-37-1
4	—	—	—	—	29812-79-1
5	—	—	—	—	17301-32-5
6	—	—	—	—	629-62-9
7	—	—	—	—	17301-23-4
8	—	—	—	—	629-50-5
9	—	—	—	—	21078-65-9

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 21:30:00 + 10:33
 SAMPLE: T-MM5-3 1/35A/1ML
 COND5.: UG/ML *1ML *1002/1002 *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30058106 # 949
 CALI: 30058106 # 3

BASE M/Z: 57
 RIC: 250096.



Library Search Data: 30068106 # 989 Base m/z: 57
 08/31/98 21:30:00 + 11:00 Cali: 30068106 # 3 RIC: 157440.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 691 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 25996 TRIDECANE, 4,8-DIMETHYL-
 3 19054 UNDECANE, 2,6-DIMETHYL-
 4 22536 TRIDECANE, 3-METHYL-
 5 37465 NONADECANE
 6 25997 PENTADECANE
 7 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 8 22530 TETRADECANE
 9 29264 TRIDECANE, 5-PROPYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	834	955	844
2	C15.H32	212	57	824	949	825
3	C13.H28	184	57	820	948	830
4	C14.H30	198	57	816	968	816
5	C19.H40	268	57	814	930	842
6	C15.H32	212	57	807	936	830
7	C21.H44	296	57	807	943	823
8	C14.H30	198	43	803	925	823
9	C16.H34	226	57	803	957	807

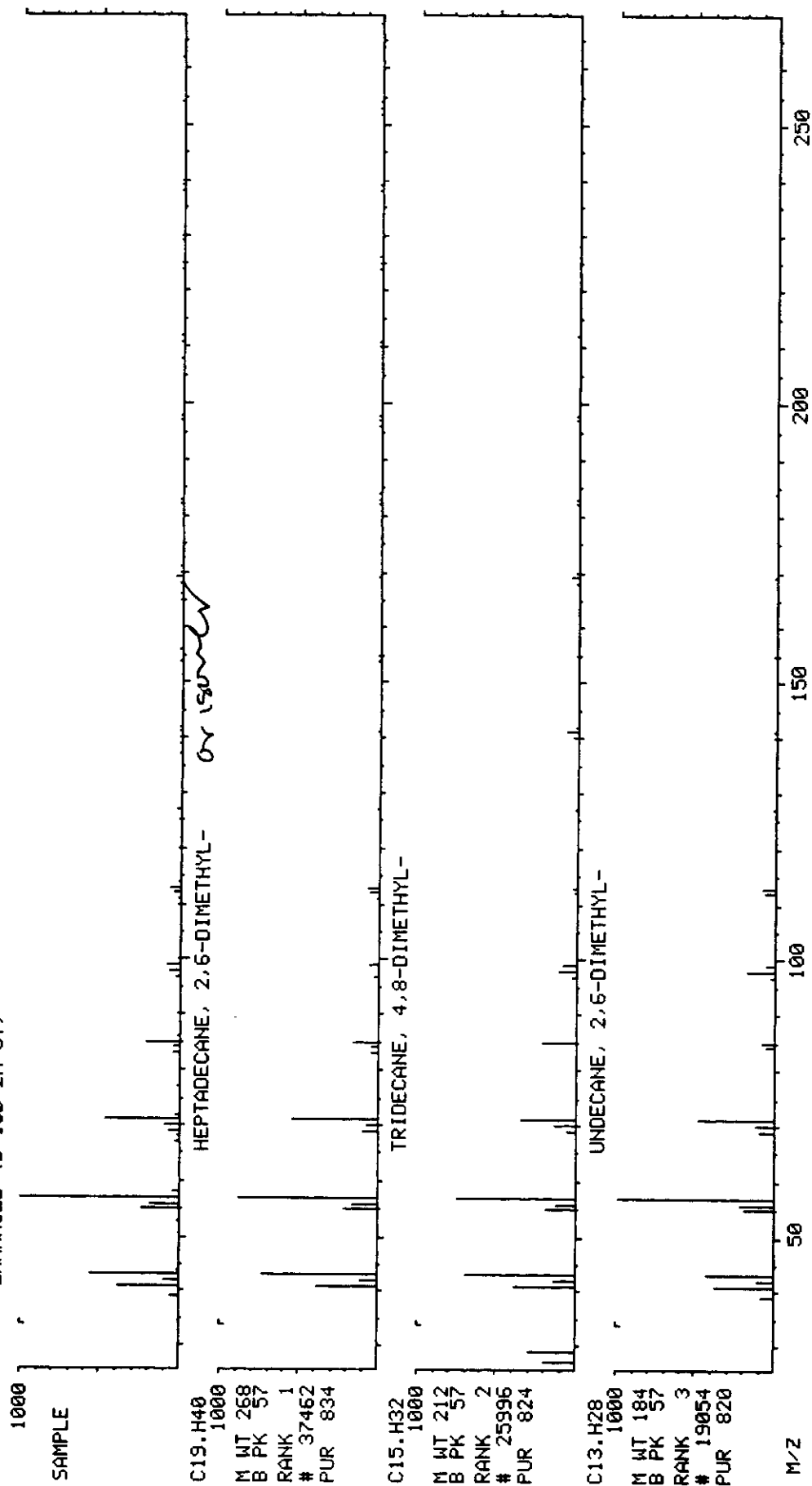
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	55030-62-1
3	—	—	—	—	17301-23-4
4	—	—	—	—	6418-41-3
5	—	—	—	—	629-92-5
6	—	—	—	—	629-62-9
7	—	—	—	—	18344-37-1
8	—	—	—	—	629-59-4
9	—	—	—	—	55045-11-9

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:30:00 + 11:00
 SAMPLE: T-MMS-3 1/35A/1ML
 CONDS.: UG/ML *1ML *100%/1002 *(NA/NA)/1/35A NA M
 ENHANCED (5 158 2N 0T)

DATA: 30068105 # 989
 CALL: 30068105 # 3

BASE M/Z: 57
 RIC: 157440.

INST. ID: F16
 NA M



Library Search Data: 30068106 #1037 Base m/z: 57
 08/31/98 21:30:00 + 11:32 Cali: 30068106 # 3 RIC: 239616.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 713 matched at least 7 of the 16 largest peaks in the unknown

Rank	In.	Name
1	37462	HEPTADECANE, 2,6-DIMETHYL-
2	42196	HEPTADECANE, 2,6,10,14-TETRAMETHYL-
3	26001	DODECANE, 2,7,10-TRIMETHYL-
4	25991	DODECANE, 2,6,10-TRIMETHYL-
5	22535	DODECANE, 4,6-DIMETHYL-
6	37466	PENTADECANE, 2,6,10,14-TETRAMETHYL-
7	37465	NONADECANE
8	25994	DODECANE, 2,6,11-TRIMETHYL-
9	22534	TRIDECANE, 7-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	872	974	880
2	C21.H44	296	57	830	945	855
3	C15.H32	212	57	818	966	824
4	C15.H32	212	57	818	968	823
5	C14.H30	198	57	818	962	819
6	C19.H40	268	71	816	944	832
7	C19.H40	268	57	812	928	836
8	C15.H32	212	57	796	941	807
9	C14.H30	198	57	792	914	805

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	18344-37-1
3	—	—	—	—	74645-98-0
4	—	—	—	—	3891-98-3
5	—	—	—	—	61141-72-8
6	—	—	—	—	1921-70-6
7	—	—	—	—	629-92-5
8	—	—	—	—	31295-56-4
9	—	—	—	—	26730-14-3

MID LIBRARY SEARCH (LIBRARYNB)

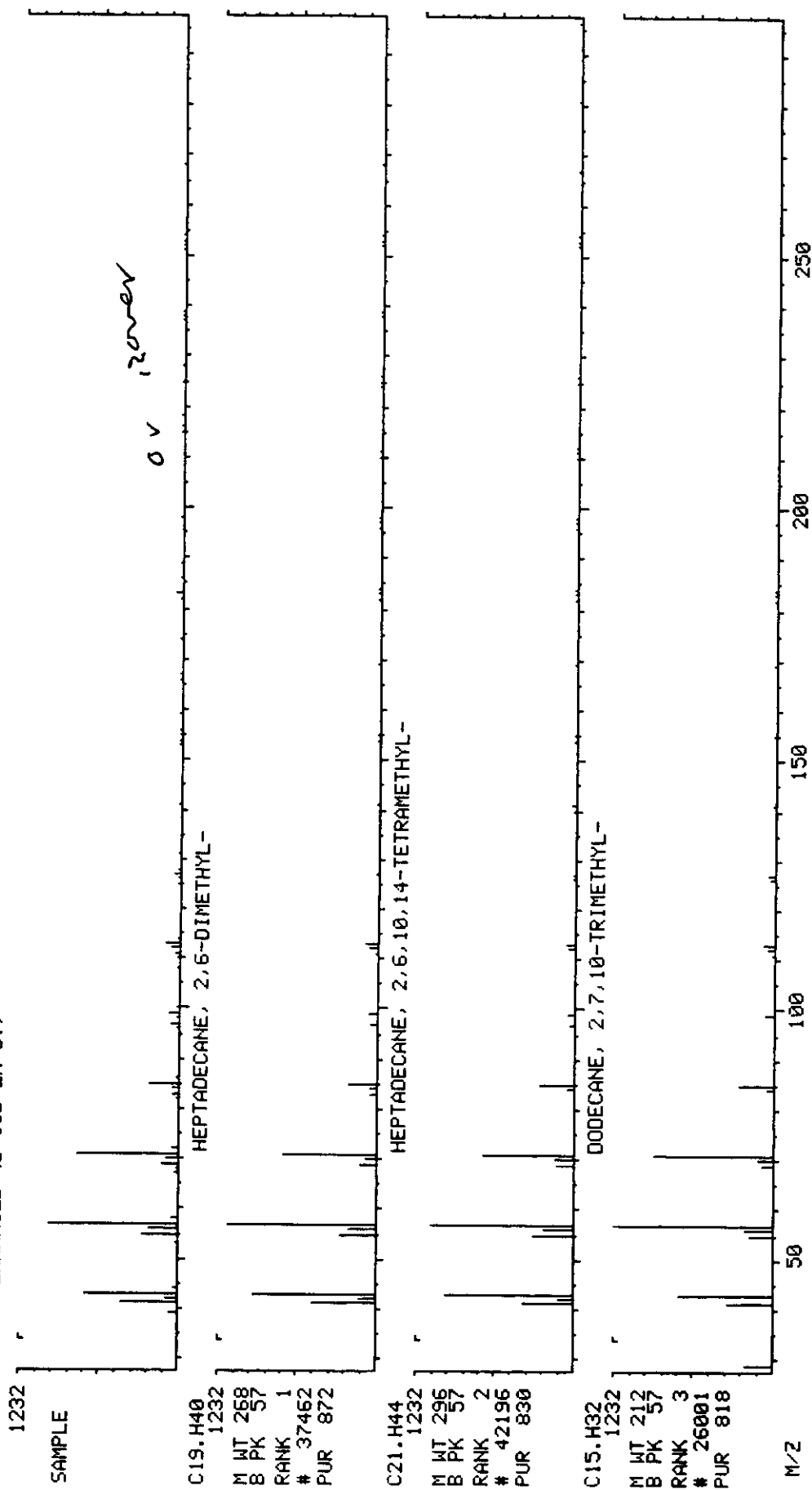
08/31/98 21:30:00 + 11:32

SAMPLE: T-MM5-3 1/35A/1ML

COND.: UG/ML *1ML *1002/1002 *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)

DATA: 30068106 #1037
CALI: 30068106 # 3

BASE M/Z: 57
RIC: 239615.



Library Search Data: 30068106 #1115 Base m/z: 57
 08/31/98 21:30:00 + 12:24 Cali: 30068106 # 3 RIC: 131328.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 934 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 22530 TETRADECANE
 3 25997 PENTADECANE
 4 19016 UNDECANE, 4,7-DIMETHYL-
 5 19006 DECANE, 6-ETHYL-2-METHYL-
 6 15969 HYDROXYLAMINE, O-DECYL-
 7 15358 UNDECANE, 3-METHYL-
 8 37462 HEPTADECANE, 2,6-DIMETHYL-
 9 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-

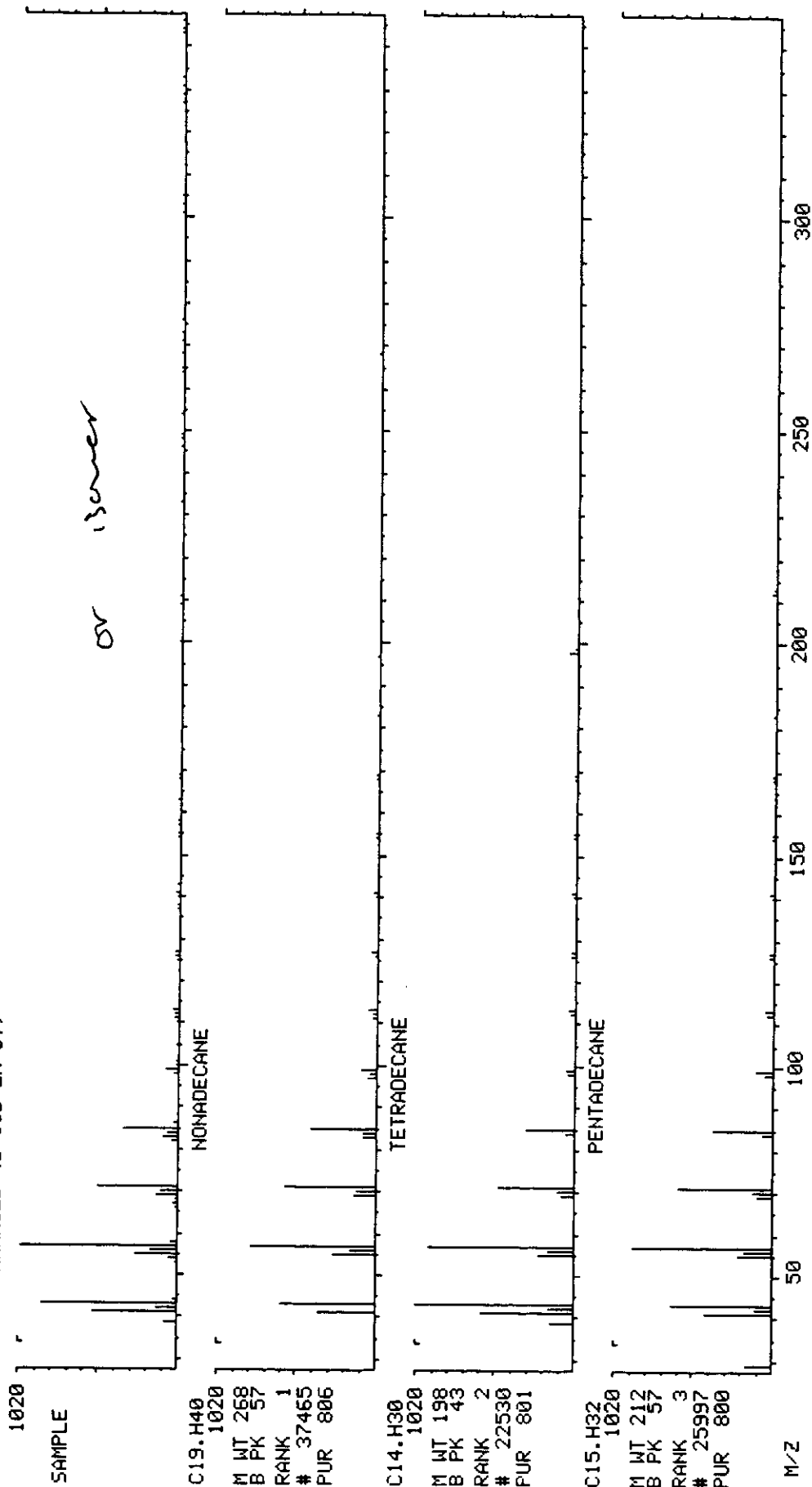
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	806	947	832
2	C14.H30	198	43	801	950	835
3	C15.H32	212	57	800	959	823
4	C13.H28	184	43	788	958	789
5	C13.H28	184	57	785	966	785
6	C10.H23.O.N	173	43	780	949	798
7	C12.H26	170	57	780	945	783
8	C19.H40	268	57	780	917	826
9	C21.H44	296	57	778	937	812

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	629-59-4
3	—	—	—	—	629-62-9
4	—	—	—	—	17301-32-5
5	—	—	—	—	62108-21-8
6	—	—	—	—	29812-79-1
7	—	—	—	—	1002-43-3
8	—	—	—	—	54105-67-8
9	—	—	—	—	18344-37-1

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:30:00 + 12:24
 SAMPLE: T-MM5-3 1/35A/1ML
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (5 15B 2N 0T)

DATA: 30068106 #1115
 CALI: 30068106 # 3

BASE M/Z: 57
 RIC: 131328.



Library Search Data: 30068106 #1120 Base m/z: 57
 08/31/98 21:30:00 + 12:27 Cali: 30068106 # 3 RIC: 135680.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 840 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 46161 TRICOSANE
 5 25991 DODECANE, 2,6,10-TRIMETHYL-
 6 42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-
 7 26001 DODECANE, 2,7,10-TRIMETHYL-
 8 29264 TRIDECANE, 5-PROPYL-
 9 39859 OCTADECANE, 2,6-DIMETHYL-

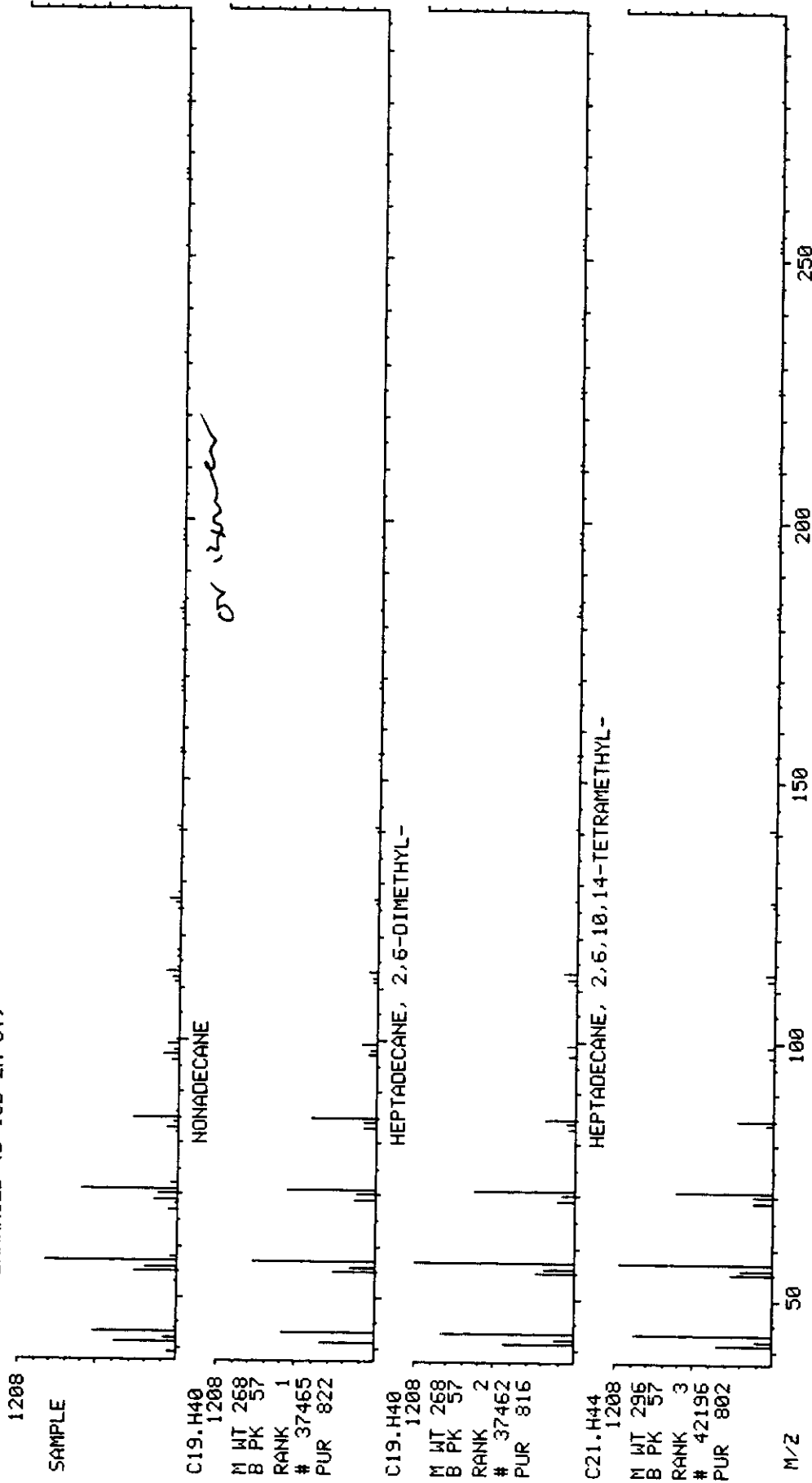
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	822	952	841
2	C19.H40	268	57	816	937	829
3	C21.H44	296	57	802	934	818
4	C23.H48	324	43	792	905	847
5	C15.H32	212	57	788	961	791
6	C21.H44	296	57	782	945	810
7	C15.H32	212	57	782	964	782
8	C16.H34	226	57	782	931	784
9	C20.H42	282	57	781	920	795

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	18344-37-1
4	—	—	—	—	638-67-5
5	—	—	—	—	3891-98-3
6	—	—	—	—	54833-48-6
7	—	—	—	—	74645-98-0
8	—	—	—	—	55045-11-9
9	—	—	—	—	75163-97-2

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:30:00 + 12:27
 SAMPLE: T-MM5-3 1/35A/1ML
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068105 #1120
 CALI: 30068105 # 3

BASE M/Z: 57
 RIC: 135680.



Library Search Data: 30068106 #1192 Base m/z: 57
 08/31/98 21:30:00 + 13:16 Cali: 30068106 # 3 RIC: 107648.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 294 matched at least 8 of the 16 largest peaks in the unknown

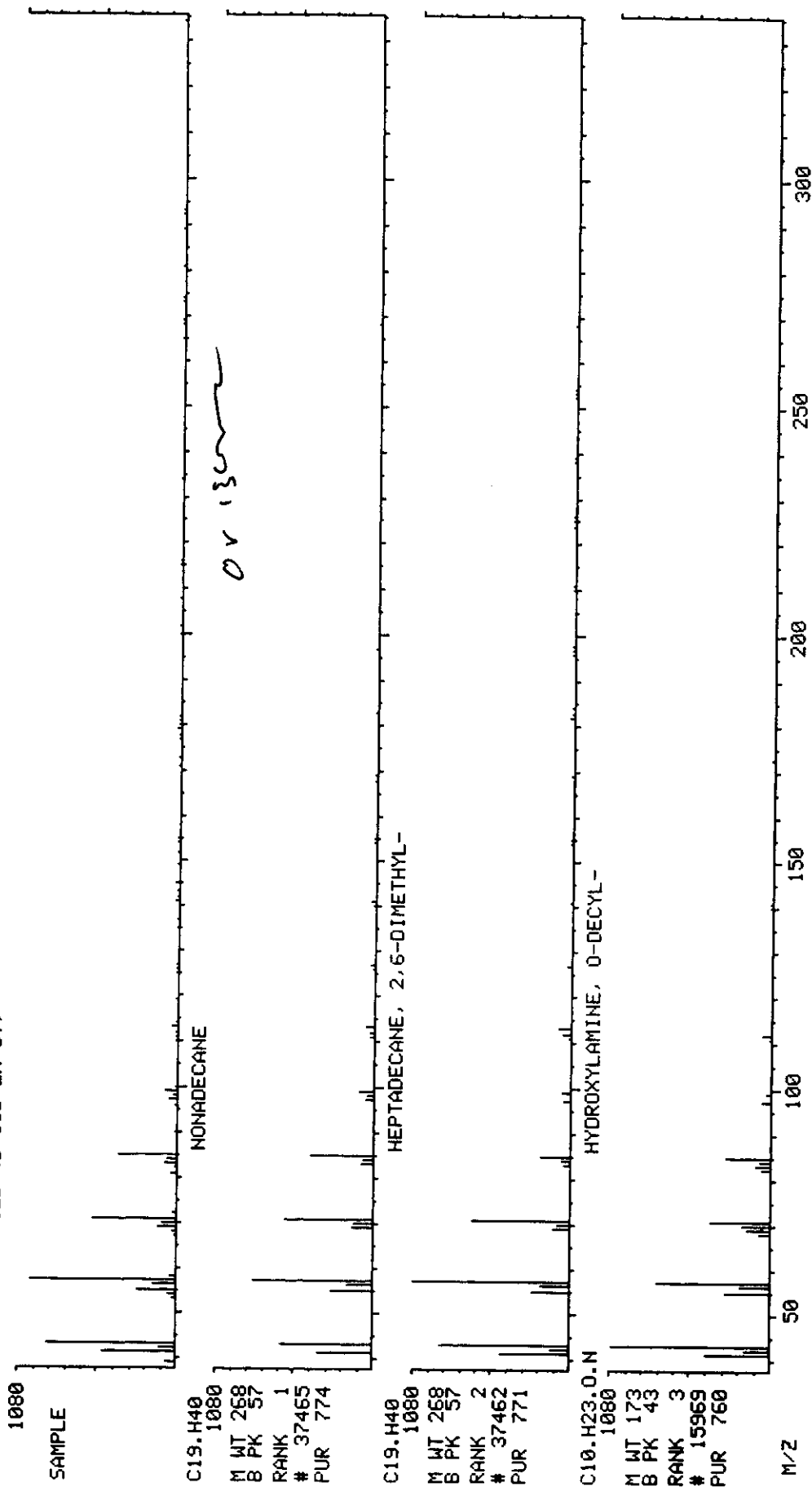
Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 15969 HYDROXYLAMINE, O-DECYL-
 4 19016 UNDECANE, 4,7-DIMETHYL-
 5 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 6 25997 PENTADECANE
 7 19523 1-DECANOL, 2-ETHYL-
 8 15353 2,6-DIMETHYLDECANE
 9 22535 DODECANE, 4,6-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	774	950	802
2	C19.H40	268	57	771	940	796
3	C10.H23.O.N	173	43	760	957	768
4	C13.H28	184	43	746	955	747
5	C21.H44	296	57	744	938	778
6	C15.H32	212	57	732	936	771
7	C12.H26.O	186	57	732	929	739
8	C12.H26	170	43	732	921	743
9	C14.H30	198	57	729	947	736

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	629-92-5
2	---	---	---	---	54105-67-8
3	---	---	---	---	29812-79-1
4	---	---	---	---	17301-32-5
5	---	---	---	---	18344-37-1
6	---	---	---	---	629-62-9
7	---	---	---	---	21078-65-9
8	---	---	---	---	13150-81-7
9	---	---	---	---	61141-72-8

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:30:00 + 13:16
 SAMPLE: T-MMS-3 1/35A/1ML
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA) 1/35A NA M
 ENHANCED (5 15B 2N 0T)

DATA: 30068106 #1192
 CALI: 30068106 # 3
 BASE M/Z: 57
 RIC: 107648.



1080

Library Search Data: 30068106 #1336 Base m/z: 57
 08/31/98 21:30:00 + 14:52 Cali: 30068106 # 3 RIC: 73216.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 429 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37456 2-METHYLOCTADECANE
 2 32058 HEXADECANE, 3-METHYL-
 3 49555 PENTACOSANE
 4 37462 HEPTADECANE, 2,6-DIMETHYL-
 5 46161 TRICOSANE
 6 29263 HEXADECANE
 7 42197 HENEICOSANE
 8 37465 NONADECANE
 9 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-

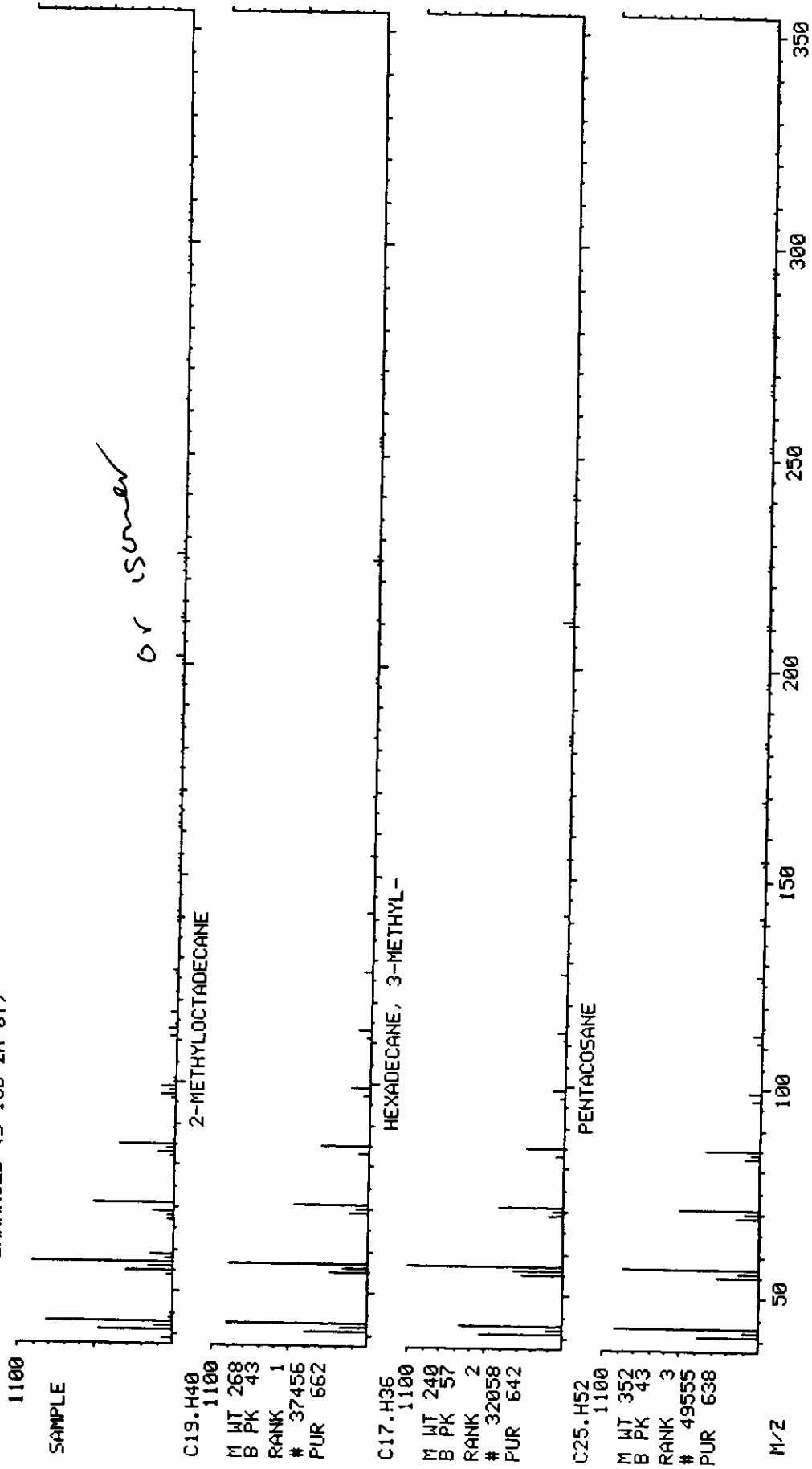
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	43	662	908	690
2	C17.H36	240	57	642	922	661
3	C25.H52	352	43	638	877	716
4	C19.H40	268	57	637	850	682
5	C23.H48	324	43	632	888	703
6	C16.H34	226	57	629	904	672
7	C21.H44	296	57	629	894	668
8	C19.H40	268	57	627	917	675
9	C21.H44	296	57	626	861	671

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	- -
2	—	—	—	—	6418-43-5
3	—	—	—	—	629-99-2
4	—	—	—	—	54105-67-8
5	—	—	—	—	638-67-5
6	—	—	—	—	544-76-3
7	—	—	—	—	629-94-7
8	—	—	—	—	629-92-5
9	—	—	—	—	18344-37-1

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:30:00 + 14:52
 SAMPLE: T-MM5-3 1/35A/1ML
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068105 #1335
 CALI: 30068105 # 3

BASE M/Z: 57
 RIC: 73216.



1020

Library Search Data: 30068106 #1531 Base m/z: 57
 08/31/98 21:30:00 + 17:02 Cali: 30068106 # 3 RIC: 37056.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 238 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 15969 HYDROXYLAMINE, O-DECYL-
 4 22536 TRIDECANE, 3-METHYL-
 5 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 6 19035 DODECANE, 2-METHYL-
 7 19007 DODECANE, 3-METHYL-
 8 46161 TRICOSANE
 9 32058 HEXADECANE, 3-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	663	935	694
2	C19.H40	268	57	651	911	681
3	C10.H23.O.N	173	43	639	934	654
4	C14.H30	198	57	634	910	658
5	C21.H44	296	57	625	922	651
6	C13.H28	184	43	622	941	624
7	C13.H28	184	57	621	907	629
8	C23.H48	324	43	618	874	696
9	C17.H36	240	57	617	894	651

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	29812-79-1
4	—	—	—	—	6418-41-3
5	—	—	—	—	18344-37-1
6	—	—	—	—	1560-97-0
7	—	—	—	—	17312-57-1
8	—	—	—	—	638-67-5
9	—	—	—	—	6418-43-5

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 21:30:00 + 17:02

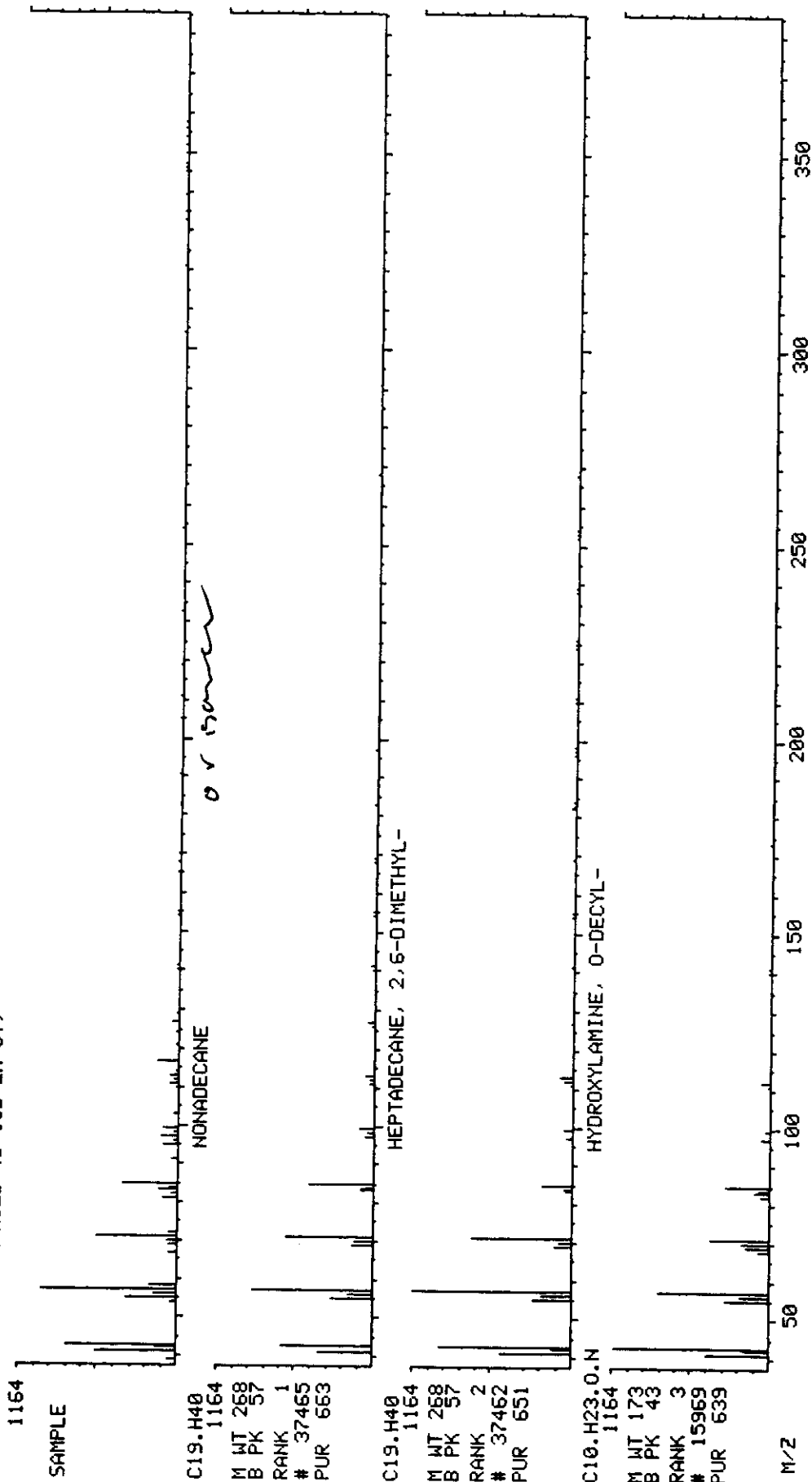
SAMPLE: T-MM5-3 1/35A/1ML

COND.: UG/ML *1ML *1002/1002 *(NA/NA) >1/35A NA M

ENHANCED (S 15B 2N 0T)

DATA: 30068106 #1531
CALI: 30068106 # 3

BASE M/Z: 57
RIC: 37056.



1000

Library Search Data: 30068106 #1587 Base m/z: 43
 08/31/98 21:30:00 + 17:39 Cali: 30068106 # 3 RIC: 106112.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 950 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 39515 9-EICOSENE, (E)-
 2 34410 3-OCTADECENE, (E)-
 3 39516 5-EICOSENE, (E)-
 4 34418 9-OCTADECENE, (E)-
 5 34411 5-OCTADECENE, (E)-
 6 31653 1-HEPTADECENE
 7 39517 3-EICOSENE, (E)-
 8 26416 4-TETRADECANOL
 9 32420 1-HEXADECANOL

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C20.H40	280	55	795	984	799
2	C18.H36	252	69	795	975	800
3	C20.H40	280	55	794	983	799
4	C18.H36	252	55	794	974	797
5	C18.H36	252	55	792	971	796
6	C17.H34	238	55	781	973	794
7	C20.H40	280	57	781	984	786
8	C14.H30.O	214	43	781	986	788
9	C16.H34.O	242	55	777	975	788

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	74685-29-3
2	—	—	—	—	7206-19-1
3	—	—	—	—	74685-30-6
4	—	—	—	—	7206-25-9
5	—	—	—	—	7206-21-5
6	—	—	—	—	6765-39-5
7	—	—	—	—	74685-33-9
8	—	—	—	—	1653-33-4
9	—	—	—	—	36653-82-4

Library Search Data: 30068106 #1809 Base m/z: 57
 08/31/98 21:30:00 + 20:07 Cali: 30068106 # 3 RIC: 18560.
 Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 334 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 52672 SILANE, TRICHLOROCTADECYL-
 2 40165 HEXADECANOIC ACID, 2-OXO-, METHYL ESTER
 3 46251 DODECANE, 1,2-DIBROMO-
 4 31486 ACETAMIDE, N-METHYL-N-[4-[4-METHOXY-1-HEXAHYDROPYRIDYL]-2-BUTYNYL]-
 5 5049 OXIRANE, (3,3-DIMETHYLBUTYL)-
 6 27504 1-BROMO-3,7-DIMETHYLOCTANE
 7 41993 1-iodo-2-methylundecane
 8 20601 2-BROMO-6-METHYLHEPTANE
 9 40866 OCTADECANE, 1-CHLORO-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C18.H37.CL3.SI	386	57	329	706	393
2	C17.H32.O3	284	57	311	873	324
3	C12.H24.BR2	326	41	302	796	353
4	C13.H22.O2.N2	238	43	292	793	349
5	C8.H16.O	128	57	284	954	288
6	C10.H21.BR	220	57	283	856	299
7	C12.H25.I	296	57	281	872	306
8	C8.H17.BR	192	57	276	921	278
9	C18.H37.CL	288	57	275	649	340

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	112-04-9
2	—	—	—	—	55836-30-1
3	—	—	—	—	55334-42-4
4	—	—	—	—	- -
5	—	—	—	—	53907-77-0
6	—	—	—	—	3383-83-3
7	—	—	—	—	73105-67-6
8	—	—	—	—	4730-24-9
9	—	—	—	—	3386-33-2

MID LIBRARY SEARCH <LIBRARYNB>

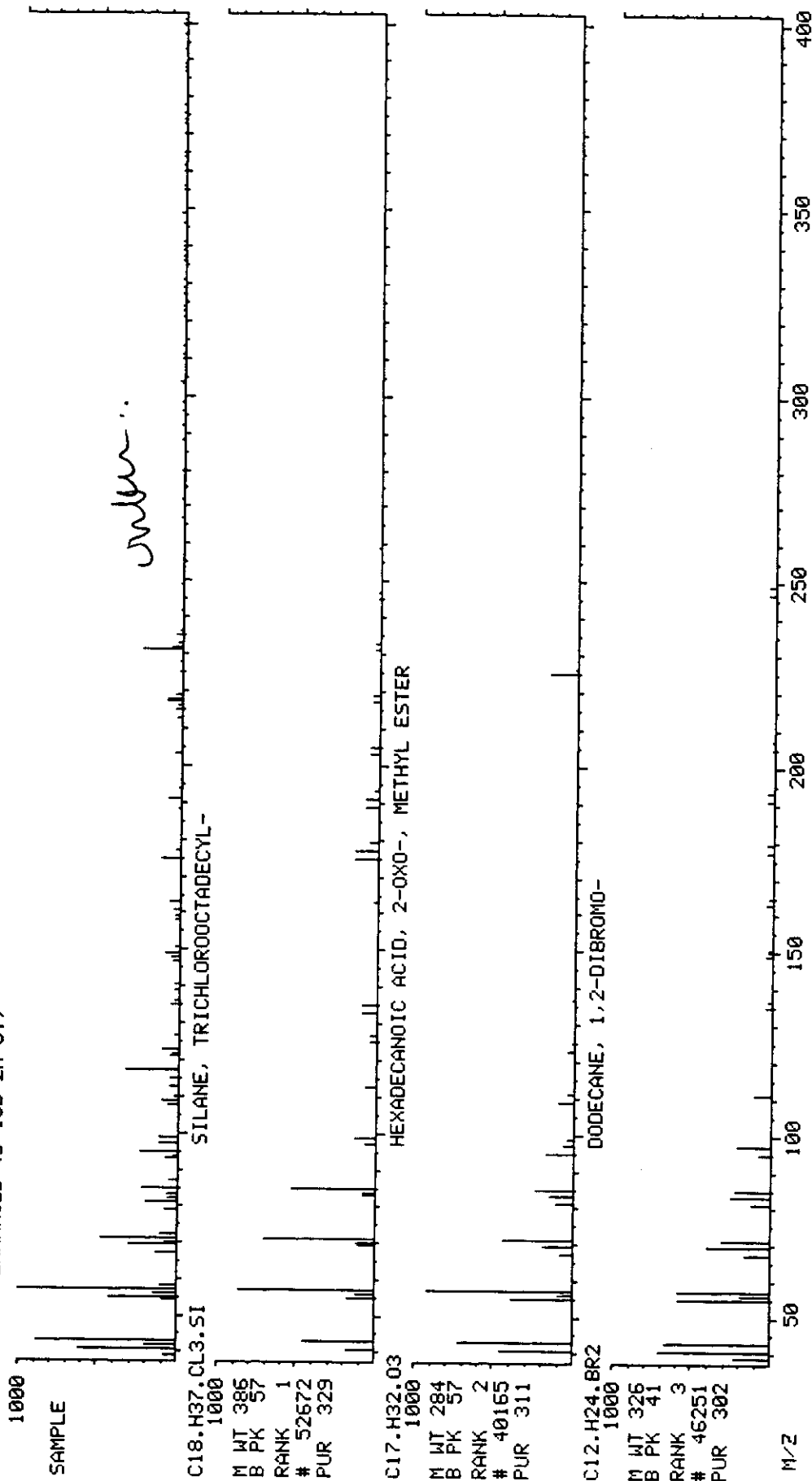
08/31/98 21:30:00 + 20:07

SAMPLE: T-MM5-3 1/35A/1ML

CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M

ENHANCED (S 158 2N 0T)

DATA: 30058105 #1809
CALI: 30058105 # 3
BASE M/Z: 57
RIC: 18550.



TIC SELECTION REPORT

DATA FILE: 30068106

THE FOLLOWING PEAKS WERE REJECTED BECAUSE
AT LEAST 40 % OF THEIR SIZE WAS ACCOUNTED FOR BY
TARGET COMPOUNDS ELUTING WITHIN 4 SCANS OF THE
PEAK TOP.

SCAN	SIZE	AMOUNT
198	280160.	30.806
341	222336.	24.448
422	289280.	31.809
531	426880.	46.939
830	363776.	40.000
1088	480256.	52.808
1386	389632.	50.148
1386	383744.	49.390
1553	310784.	40.000

TOTAL NUMBER OF UNIDENTIFIED PEAKS WITH SIZE
GREATER THAN 10 % OF THE CLOSEST INTERNAL STANDARD
THAT DOES NOT HAVE INTERFERENCES = 123

INTERNAL STANDARDS THAT HAVE RIC SIZE LESS THAN
50 % OR GREATER THAN 200 % OF THE ESTIMATED RIC SIZE
ARE CONSIDERED TO HAVE INTERFERENCES AND WILL NOT BE USED
FOR QUANTITATION.

#	INTERNAL STANDARD	RIC SIZE	PERCENT OF ESTIMATED	
			RIC SIZE	SCAN
1	C140 NAPHTHALENE-D8	426880.	222 *	531
2	C150 ACENAPHTHENE-D10	363776.	163	830
3	C160 PHENANTHRENE-D10	480256.	207 *	1088
4	C170 CHRYSENE-D12	310784.	124	1553
5	C175 PERYLENE-D12	270816.	105	1784

* INDICATES INTERFERENCE

SIZE = AREA

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 300681RI06
Std Id: ST16980902

Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 09/02/98 20:29
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Surrogate Spike Recoveries
8270-G, LIMS 10/10/96

Surrogate	Surrogate Amount (ug)		% Recovery	
	Spiked	Measured	Measured	QC limits
CS20 NITROBENZENE-D5	50.00	39.60	79.2	45 107
CS25 2-FLUOROBIPHENY	50.00	54.20	108.	62 110
CS30 TERPHENYL-D14	50.00	41.40	82.8	58 135
CS45 PHENOL-D5	100.0	63.90	63.9	43 130
CS50 2-FLUOROPHENOL	100.0	57.20	57.2	36 111
CS55 2,4,6-TRIBROMOP	100.0	49.30	49.3	* 58 131

confirming

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
CS70 2-CHLOROPHENOL-D4		ND	UG/A	30.0
CS75 1,2-DICHLOROBNZN-D4	342	130.	UG/A	30.0
HEXACHLOROBENZENE-C13		ND	UG/A	30.0
C310 N-NITROSODIMETHYLAMINE		ND	UG/A	30.0
PYRIDINE		ND	UG/A	60.0
2-PICOLINE		ND	UG/A	30.0
N-NITROSOMETHYLETHYLAMINE		ND	UG/A	30.0
METHYLMETHANESULFONATE		ND	UG/A	30.0
N-NITROSODIETHYLAMINE		ND	UG/A	30.0
ETHYLMETHANE SULFONATE		ND	UG/A	30.0
PENTACHLOROETHANE		ND	UG/A	30.0
C320 ANILINE		ND	UG/A	30.0
C315 PHENOL	313	49.	UG/A	30.0
C325 BIS(2-CHLOROETHYL) ETHER		ND	UG/A	30.0
C330 2-CHLOROPHENOL		ND	UG/A	30.0
C335 1,3-DICHLOROBENZENE		ND	UG/A	30.0
C340 1,4-DICHLOROBENZENE		ND	UG/A	30.0
C345 BENZYL ALCOHOL		ND	UG/A	30.0
C350 1,2-DICHLOROBENZENE		ND	UG/A	30.0
C355 2-METHYLPHENOL		ND	UG/A	30.0
C360 2,2'-OXYBIS(1-CLPROPAN)		ND	UG/A	30.0
C361 ACETOPHENONE	401	30.	UG/A	30.0
N-NITROSOPYRROLIDINE		ND	UG/A	30.0
N-NITROSOMORPHOLINE		ND	UG/A	30.0
3-METHYL PHENOL		ND	UG/A	30.0
C365 4-METHYLPHENOL		ND	UG/A	30.0
C370 N-NITROSO-DI-N-PROPYLAM	423	31.	UG/A	30.0
O-TOLUIDINE		ND	UG/A	30.0
C375 HEXACHLOROETHANE		ND	UG/A	30.0

1098

Reviewed by: _____

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 300681RI06
Std Id: ST16980902

Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 09/02/98 20:29
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
=====				
C410 NITROBENZENE		ND	UG/A	30.0
N-NITROSOPIPERIDINE		ND	UG/A	30.0
C415 ISOPHORONE		ND	UG/A	30.0
C420 2-NITROPHENOL		ND	UG/A	30.0
C425 2,4-DIMETHYLPHENOL		ND	UG/A	30.0
C435 BIS(2-CHLOROETHOXY)METH		ND	UG/A	30.0
C440 2,4-DICHLOROPHENOL		ND	UG/A	30.0
C430 BENZOIC ACID		ND	UG/A	150.0
AA-DIMETHYLPHENETHYLAMINE		ND	UG/A	30.0
C445 1,2,4-TRICHLOROBENZENE		ND	UG/A	30.0
C450 NAPHTHALENE	535	42.	UG/A	30.0
C455 4-CHLOROANALINE		ND	UG/A	30.0
2,6-DICHLOROPHENOL		ND	UG/A	30.0
HEXACHLOROPROPENE		ND	UG/A	30.0
C460 HEXACHLOROBUTADIENE		ND	UG/A	30.0
P-PHENYLENE DIAMINE		ND	UG/A	30.0
N-NITROSODI-N-BUTYLAMINE		ND	UG/A	30.0
C465 4-CHLORO-3-METHYLPHENO		ND	UG/A	30.0
SAFROLE		ND	UG/A	30.0
C470 2-METHYLNAPHTHALENE	654	77.	UG/A	30.0
1,2,4,5-TETRACHLOROBENZENE		ND	UG/A	30.0
ISOSAFROLE (#1)		ND	UG/A	60.0
C510 HEXACHLOROCYCLOPENTADI		ND	UG/A	30.0
C515 2,4,6-TRICHLOROPHENOL		ND	UG/A	30.0
C520 2,4,5-TRICHLOROPHENOL		ND	UG/A	30.0
ISOSAFROLE (#2)		ND	UG/A	60.0
C525 2-CHLORONAPHTHALENE		ND	UG/A	30.0
1-CHLORONAPHTHALENE		ND	UG/A	15.0
C530 2-NITROANALINE		ND	UG/A	150.0
1,4-NAPHTHOQUINONE		ND	UG/A	30.0
C535 DIMETHYLPHTHALATE		ND	UG/A	30.0
1,3-DINITROBENZENE		ND	UG/A	30.0
C540 ACENAPHTHYLENE		ND	UG/A	30.0
C543 2,6-DINITROTOLUENE		ND	UG/A	150.0
C545 3-NITROANILINE		ND	UG/A	30.0
C550 ACENAPHTHENE		ND	UG/A	150.0
C555 2,4-DINITROPHENOL		ND	UG/A	30.0
C565 DIBENZOFURAN		ND	UG/A	150.0
C560 4-NITROPHENOL		ND	UG/A	30.0
PENTACHLOROBENZENE		ND	UG/A	30.0
C570 2,4-DINITROTOLUENE		ND	UG/A	30.0
1-NAPHTHYLAMINE		ND	UG/A	30.0
2-NAPHTHYLAMINE		ND	UG/A	30.0
2,3,4,6-TETRACHLOROPHENOL		ND	UG/A	60.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 300681RI06
Std Id: ST16980902

Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 09/02/98 20:29
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C580 DIETHYLPHTHALATE		ND	UG/A	30.0
C590 FLUORENE		ND	UG/A	30.0
C585 4-CHLOROPHENYL-PHENYLE		ND	UG/A	30.0
5-NITRO-O-TOLUIDINE		ND	UG/A	30.0
C595 4-NITROANALINE		ND	UG/A	150.0
C610 4,6-DINITRO-2-METHYLPH		ND	UG/A	150.0
C615 N-NITROSODIPHENYLAMINE		ND	UG/A	30.0
C620 AZOBENZENE		ND	UG/A	30.0
SYM-TRINITROBENZENE		ND	UG/A	30.0
C625 4-BROMOPHENYL-PHENYLET		ND	UG/A	30.0
PHENACETIN		ND	UG/A	30.0
DIALATE	234	ND	UG/A	30.0
C630 HEXACHLOROBENZENE		ND	UG/A	30.0
4-AMINOBIIPHENYL		ND	UG/A	30.0
C635 PENTACHLOROPHENOL		ND	UG/A	150.0
PRONAMIDE		ND	UG/A	30.0
PENTACHLORONITROBENZENE		ND	UG/A	150.0
C640 PHENANTHRENE		ND	UG/A	30.0
C645 ANTHRACENE		ND	UG/A	30.0
2SECBUTYL-4,6-DINITROPHENOL		ND	UG/A	30.0
C647 CARBAZOLE		ND	UG/A	30.0
C650 DI-N-BUTYLPHTHALATE		ND	UG/A	30.0
4-NITROQUINOLINE-1-OXIDE		ND	UG/A	30.0
METHAPYRILENE		ND	UG/A	30.0
ISODRIN		ND	UG/A	30.0
C655 FLUORANTHENE		ND	UG/A	30.0
CHLOROBENZILATE		ND	UG/A	30.0
C710 BENZIDINE		ND	UG/A	300.0
C715 PYRENE		ND	UG/A	30.0
ARAMITE (#1)		ND	UG/A	30.0
ARAMITE (#2)		ND	UG/A	30.0
P-DIMETHYLAMINOAZOBENZENE		ND	UG/A	30.0
3,3'-DIMETHYLBENZIDINE		ND	UG/A	30.0
KEPONE		ND	UG/A	150.0
C720 BUTYLBENZYLPHTHALATE		ND	UG/A	30.0
2-ACETYLAMINOFLUORENE		ND	UG/A	30.0
C730 BENZO (A) ANTHRACENE		ND	UG/A	30.0
C725 3,3'-DICHLOROBENZIDINE		ND	UG/A	60.0
C740 CHRYSENE		ND	UG/A	30.0
C745 BIS (2-ETHYLHEXYL) PHTHA		ND	UG/A	30.0
3-METHYLCHOLANTHRENE		ND	UG/A	30.0
C760 DI-N-OCTYL PHTHALATE		ND	UG/A	30.0
C765 BENZO (B) FLUORANTHENE		ND	UG/A	30.0
7,12-DIMETHYLBENZANTHRACENE		ND	UG/A	30.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

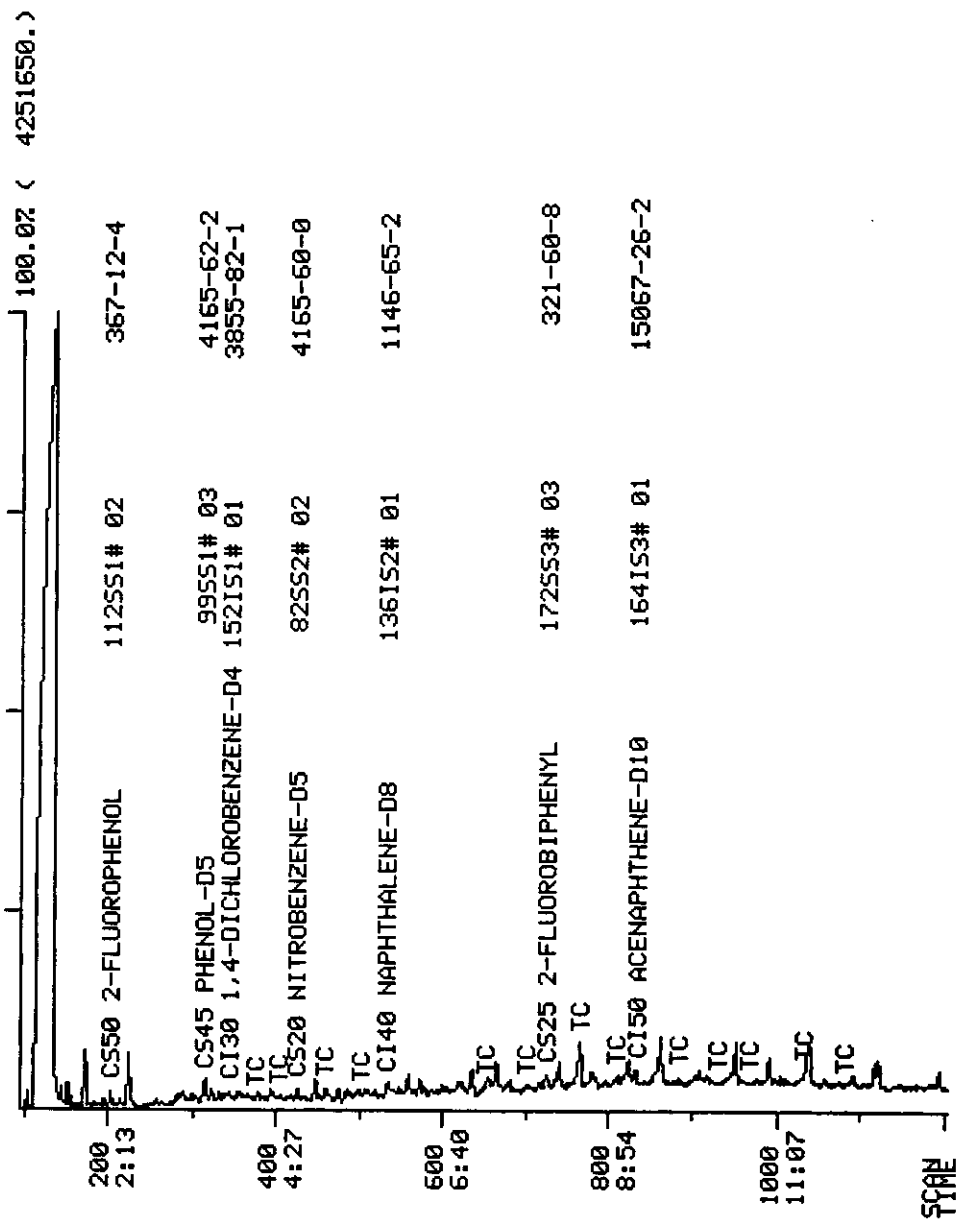
Data File: 300681RI06
Std Id: ST16980902

Sample: T-MM5-3 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 09/02/98 20:29
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

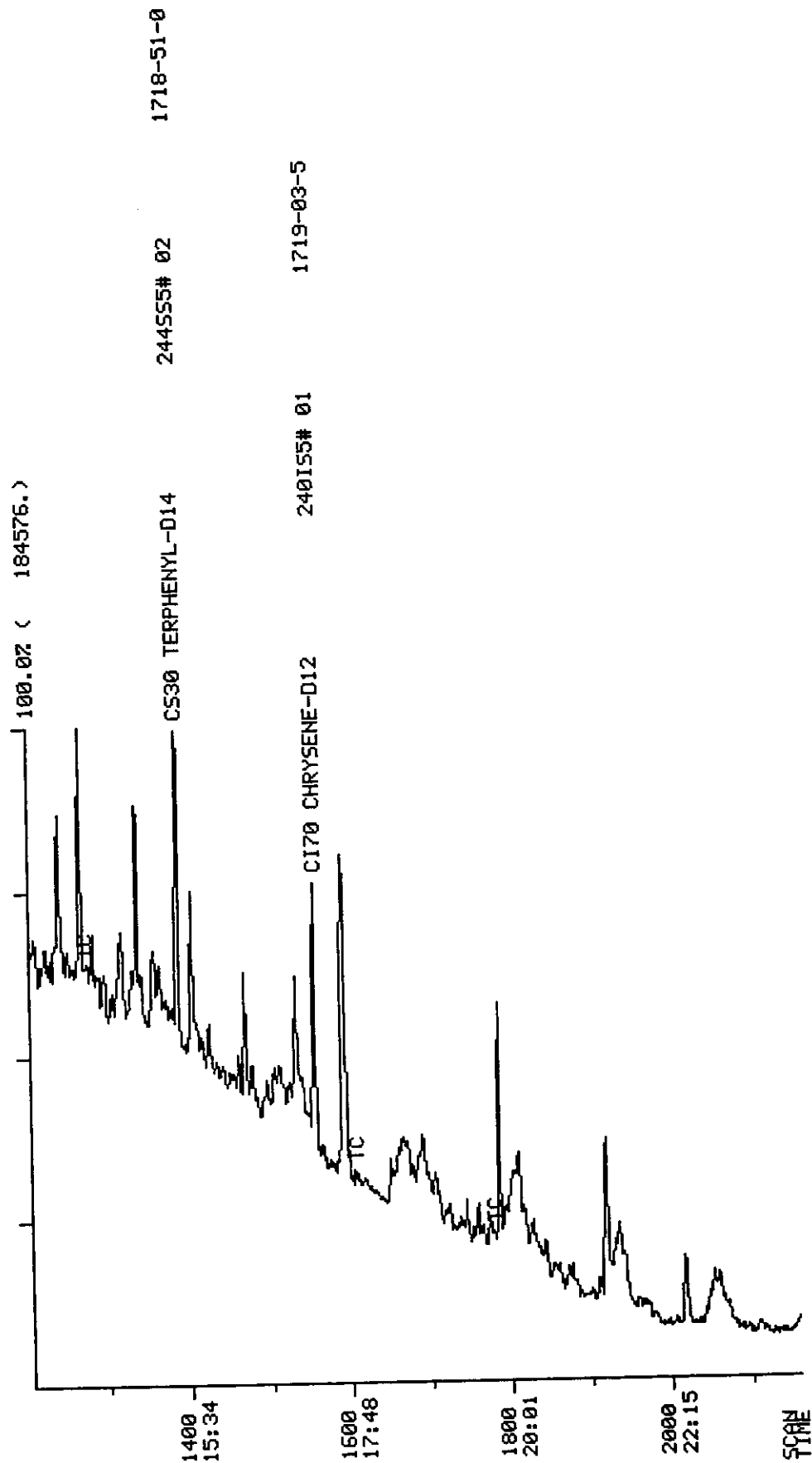
Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C770 BENZO (K) FLUORANTHENE		ND	UG/A	30.0
HEXACHLOROPHENE	1777	40.	UG/A	30.0
C775 BENZO (A) PYRENE		ND	UG/A	30.0
C780 INDENO (1,2,3-CD) PYRENE		ND	UG/A	30.0
C785 DIBENZ (A,H) ANTHRACENE		ND	UG/A	30.0
C790 BENZO (G,H,I) PERYLENE		ND	UG/A	30.0

DATA FROM FILE: 300681R106 SCANS 94 TO 1204 ACQUIRED: 09/02/98 20:29:00
 SAMPLE: T-MM5-3 1/35A/1ML CALI: 300681R106 #3
 COND.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M INST. ID: F16



DATA FROM FILE: 300681RI06 SCANS 1204 TO 2158 ACQUIRED: 09/02/98 20:29:00
 CALI: 300681RI06 #3
 SAMPLE: T-MM5-3 1/35A/1ML INST. ID: F16
 CONDS.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M



ADDITIONAL STANDARD USED FOR THIS REPORT IS: ST16980902A

QUANTERRA QUANTITATION SUMMARY

File: 300681RI06

Sample: T-MM5-3 1/3SA/1ML INST. ID: F16

Analyst: DAT

Instrument Id: F16

Analyzed: 09/02/98 20:29

Compounds with amounts less than 0.20 reported as NOT FOUND

Library									Units: UG/ML
No	Entry	Name	Mass	Meth	Scan	Ref	Area	RRF(L)	Amount
1	S1# 1	CI30 1,4-DICHLOROBENZENE	152	A BV	342	1	27901.	1.000	40.000
2	S2# 1	CI40 NAPHTHALENE-D8	136	A BB	531	2	94024.	1.000	40.000
3	S3# 1	CI50 ACENAPHTHENE-D10	164	A BB	831	3	49022.	1.000	40.000
4	S4# 1	CI60 PHENANTHRENE-D10	188	A BB	1089	4	82021.	1.000	40.000
5	S5# 1	CI70 CHRYSENE-D12	240	A BB	1554	5	80400.	1.000	40.000
6	S6# 1	CI75 PERYLENE-D12	264	A VB	1785	6	74618.	1.000	40.000
7	S2# 2	CS20 NITROBENZENE-D5	82	A BV	423	2	43753.	0.470	39.640
8	S3# 3	CS25 2-FLUOROBIPHENYL	172	A BB	725	3	82055.	1.234	54.245
9	S5# 2	CS30 TERPHENYL-D14	244	A BB	1386	5	82840.	0.995	41.412
10	S1# 3	CS45 PHENOL-D5	99	A BB	312	1	90260.	2.025	63.885
11	S1# 2	CS50 2-FLUOROPHENOL	112	A BV	200	1	54614.	1.368	57.224
12	S3# 2	CS55 2,4,6-TRIBROMOPHENO	330	A BB	973	3	14231.	0.236	49.275
13	S1# 5	CS70 2-CHLOROPHENOL-D4	132		1	NOT FOUND			
14	S1# 6	CS75 1,2-DICHLOROBNZN-D4	152	A BV	342	1	27901.	0.920	43.476
15	S4# 4	HEXACHLOROBENZENE-C13	294		4	NOT FOUND			
16	S1# 15	C310 N-NITROSODIMETHYLAM	74		1	NOT FOUND			
17	S1# 20	PYRIDINE	79		1	NOT FOUND			
18	S1# 30	2-PICOLINE	93		1	NOT FOUND			
19	S1# 40	N-NITROSOMETHYLETHYLAMIN	42		1	NOT FOUND			
20	S1# 60	METHYLMETHANESULFONATE	80		1	NOT FOUND			
21	S1# 70	N-NITROSODIETHYLAMINE	102		1	NOT FOUND			
22	S1# 85	ETHYLMETHANE SULFONATE	79		1	NOT FOUND			
23	S1# 95	PENTACHLOROETHANE	117		1	NOT FOUND			
24	S1#100	C320 ANILINE	93		1	NOT FOUND			
25	S1#105	C315 PHENOL	94	A BB	313	1	22451.	1.968	16.356
26	S1#110	C325 BIS(2-CHLOROETHYL)E	93	A BB	320	1	3280.	1.368	3.438
27	S1#115	C330 2-CHLOROPHENOL	128	A VB	320	1	396.	1.497	0.379
28	S1#125	C335 1,3-DICHLOROBENZENE	146		1	NOT FOUND			
29	S1#130	C340 1,4-DICHLOROBENZENE	146	A BB	344	1	388.	1.617	0.344
30	S1#145	C345 BENZYL ALCOHOL	108	A VB	372	1	5339.	0.854	8.959
31	S1#150	C350 1,2-DICHLOROBENZENE	146		1	NOT FOUND			
32	S1#160	C355 2-METHYLPHENOL	108		1	NOT FOUND			
33	S1#165	C360 2,2'-OXYBIS(1-CLPRO	45		1	NOT FOUND			
34	S1#170	C361 ACETOPHENONE	105	A BB	401	1	13534.	1.934	10.033
35	S1#175	N-NITROSOPYRROLIDINE	100	A BB	404	1	120.	0.553	0.311
36	S1#180	N-NITROSOMORPHOLINE	56	A BV	411	1	4290.	0.821	7.494
37	S1#182	3-METHYL PHENOL	108		1	NOT FOUND			
38	S1#185	C365 4-METHYLPHENOL	108		1	NOT FOUND			
39	S1#190	C370 N-NITROSO-DI-N-PROP	70	A BV	423	1	6576.	0.906	10.407
40	S1#195	O-TOLUIDINE	106		1	NOT FOUND			
41	S1#200	C375 HEXACHLOROETHANE	117		1	NOT FOUND			
42	S2# 10	C410 NITROBENZENE	77		2	NOT FOUND			
43	S2# 15	N-NITROSOPIPERIDINE	42	A BB	457	2	1528.	0.320	2.028
44	S2# 20	C415 ISOPHORONE	82		2	NOT FOUND			
45	S2# 25	C420 2-NITROPHENOL	139		2	NOT FOUND			

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46	S2#	30	C425	2,4-DIMETHYLPHENOL	107	A	BB	501	2	1507.	0.330	1.943
47	S2#	35	C435	BIS(2-CHLOROETHOXY)	93			2	NOT FOUND			
48	S2#	40	C440	2,4-DICHLOROPHENOL	162	A	BB	520	2	560.	0.308	0.775
49	S2#	45	C430	BENZOIC ACID	122			2	NOT FOUND			

50	S2# 50	AA-DIMETHYLPHENETHYLAMIN	58	A BB	707	2	2424.	1.285	0.803
51	S2# 55	C445 1,2,4-TRICHLOROBENZ	180	A BV	525	2	185.	0.331	0.238
52	S2# 60	C450 NAPHTHALENE	128	A BV	535	2	33489.	1.022	13.937
53	S2# 80	C455 4-CHLOROANALINE	127	A BB	557	2	848.	0.454	0.795
54	S2# 85	2,6-DICHLOROPHENOL	162		2	NOT FOUND			
55	S2# 90	HEXACHLOROPROPENE	213		2	NOT FOUND			
56	S2# 95	C460 HEXACHLOROBUTADIENE	225		2	NOT FOUND			
57	S2#115	P-PHENYLENE DIAMINE	108		2	NOT FOUND			
58	S2#120	N-NITROSODI-N-BUTYLAMINE	84		2	NOT FOUND			
59	S2#130	C465 4-CHLORO-3-METHYLPH	107		2	NOT FOUND			
60	S2#140	SAFROLE	162	A BB	650	2	1794.	0.248	3.081
61	S2#145	C470 2-METHYLNAPHTHALENE	142	A BV	654	2	45517.	0.751	25.775
62	S3# 10	1,2,4,5-TETRACHLOROBENZE	214		3	NOT FOUND			
63	S3# 15	ISOSAFROLE (#1)	162	A BB	699	3	296.	0.044	5.533
64	S3# 20	C510 HEXACHLOROCYCLOPENT	237		3	NOT FOUND			
65	S3# 25	C515 2,4,6-TRICHLOROPHEN	196		3	NOT FOUND			
66	S3# 30	C520 2,4,5-TRICHLOROPHEN	196		3	NOT FOUND			
67	S3# 35	ISOSAFROLE (#2)	104		3	NOT FOUND			
68	S3# 40	C525 2-CHLORONAPHTHALENE	162		3	NOT FOUND			
69	S3# 42	1-CHLORONAPHTHALEN	162		3	NOT FOUND			
70	S3# 45	C530 2-NITROANALINE	65	A BB	766	3	721.	0.502	1.173
71	S3# 50	1,4-NAPHTHOQUINONE	158		3	NOT FOUND			
72	S3# 55	C535 DIMETHYLPHTHALATE	163	A BB	806	3	905.	1.341	0.551
73	S3# 60	1,3-DINITROBENZENE	168	A BB	809	3	261.	0.197	1.078
74	S3# 65	C540 ACENAPHTHYLENE	152		3	NOT FOUND			
75	S3# 70	C543 2,6-DINITROTOLUENE	165		3	NOT FOUND			
76	S3# 75	C545 3-NITROANILINE	138		3	NOT FOUND			
77	S3# 80	C550 ACENAPHTHENE	153	A VB	836	3	1302.	1.159	0.917
78	S3# 85	C555 2,4-DINITROPHENOL	184	A VB	860	3	150.	0.189	0.647
79	S3# 90	C565 DIBENZOFURAN	168	A VB	868	3	1263.	1.669	0.617
80	S3# 95	C560 4-NITROPHENOL	109		3	NOT FOUND			
81	S3#100	PENTACHLOROBENZENE	250		3	NOT FOUND			
82	S3#105	C570 2,4-DINITROTOLUENE	165	A VB	883	3	549.	0.397	1.127
83	S3#110	1-NAPHTHYLAMINE	143		3	NOT FOUND			
84	S3#115	2-NAPHTHYLAMINE	143		3	NOT FOUND			
85	S3#120	2,3,4,6-TETRACHLOROPHENO	232	A VB	896	3	298.	0.297	0.818
86	S3#130	C580 DIETHYLPHTHALATE	149	A VB	934	3	948.	1.369	0.565
87	S3#135	C590 FLUORENE	166	A BB	928	3	4654.	1.298	2.926
88	S3#140	C585 4-CHLOROPHENYL-PHEN	204		3	NOT FOUND			
89	S3#145	5-NITRO-O-TOLUIDINE	152	A BB	943	3	3609.	0.356	8.284
90	S3#150	C595 4-NITROANALINE	138		3	NOT FOUND			
91	S4# 10	C610 4,6-DINITRO-2-METHY	198		4	NOT FOUND			
92	S4# 15	C615 N-NITROSODIPHENYLAM	169	A VB	965	4	4111.	0.551	3.641
93	S4# 20	C620 AZOBENZENE	77	A VB	968	4	512.	1.171	0.213
94	S4# 25	SYM-TRINITROBENZENE	75	A BV	1031	4	900.	0.162	2.718
95	S4# 30	C625 4-BROMOPHENYL-PHENY	248		4	NOT FOUND			
96	S4# 35	PHENACETIN	108		4	NOT FOUND			
97	S4# 37	DIALATE	234		4	NOT FOUND			
98	S4# 40	C630 HEXACHLOROBENZENE	284		4	NOT FOUND			
99	S4# 45	4-AMINOBIIPHENYL	169	A BB	1078	4	2522.	0.609	2.020
100	S4# 50	C635 PENTACHLOROPHENOL	266		4	NOT FOUND			
101	S4# 55	PRONAMIDE	173		4	NOT FOUND			
102	S4# 60	PENTACHLORONITROBENZENE	237	A VB	1063	4	53.	0.080	0.322
103	S4# 65	C640 PHENANTHRENE	178	A BB	1093	4	14391.	1.033	6.797
104	S4# 70	C645 ANTHRACENE	178		4	NOT FOUND			
105	S4# 75	2SECBUTYL-4,6-DINITROPHE	211		4	NOT FOUND			
106	S4# 80	C647 CARBAZOLE	167		4	NOT FOUND			
107	S4# 85	C650 DI-N-BUTYLPHTHALATE	149		4	NOT FOUND			
108	S4#100	4-NITROQUINOLINE-1-OXIDE	190	A BB	1241	4	114.	0.073	0.770
109	S4#105	METHAPYRILENE	58		4	NOT FOUND			
110	S4#106	ISODRIN	193	A VV	1272	4	572.	0.127	2.201

111	S4#110	C655 FLUORANTHENE	202	A BB	1301	4	410.	0.900	0.222
112	S4#120	CHLOROBENZILATE	139		4	NOT FOUND			
113	S5# 10	C710 BENZIDINE	184		5	NOT FOUND			
114	S5# 15	C715 PYRENE	202	A BB	1337	5	1903.	1.323	0.716
115	S5# 20	ARAMITE (#1)	185		5	NOT FOUND			
116	S5# 25	ARAMITE (#2)	185		5	NOT FOUND			
117	S5# 30	P-DIMETHYLAMINOAZOBENZEN	120		5	NOT FOUND			
118	S5# 35	3,3'-DIMETHYLBENZIDINE	212		5	NOT FOUND			
119	S5# 37	KEPONE	272	A BB	1508	5	53.	0.082	0.322
120	S5# 40	C720 BUTYLBENZYLPHTHALAT	149		5	NOT FOUND			
121	S5# 45	2-ACETYLAMINOFLUORENE	181		5	NOT FOUND			
122	S5# 50	C730 BENZO(A)ANTHRACENE	228	A BB	1557	5	1764.	1.125	0.780
123	S5# 55	C725 3,3'-DICHLOOROBENZID	252		5	NOT FOUND			
124	S5# 60	C740 CHRYSENE	228	A BB	1557	5	1764.	1.014	0.865
125	S5# 65	C745 BIS(2-ETHYLHEXYL) PH	149	A BB	1605	5	5223.	0.980	2.652
126	S5# 85	3-METHYLCHOLANTHRENE	268	A VV	1831	5	300.	0.548	0.273
127	S6# 10	C760 DI-N-OCTYL PHTHALAT	149		6	NOT FOUND			
128	S6# 15	C765 BENZO(B) FLUORANTHEN	252		6	NOT FOUND			
129	S6# 20	7,12-DIMETHYLBENZANTHRAC	256		6	NOT FOUND			
130	S6# 25	C770 BENZO(K) FLUORANTHEN	252		6	NOT FOUND			
131	S6# 30	HEXACHLOROPHENE	196	A BB	1777	6	101.	0.004	13.492
132	S6# 35	C775 BENZO(A) PYRENE	252	A VB	1775	6	519.	1.182	0.235
133	S6# 55	C780 INDENO(1,2,3-CD) PYR	276		6	NOT FOUND			
134	S6# 60	C785 DIBENZ(A,H)ANTHRACE	278		6	NOT FOUND			
135	S6# 65	C790 BENZO(G,H,I) PERYLEN	276		6	NOT FOUND			

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14/7/54

QUANTERRA Internal Standard Check

Standard Filename: ST16980902

Analyzed: 09/02/98 14:53

Standard	IS# 1			IS# 2			IS# 3		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980902	23319.		3.78	79260.		5.91	43605.		9.22
Upper Limit	46638.=+200		4.28	158519.=+200		6.41	87211.=+200		9.72
Lower Limit	11659.=+ 50		3.28	39630.=+ 50		5.41	21803.=+ 50		8.72
Filename									
1 300681RI06	27901.= 119		3.80	94024.= 118		5.91	49022.= 112		9.24

IS# 1 = CI30 1,4-DICHLOROBENZENE-D4

IS# 2 = CI40 NAPHTHALENE-D8

IS# 3 = CI50 ACENAPHTHENE-D10

* - indicates an I.S. is outside QC limit(s)

Standard	IS# 4			IS# 5			IS# 6		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980902	76218.		12.09	51424.		17.27	44306.		19.84
Upper Limit	152436.=+200		12.59	102848.=+200		17.77	88612.=+200		20.34
Lower Limit	38109.=+ 50		11.59	25712.=+ 50		16.77	22153.=+ 50		19.34
Filename									
1 300681RI06	82021.= 107		12.11	80400.= 156		17.29	74618.= 168		19.85

IS# 4 = CI60 PHENANTHRENE-D10

IS# 5 = CI70 CHRYSENE-D12

IS# 6 = CI75 PERYLENE-D12

* - indicates an I.S. is outside QC limit(s)

Semivolatile Organics
Method 0010/8270

Client Name: Pacific Environmental Services
Client ID: S-MM5-3-F, FH, XAD, COND, BH
LAB ID: 300681-0007-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 27 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 100

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
Phenol	ND	ug/Sample	3000	GR
bis(2-Chloroethyl)ether	ND	ug/Sample	3000	
2-Chlorophenol	ND	ug/Sample	3000	
1,3-Dichlorobenzene	ND	ug/Sample	3000	
1,4-Dichlorobenzene	ND	ug/Sample	3000	
Benzyl alcohol	ND	ug/Sample	3000	
1,2-Dichlorobenzene	ND	ug/Sample	3000	
2-Methylphenol	ND	ug/Sample	3000	
2,2'-Oxybis(1-chloropropane)	ND	ug/Sample	3000	
3/4-Methylphenol	ND	ug/Sample	3000	
N-Nitroso-di-n-propylamine	ND	ug/Sample	3000	
Hexachloroethane	ND	ug/Sample	3000	
Nitrobenzene	ND	ug/Sample	3000	
Isophorone	ND	ug/Sample	3000	
2-Nitrophenol	ND	ug/Sample	3000	
2,4-Dimethylphenol	ND	ug/Sample	3000	
Benzoic acid	ND	ug/Sample	15000	
bis(2-Chloroethoxy)-methane	ND	ug/Sample	3000	
2,4-Dichlorophenol	ND	ug/Sample	3000	
1,2,4-Trichlorobenzene	ND	ug/Sample	3000	
Naphthalene	3200	ug/Sample	3000	
4-Chloroaniline	ND	ug/Sample	3000	
Hexachlorobutadiene	ND	ug/Sample	3000	
4-Chloro-3-methylphenol	ND	ug/Sample	3000	
2-Methylnaphthalene	6300	ug/Sample	3000	
Hexachlorocyclopentadiene	ND	ug/Sample	3000	
2,4,6-Trichlorophenol	ND	ug/Sample	3000	
2,4,5-Trichlorophenol	ND	ug/Sample	15000	
2-Chloronaphthalene	ND	ug/Sample	3000	
2-Nitroaniline	ND	ug/Sample	3000	
Dimethyl phthalate	ND	ug/Sample	3000	
Acenaphthylene	ND	ug/Sample	3000	
3-Nitroaniline	ND	ug/Sample	15000	
Acenaphthene	ND	ug/Sample	3000	
2,4-Dinitrophenol	ND	ug/Sample	15000	
4-Nitrophenol	ND	ug/Sample	15000	

Note G = Reporting limit(s) raised due to matrix interference.
Note R = Reporting limit(s) raised due to sample volume limitations.
ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Mike Orbanosky

The cover letter is an integral part of this report.
Rev 230787

1107

Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: S-MM5-3-F, FH, XAD, COND, BH
LAB ID: 300681-0007-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 27 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 100

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
Dibenzofuran	ND	ug/Sample	3000	
2,4-Dinitrotoluene	ND	ug/Sample	3000	
2,6-Dinitrotoluene	ND	ug/Sample	3000	
Diethyl phthalate	ND	ug/Sample	3000	
4-Chlorophenyl phenyl ether	ND	ug/Sample	3000	
Fluorene	ND	ug/Sample	3000	
4-Nitroaniline	ND	ug/Sample	15000	
4,6-Dinitro-2-methylphenol	ND	ug/Sample	15000	
N-Nitrosodiphenylamine	ND	ug/Sample	3000	
4-Bromophenyl phenyl ether	ND	ug/Sample	3000	
Hexachlorobenzene	ND	ug/Sample	3000	
Pentachlorophenol	ND	ug/Sample	15000	
Phenanthrene	ND	ug/Sample	3000	
Anthracene	ND	ug/Sample	3000	
Di-n-butyl phthalate	ND	ug/Sample	3000	
Fluoranthene	ND	ug/Sample	3000	
Pyrene	ND	ug/Sample	3000	
Butyl benzyl phthalate	ND	ug/Sample	3000	
3,3'-Dichlorobenzidine	ND	ug/Sample	6000	
Benzo(a)anthracene	ND	ug/Sample	3000	
bis(2-Ethylhexyl)-phthalate	ND	ug/Sample	3000	
Chrysene	ND	ug/Sample	3000	
Di-n-octyl phthalate	ND	ug/Sample	3000	
Benzo(b)fluoranthene	ND	ug/Sample	3000	
Benzo(k)fluoranthene	ND	ug/Sample	3000	
Benzo(a)pyrene	ND	ug/Sample	3000	
Indeno(1,2,3-cd)pyrene	ND	ug/Sample	3000	
Dibenz(a,h)anthracene	ND	ug/Sample	3000	
Benzo(g,h,i)perylene	ND	ug/Sample	3000	
Acetophenone	ND	ug/Sample	3000	
4-Aminobiphenyl	ND	ug/Sample	15000	
Aniline	ND	ug/Sample	3000	
Benzidine	ND	ug/Sample	30000	
3,3'-Dimethylbenzidine	ND	ug/Sample	6000	
N-Nitrosodimethylamine	ND	ug/Sample	3000	
N-Nitrosomorpholine	ND	ug/Sample	3000	
Pentachloronitrobenzene (PCNB)	ND	ug/Sample	15000	
o-Toluidine	ND	ug/Sample	6000	

ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Mike Orbanosky

The cover letter is an integral part of this report.

Rev 230787

1108

Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: S-MM5-3-F, FH, XAD, COND, BH
LAB ID: 300681-0007-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 27 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 100

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
2-Methoxybenzenamine	ND	ug/Sample	--	
Biphenyl	ND	ug/Sample	--	
Chloroacetophenone	ND	ug/Sample	--	
Cumene	ND	ug/Sample	--	
DBCP (1,2-Dibromo-3-chloropropane)	ND	ug/Sample	--	
Benzo(e)pyrene	ND	ug/Sample	--	
N-N-Diethylaniline	ND	ug/Sample	--	
Dimethylaniline	ND	ug/Sample	--	
3,3'-Dimethoxybenzidine	ND	ug/Sample	--	
Hydroquinone	ND	ug/Sample	--	
4,4'-Methyl-bis(2-chloroaniline)	ND	ug/Sample	--	
4-Nitrodiphenyl	ND	ug/Sample	--	
Trifluralin	ND	ug/Sample	--	

Surrogate	Recovery	Acceptable Range	
Nitrobenzene-d5	ND %	45 - 107	H
2-Fluorobiphenyl	ND %	62 - 110	
Terphenyl-d14	ND %	58 - 135	
Phenol-d5	ND %	43 - 130	
2-Fluorophenol	ND %	36 - 111	
2,4,6-Tribromophenol	ND %	58 - 131	

Note H = Spiked analyte not detected because of required sample dilution.
ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Mike Orbanosky

The cover letter is an integral part of this report.
Rev 230787

1108

Semivolatiles Library Search (20 Compound TID)
Method 8270

Client Name: Pacific Environmental Services

Client ID: S-MM5-3-F, FH, XAD, COND, BH

LAB ID: 300681-0007-SA

Matrix: AIRTRAIN

Authorized: 30 JUL 98

Sampled: 27 JUL 98

Prepared: NA

Received: 30 JUL 98

Analyzed: 31 AUG 98

Dilution Factor: 100

Parameter	Result	Units	Reporting Limit	Qualifier
Undecane	10000	ug/Sample	--	0
Decane, 2,5,6-trimethyl-	12000	ug/Sample	--	0
Undecane, 2,6-dimethyl-	9900	ug/Sample	--	0
Unknown	11000	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	19000	ug/Sample	--	0
Unknown	13000	ug/Sample	--	0
Decane, 2,4-dimethyl	19000	ug/Sample	--	0
Unknown	9100	ug/Sample	--	0
Unknown alkane	22000	ug/Sample	--	
Unknown alkane	36000	ug/Sample	--	
Naphthalene, 1,3-dimethyl-	16000	ug/Sample	--	0
Heptadecane, 2,6,10,14 -tetramethyl-	26000	ug/Sample	--	0
Dodecane, 3-methyl-	35000	ug/Sample	--	0
Unknown alkane	11000	ug/Sample	--	0
Nonadecane	35000	ug/Sample	--	0
Nonadecane	22000	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	46000	ug/Sample	--	0
Unknown alkane	14000	ug/Sample	--	0
Nonadecane	17000	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	10000	ug/Sample	--	0

Note 0 = Or structurally similar compound (isomer).

NA = Not Applicable

Reported By: Emily Uebelhoer

Approved By: Mike Orbanosky

The cover letter is an integral part of this report.
Rev 230787

1110

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068107
Std Id: ST16980831

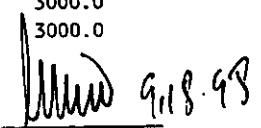
Sample: S-MM5-3 1/3SA/100M INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 21:59
Analyst: DAT Instrument ID: F16 Run Factor: 300.
Quan List Threshold: 0.95 Surrogate Vol.: 100.

Surrogate Spike Recoveries
8270-G, LIMS 10/10/96

Surrogate	Surrogate Amount (ug)		% Recovery		QC limits
	Spiked	Measured	Measured		
CS20 NITROBENZENE-D5	50.0	BDL	0*	H 45	107
CS25 2-FLUOROBIPHENY	50.00	95.10	190.	*	62 110
CS30 TERPHENYL-D14	50.00	101.0	202.	*	58 135
CS45 PHENOL-D5	100.	BDL	0*	43	130
CS50 2-FLUOROPHENOL	100.0	153.0	153.	*	36 111
CS55 2,4,6-TRIBROMOP	100.0	123.0	123.		58 131

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
HEXACHLOROBENZENE-C13		ND	UG/A	3000.0
C310 N-NITROSODIMETHYLAMINE		ND	UG/A	3000.0
PYRIDINE		ND	UG/A	6000.0
2-PICOLINE		ND	UG/A	3000.0
N-NITROSOMETHYLETHYLAMINE		ND	UG/A	3000.0
METHYLMETHANESULFONATE		ND	UG/A	3000.0
N-NITROSODIETHYLAMINE		ND	UG/A	3000.0
ETHYLMETHANE SULFONATE		ND	UG/A	3000.0
PENTACHLOROETHANE		ND	UG/A	3000.0
C320 ANILINE		ND	UG/A	3000.0
C315 PHENOL		ND	UG/A	3000.0
C325 BIS(2-CHLOROETHYL)ETHER		ND	UG/A	3000.0
C330 2-CHLOROPHENOL		ND	UG/A	3000.0
C335 1,3-DICHLOROBENZENE		ND	UG/A	3000.0
C340 1,4-DICHLOROBENZENE		ND	UG/A	3000.0
C345 BENZYL ALCOHOL		ND	UG/A	3000.0
C350 1,2-DICHLOROBENZENE		ND	UG/A	3000.0
C355 2-METHYLPHENOL		ND	UG/A	3000.0
C360 2,2'-OXYBIS(1-CLPROPAN)		ND	UG/A	3000.0
C361 ACETOPHENONE		ND	UG/A	3000.0
N-NITROSOPYRROLIDINE		ND	UG/A	3000.0
N-NITROSOMORPHOLINE		ND	UG/A	3000.0
3-METHYL PHENOL		ND	UG/A	3000.0
C365 4-METHYLPHENOL		ND	UG/A	3000.0
C370 N-NITROSO-DI-N-PROPYLAM		ND	UG/A	3000.0
O-TOLUIDINE		ND	UG/A	3000.0
C375 HEXACHLOROETHANE		ND	UG/A	3000.0
C410 NITROBENZENE		ND	UG/A	3000.0
N-NITROSOPIPERIDINE		ND	UG/A	3000.0

Reviewed by:  9.18.98

1111

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068107
Std Id: ST16980831

Sample: S-MM5-3 1/3SA/100M INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 21:59
Analyst: DAT Instrument ID: F16 Run Factor: 300.
Quan List Threshold: 0.95 Surrogate Vol.: 100.

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C415 ISOPHORONE		ND	UG/A	3000.0
C420 2-NITROPHENOL		ND	UG/A	3000.0
C425 2,4-DIMETHYLPHENOL		ND	UG/A	3000.0
C435 BIS(2-CHLOROETHOXY)METH		ND	UG/A	3000.0
C440 2,4-DICHLOROPHENOL		ND	UG/A	3000.0
C430 BENZOIC ACID		ND	UG/A	15000.0
AA-DIMETHYLPHENETHYLAMINE		ND	UG/A	3000.0
C445 1,2,4-TRICHLOROBENZENE		ND	UG/A	3000.0
C450 NAPHTHALENE	533	3200.	UG/A	3000.0
C455 4-CHLOROANILINE		ND	UG/A	3000.0
2,6-DICHLOROPHENOL		ND	UG/A	3000.0
HEXACHLOROPROPENE		ND	UG/A	3000.0
C460 HEXACHLOROBUTADIENE		ND	UG/A	3000.0
P-PHENYLENE DIAMINE		ND	UG/A	3000.0
N-NITROSODI-N-BUTYLAMINE		ND	UG/A	3000.0
C465 4-CHLORO-3-METHYLPHENO		ND	UG/A	3000.0
SAFROLE		ND	UG/A	3000.0
C470 2-METHYLNAPHTHALENE	652	6300.	UG/A	3000.0
1,2,4,5-TETRACHLOROBENZENE		ND	UG/A	3000.0
ISOSAFROLE (#1)		ND	UG/A	6000.0
C510 HEXACHLOROCYCLOPENTADI		ND	UG/A	3000.0
C515 2,4,6-TRICHLOROPHENOL		ND	UG/A	3000.0
C520 2,4,5-TRICHLOROPHENOL		ND	UG/A	3000.0
ISOSAFROLE (#2)		ND	UG/A	6000.0
C525 2-CHLORONAPHTHALENE		ND	UG/A	3000.0
1-CHLORONAPHTHALENE		ND	UG/A	1500.0
C530 2-NITROANILINE		ND	UG/A	15000.0
1,4-NAPHTHOQUINONE		ND	UG/A	3000.0
C535 DIMETHYLPHTHALATE		ND	UG/A	3000.0
1,3-DINITROBENZENE		ND	UG/A	3000.0
C540 ACENAPHTHYLENE		ND	UG/A	3000.0
C543 2,6-DINITROTOLUENE		ND	UG/A	3000.0
C545 3-NITROANILINE		ND	UG/A	15000.0
C550 ACENAPHTHENE		ND	UG/A	3000.0
C555 2,4-DINITROPHENOL		ND	UG/A	15000.0
C565 DIBENZOFURAN		ND	UG/A	3000.0
C560 4-NITROPHENOL		ND	UG/A	15000.0
PENTACHLOROBENZENE		ND	UG/A	3000.0
C570 2,4-DINITROTOLUENE		ND	UG/A	3000.0
1-NAPHTHYLAMINE		ND	UG/A	3000.0
2-NAPHTHYLAMINE		ND	UG/A	3000.0
2,3,4,6-TETRACHLOROPHENOL		ND	UG/A	6000.0
C580 DIETHYLPHTHALATE		ND	UG/A	3000.0
C590 FLUORENE		ND	UG/A	3000.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068107
Std Id: ST16980831

Sample: S-MM5-3 1/3SA/100M INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 21:59
Analyst: DAT Instrument ID: F16 Run Factor: 300.
Quan List Threshold: 0.95 Surrogate Vol.: 100.

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C585 4-CHLOROPHENYL-PHENYLE		ND	UG/A	3000.0
5-NITRO-O-TOLUIDINE		ND	UG/A	3000.0
C595 4-NITROANALINE		ND	UG/A	15000.0
C610 4,6-DINITRO-2-METHYLPH		ND	UG/A	15000.0
C615 N-NITROSODIPHENYLAMINE		ND	UG/A	3000.0
C620 AZOBENZENE		ND	UG/A	3000.0
SYM-TRINITROBENZENE		ND	UG/A	3000.0
C625 4-BROMOPHENYL-PHENYLET		ND	UG/A	3000.0
PHENACETIN		ND	UG/A	3000.0
DIALATE	234	ND	UG/A	3000.0
C630 HEXACHLOROBENZENE		ND	UG/A	3000.0
4-AMINOBIIPHENYL		ND	UG/A	3000.0
C635 PENTACHLOROPHENOL		ND	UG/A	15000.0
PRONAMIDE		ND	UG/A	3000.0
PENTACHLORONITROBENZENE		ND	UG/A	15000.0
C640 PHENANTHRENE		ND	UG/A	3000.0
C645 ANTHRACENE		ND	UG/A	3000.0
2SECBUTYL-4,6-DINITROPHENOL		ND	UG/A	3000.0
C647 CARBAZOLE		ND	UG/A	3000.0
C650 DI-N-BUTYLPHTHALATE		ND	UG/A	3000.0
4-NITROQUINOLINE-1-OXIDE		ND	UG/A	3000.0
METHAPYRILENE		ND	UG/A	3000.0
ISODRIN		ND	UG/A	3000.0
C655 FLUORANTHENE		ND	UG/A	3000.0
CHLOROBENZILATE		ND	UG/A	3000.0
C710 BENZIDINE		ND	UG/A	30000.0
C715 PYRENE		ND	UG/A	3000.0
ARAMITE (#1)		ND	UG/A	3000.0
ARAMITE (#2)		ND	UG/A	3000.0
P-DIMETHYLAMINOAZOBENZENE		ND	UG/A	3000.0
3,3'-DIMETHYLBENZIDINE		ND	UG/A	3000.0
KEPONE		ND	UG/A	15000.0
C720 BUTYLBENZYLPHTHALATE		ND	UG/A	3000.0
2-ACETYLAMINOFLUORENE		ND	UG/A	3000.0
C730 BENZO(A)ANTHRACENE		ND	UG/A	3000.0
C725 3,3'-DICHLOROBENZIDINE		ND	UG/A	6000.0
C740 CHRYSENE		ND	UG/A	3000.0
C745 BIS(2-ETHYLHEXYL)PHTHA		ND	UG/A	3000.0
3-METHYLCHOLANTHRENE		ND	UG/A	3000.0
C760 DI-N-OCTYL PHTHALATE		ND	UG/A	3000.0
C765 BENZO(B)FLUORANTHENE		ND	UG/A	3000.0
7,12-DIMETHYLBENZANTHRACENE		ND	UG/A	3000.0
C770 BENZO(K)FLUORANTHENE		ND	UG/A	3000.0
HEXACHLOROPHENE		ND	UG/A	3000.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

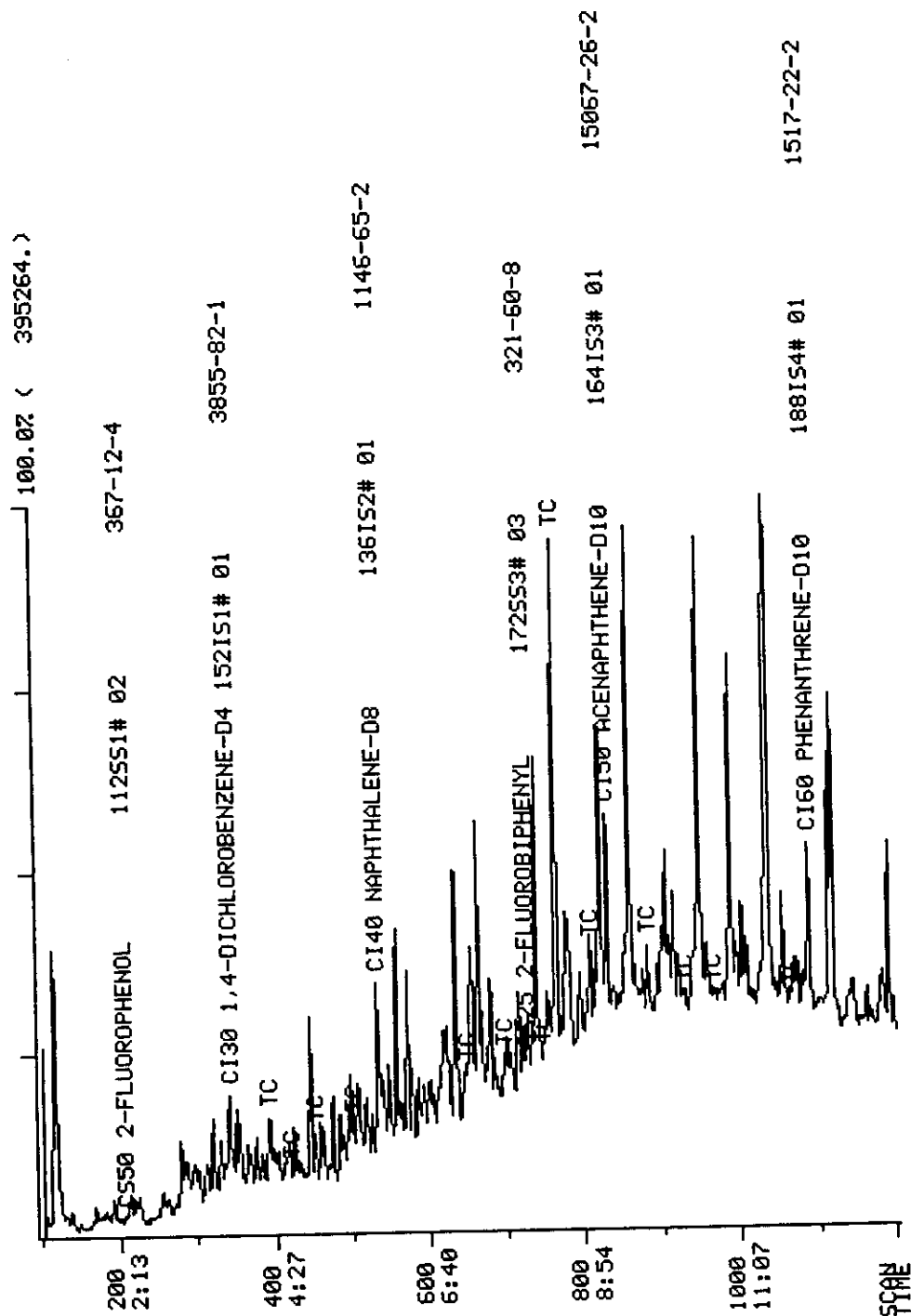
Data File: 30068107
Std Id: ST16980831

Sample: S-MM5-3 1/3SA/100M INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 21:59
Analyst: DAT Instrument ID: F16 Run Factor: 300.
Quan List Threshold: 0.95 Surrogate Vol.: 100.

Target Compounds: SAP9

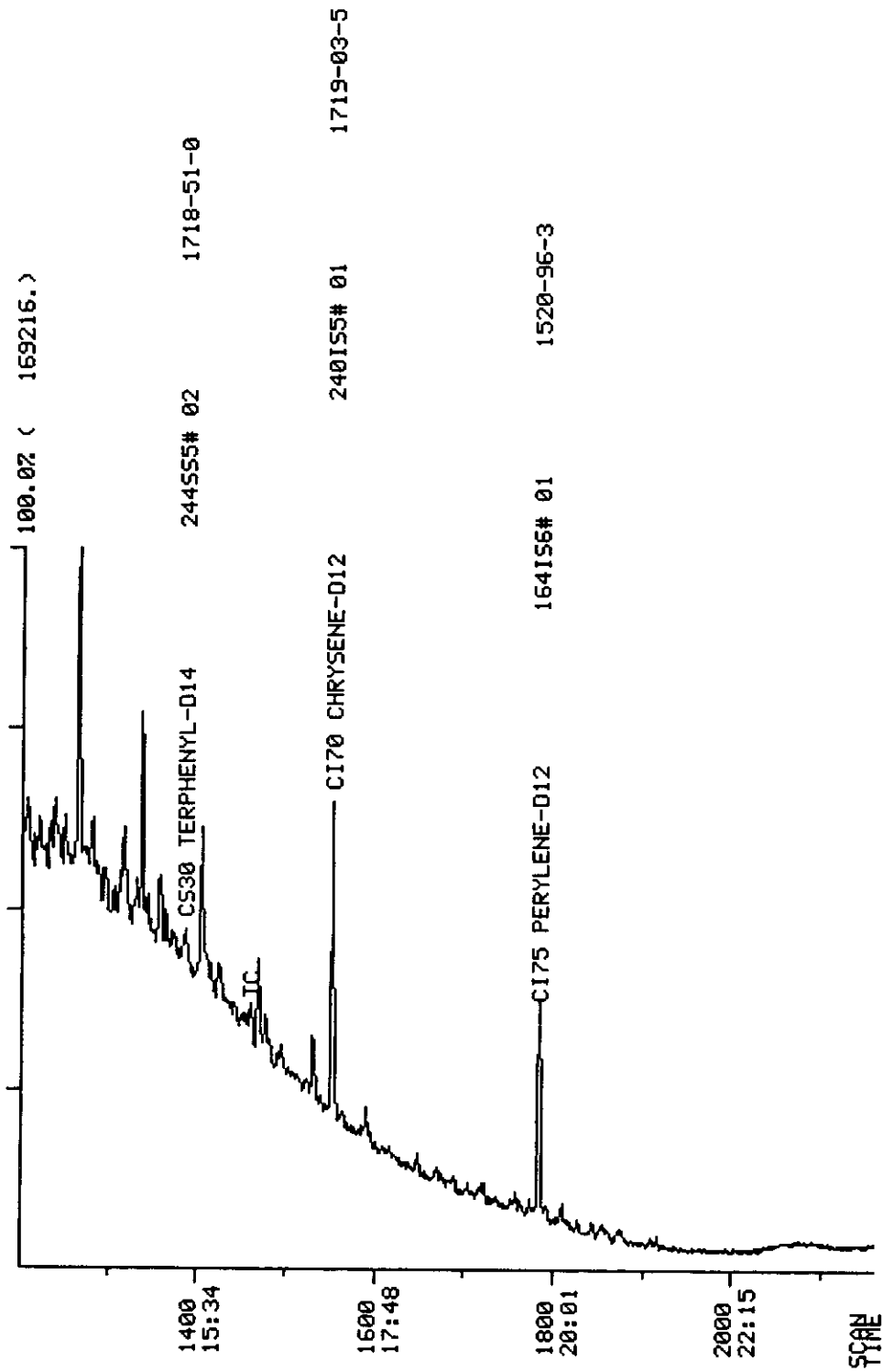
Parameter	Scan	Result	Units	Reporting Limit
C775 BENZO(A)PYRENE		ND	UG/A	3000.0
C780 INDENO(1,2,3-CD)PYRENE		ND	UG/A	3000.0
C785 DIBENZ(A,H)ANTHRACENE		ND	UG/A	3000.0
C790 BENZO(G,H,I)PERYLENE		ND	UG/A	3000.0

DATA FROM FILE: 30068107 SCANS 93 TO 1203 ACQUIRED: 08/31/98 21:59:00
 CALI: 30068107 #3
 SAMPLE: S-MMS-3 1/35A/100M INST. ID: F16
 CONDS.: UG/ML *100Z/100Z *(NA/NA >1/35A NA M



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1
1
C11

DATA FROM FILE: 30068107 SCANS 1203 TO 2158 ACQUIRED: 08/31/98 21:59:00
 CALI: 30068107 #3
 SAMPLE: S-MM5-3 1/35A/100M INST. ID: F16
 CONDS.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M



1-1-18

ADDITIONAL STANDARD USED FOR THIS REPORT IS: ST16980831A

QUANTERRA QUANTITATION SUMMARY

File: 30068107

Sample: S-MM5-3 1/3SA/100M INST. ID: F16
Analyst: DAT Instrument Id: F16 Analyzed: 08/31/98 21:59

Compounds with amounts less than 0.20 reported as NOT FOUND

Library										Units: UG/ML	
No	Entry	Name	Mass	Meth	Scan	Ref	Area	RRF(L)	Amount		
1	S1#	1 C130 1,4-DICHLOROBENZENE	152	A	BB	340	1	30514. 1.000	40.000		
2	S2#	1 C140 NAPHTHALENE-D8	136	A	BB	530	2	110082. 1.000	40.000		
3	S3#	1 C150 ACENAPHTHENE-D10	164	A	BB	829	3	60039. 1.000	40.000		
4	S4#	1 C160 PHENANTHRENE-D10	188	A	BB	1088	4	106261. 1.000	40.000		
5	S5#	1 C170 CHRYSENE-D12	240	A	BB	1552	5	67833. 1.000	40.000		
6	S6#	1 C175 PERYLENE-D12	264	A	BV	1782	6	51078. 1.000	40.000		
7	S2#	2 CS20 NITROBENZENE-D5	82			2	NOT FOUND				
8	S3#	3 CS25 2-FLUOROBIPHENYL	172	A	BB	723	3	1762. 1.234	0.951		
9	S5#	2 CS30 TERPHENYL-D14	244	A	BB	1384	5	1708. 0.995	1.012		
10	S1#	3 CS45 PHENOL-D5	99			1	NOT FOUND				
11	S1#	2 CS50 2-FLUOROPHENOL	112	A	VV	205	1	1595. 1.368	1.528		
12	S3#	2 CS55 2,4,6-TRIBROMOPHENO	330	A	BB	971	3	436. 0.236	1.233		
13	S4#	4 HEXACHLOROBENZENE-C13	294			4	NOT FOUND				
14	S1#	15 C310 N-NITROSODIMETHYLAM	74			1	NOT FOUND				
15	S1#	20 PYRIDINE	79			1	NOT FOUND				
16	S1#	30 2-PICOLINE	93			1	NOT FOUND				
17	S1#	40 N-NITROSOMETHYLETHYLAMIN	42			1	NOT FOUND				
18	S1#	60 METHYLMETHANESULFONATE	80			1	NOT FOUND				
19	S1#	70 N-NITROSODIETHYLAMINE	102			1	NOT FOUND				
20	S1#	85 ETHYLMETHANE SULFONATE	79			1	NOT FOUND				
21	S1#	95 PENTACHLOROETHANE	117	A	BB	305	1	157. 0.602	0.343		
22	S1#	100 C320 ANILINE	93			1	NOT FOUND				
23	S1#	105 C315 PHENOL	94			1	NOT FOUND				
24	S1#	110 C325 BIS(2-CHLOROETHYL)E	93			1	NOT FOUND				
25	S1#	115 C330 2-CHLOROPHENOL	128			1	NOT FOUND				
26	S1#	125 C335 1,3-DICHLOROBENZENE	146			1	NOT FOUND				
27	S1#	130 C340 1,4-DICHLOROBENZENE	146			1	NOT FOUND				
28	S1#	145 C345 BENZYL ALCOHOL	108			1	NOT FOUND				
29	S1#	150 C350 1,2-DICHLOROBENZENE	146			1	NOT FOUND				
30	S1#	160 C355 2-METHYLPHENOL	108	A	BB	392	1	2412. 1.175	2.692		
31	S1#	165 C360 2,2'-OXYBIS(1-CLPRO	45	A	BB	394	1	3276. 2.539	1.691		
32	S1#	170 C361 ACETOPHENONE	105	A	BB	400	1	8775. 1.934	5.948		
33	S1#	175 N-NITROSOPYRROLIDINE	100			1	NOT FOUND				
34	S1#	180 N-NITROSOMORPHOLINE	56			1	NOT FOUND				
35	S1#	182 3-METHYL PHENOL	108	A	BB	418	1	2022. 2.252	1.177		
36	S1#	185 C365 4-METHYLPHENOL	108	A	BB	418	1	2003. 1.276	2.059		
37	S1#	190 C370 N-NITROSO-DI-N-PROP	70			1	NOT FOUND				
38	S1#	195 O-TOLUIDINE	106			1	NOT FOUND				
39	S1#	200 C375 HEXACHLOROETHANE	117			1	NOT FOUND				
40	S2#	10 C410 NITROBENZENE	77	A	VB	422	2	2364. 0.445	1.930		
41	S2#	15 N-NITROSOPIPERIDINE	42	A	BB	449	2	1186. 0.320	1.345		
42	S2#	20 C415 ISOPHORONE	82			2	NOT FOUND				
43	S2#	25 C420 2-NITROPHENOL	139			2	NOT FOUND				
44	S2#	30 C425 2,4-DIMETHYLPHENOL	107	A	BB	500	2	3534. 0.330	3.893		
45	S2#	35 C435 BIS(2-CHLOROETHOXY)	93			2	NOT FOUND				

46	S2# 40	C440 2,4-DICHLOROPHENOL	162 A BB	519 2	256. 0.308	0.302
47	S2# 45	C430 BENZOIC ACID	122 A BB	548 2	438. 0.276	0.577
48	S2# 50	AA-DIMETHYLPHENETHYLAMIN	58	2 NOT FOUND		
49	S2# 55	C445 1,2,4-TRICHLOROBENZ	180 A*BB	523 2	252. 0.331	0.277

50	S2# 60	C450 NAPHTHALENE	128 A BV	533 2	29645. 1.022	10.538
51	S2# 80	C455 4-CHLOROANALINE	127 A BB	556 2	677. 0.454	0.542
52	S2# 85	2,6-DICHLOROPHENOL	162	2	NOT FOUND	
53	S2# 90	HEXACHLOROPROPENE	213	2	NOT FOUND	
54	S2# 95	C460 HEXACHLOROBUTADIENE	225	2	NOT FOUND	
55	S2#115	P-PHENYLENE DIAMINE	108	2	NOT FOUND	
56	S2#120	N-NITROSODI-N-BUTYLAMINE	84	2	NOT FOUND	
57	S2#130	C465 4-CHLORO-3-METHYLPH	107	2	NOT FOUND	
58	S2#140	SAFROLE	162 A BB	648 2	1458. 0.248	2.140
59	S2#145	C470 2-METHYLNAPHTHALENE	142 A BB	652 2	43357. 0.751	20.970
60	S3# 10	1,2,4,5-TETRACHLOROBENZE	214	3	NOT FOUND	
61	S3# 15	ISOSAFROLE (#1)	162 A BV	697 3	592. 0.044	9.033
62	S3# 20	C510 HEXACHLOROCYCLOPENT	237	3	NOT FOUND	
63	S3# 25	C515 2,4,6-TRICHLOROPHEN	196	3	NOT FOUND	
64	S3# 30	C520 2,4,5-TRICHLOROPHEN	196	3	NOT FOUND	
65	S3# 35	ISOSAFROLE (#2)	104 A BB	746 3	1306. 0.195	4.467
66	S3# 40	C525 2-CHLORONAPHTHALENE	162	3	NOT FOUND	
67	S3# 42	1-CHLORONAPHTHALENE	162	3	NOT FOUND	
68	S3# 45	C530 2-NITROANALINE	65 A BB	764 3	838. 0.502	1.112
69	S3# 50	1,4-NAPHTHOQUINONE	158	3	NOT FOUND	
70	S3# 55	C535 DIMETHYLPHTHALATE	163	3	NOT FOUND	
71	S3# 60	1,3-DINITROBENZENE	168 A BB	807 3	340. 0.197	1.148
72	S3# 65	C540 ACENAPHTHYLENE	152	3	NOT FOUND	
73	S3# 70	C543 2,6-DINITROTOLUENE	165 A BB	817 3	429. 0.324	0.883
74	S3# 75	C545 3-NITROANILINE	138	3	NOT FOUND	
75	S3# 80	C550 ACENAPHTHENE	153 A BV	834 3	1539. 1.159	0.885
76	S3# 85	C555 2,4-DINITROPHENOL	184 A BB	858 3	129. 0.189	0.456
77	S3# 90	C565 DIBENZOFURAN	168 A BB	867 3	1304. 1.669	0.521
78	S3# 95	C560 4-NITROPHENOL	109	3	NOT FOUND	
79	S3#100	PENTACHLOROBENZENE	250	3	NOT FOUND	
80	S3#105	C570 2,4-DINITROTOLUENE	165 A VB	881 3	770. 0.397	1.291
81	S3#110	1-NAPHTHYLAMINE	143 A BV	882 3	547. 0.837	0.435
82	S3#115	2-NAPHTHYLAMINE	143 A BB	902 3	425. 0.966	0.293
83	S3#120	2,3,4,6-TETRACHLOROPHENO	232	3	NOT FOUND	
84	S3#130	C580 DIETHYLPHTHALATE	149	3	NOT FOUND	
85	S3#135	C590 FLUORENE	166 A BB	926 3	4384. 1.298	2.251
86	S3#140	C585 4-CHLOROPHENYL-PHEN	204	3	NOT FOUND	
87	S3#145	5-NITRO-O-TOLUIDINE	152 A BB	942 3	4181. 0.356	7.835
88	S3#150	C595 4-NITROANALINE	138 A VB	949 3	213. 0.307	0.463
89	S4# 10	C610 4,6-DINITRO-2-METHY	198	4	NOT FOUND	
90	S4# 15	C615 N-NITROSODIPHENYLAM	169 A VB	964 4	5183. 0.551	3.543
91	S4# 20	C620 AZOBENZENE	77 A VV	966 4	1221. 1.171	0.392
92	S4# 25	SYM-TRINITROBENZENE	75	4	NOT FOUND	
93	S4# 30	C625 4-BROMOPHENYL-PHENY	248	4	NOT FOUND	
94	S4# 35	PHENACETIN	108	4	NOT FOUND	
95	S4# 37	DIALATE	234	4	NOT FOUND	
96	S4# 40	C630 HEXACHLOROBENZENE	284	4	NOT FOUND	
97	S4# 45	4-AMINOBIIPHENYL	169	4	NOT FOUND	
98	S4# 50	C635 PENTACHLOROPHENOL	266	4	NOT FOUND	
99	S4# 55	PRONAMIDE	173	4	NOT FOUND	
100	S4# 60	PENTACHLORONITROBENZENE	237 A BB	1062 4	250. 0.080	1.171
101	S4# 65	C640 PHENANTHRENE	178 A BB	1091 4	13289. 1.033	4.845
102	S4# 70	C645 ANTHRACENE	178	4	NOT FOUND	
103	S4# 75	2SECBUTYL-4,6-DINITROPHE	211	4	NOT FOUND	
104	S4# 80	C647 CARBAZOLE	167 A BB	1148 4	571. 0.764	0.281
105	S4# 85	C650 DI-N-BUTYLPHTHALATE	149	4	NOT FOUND	
106	S4#100	4-NITROQUINOLINE-1-OXIDE	190	4	NOT FOUND	
107	S4#105	METHAPYRILENE	58	4	NOT FOUND	
108	S4#106	ISODRIN	193	4	NOT FOUND	
109	S4#110	C655 FLUORANTHENE	202	4	NOT FOUND	
110	S4#120	CHLOROBENZILATE	139	4	NOT FOUND	

111 S5# 10 C710 BENZIDINE	184	5	NOT FOUND		
112 S5# 15 C715 PYRENE	202	A BB 1335	5	1633.	1.323 0.728
113 S5# 20 ARAMITE (#1)	185	5	NOT FOUND		
114 S5# 25 ARAMITE (#2)	185	5	NOT FOUND		
115 S5# 30 P-DIMETHYLAMINOAZOBENZEN	120	5	NOT FOUND		
116 S5# 35 3,3'-DIMETHYLBENZIDINE	212	5	NOT FOUND		
117 S5# 37 KEPONE	272	A BV 1460	5	234.	0.082 1.684
118 S5# 40 C720 BUTYLBENZYLPHTHALAT	149	5	NOT FOUND		
119 S5# 45 2-ACETYLAMINOFLUORENE	181	5	NOT FOUND		
120 S5# 50 C730 BENZO(A)ANTHRACENE	228	5	NOT FOUND		
121 S5# 55 C725 3,3'-DICHLOROBENZID	252	5	NOT FOUND		
122 S5# 60 C740 CHRYSENE	228	A BB 1555	5	916.	1.014 0.533
123 S5# 65 C745 BIS(2-ETHYLHEXYL)PH	149	5	NOT FOUND		
124 S5# 85 3-METHYLCHOLANTHRENE	268	5	NOT FOUND		
125 S6# 10 C760 DI-N-OCTYL PHTHALAT	149	6	NOT FOUND		
126 S6# 15 C765 BENZO(B)FLUORANTHEN	252	6	NOT FOUND		
127 S6# 20 7,12-DIMETHYLBENZANTHRAC	256	A BB 1741	6	218.	0.610 0.280
128 S6# 25 C770 BENZO(K)FLUORANTHEN	252	6	NOT FOUND		
129 S6# 30 HEXACHLOROPHENE	196	6	NOT FOUND		
130 S6# 35 C775 BENZO(A)PYRENE	252	6	NOT FOUND		
131 S6# 55 C780 INDENO(1,2,3-CD)PYR	276	6	NOT FOUND		
132 S6# 60 C785 DIBENZ(A,H)ANTHRACE	278	6	NOT FOUND		
133 S6# 65 C790 BENZO(G,H,I)PERYLEN	276	6	NOT FOUND		

QUANTERRA Internal Standard Check

Standard Filename: ST16980831

Analyzed: 08/31/98 12:31

Standard	IS# 1			IS# 2			IS# 3		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	28649.		3.78	100432.		5.90	54347.		9.21
Upper Limit	57297.=+200		4.28	200864.=+200		6.40	108695.=+200		9.71
Lower Limit	14324.=+ 50		3.28	50216.=+ 50		5.40	27174.=+ 50		8.71
Filename									
1 30068107	30514.= 106		3.78	110082.= 109		5.90	60039.= 110		9.22

IS# 1 = C130 1,4-DICHLOROBENZENE-D4

IS# 2 = C140 NAPHTHALENE-D8

IS# 3 = C150 ACENAPHTHENE-D10

* - indicates an I.S. is outside QC limit(s)

Standard	IS# 4			IS# 5			IS# 6		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	98520.		12.08	62289.		17.26	50750.		19.83
Upper Limit	197039.=+200		12.58	124578.=+200		17.76	101500.=+200		20.33
Lower Limit	49260.=+ 50		11.58	31145.=+ 50		16.76	25375.=+ 50		19.33
Filename									
1 30068107	106261.= 107		12.10	67833.= 108		17.26	51078.= 100		19.82

IS# 4 = C160 PHENANTHRENE-D10

IS# 5 = C170 CHRYSENE-D12

IS# 6 = C175 PERYLENE-D12

* - indicates an I.S. is outside QC limit(s)

TARGET COMPOUND COMPARISON

COMPOUND: C450 NAPHTHALENE

128 S2# 60

91-20-3

RAW DATA: 30068107 #533

08/31/98 21:59

SAMPLE: S-MM5-3 1/35A/100M

INST. ID: F16

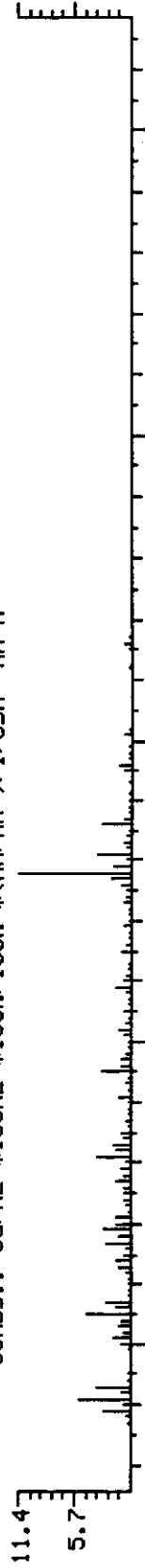
CONDS.: UG/ML *100ML *100%/100% *(NA/NA)/1/35A NA M

RIC: 88448.

BASE M/Z: 128

128

9072.



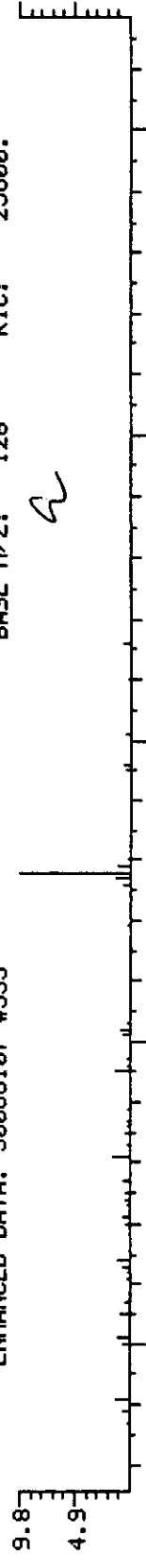
ENHANCED DATA: 30068107 #533

RIC: 25600.

BASE M/Z: 128

128

7768.



STANDARD FILE: ST16980831 #533

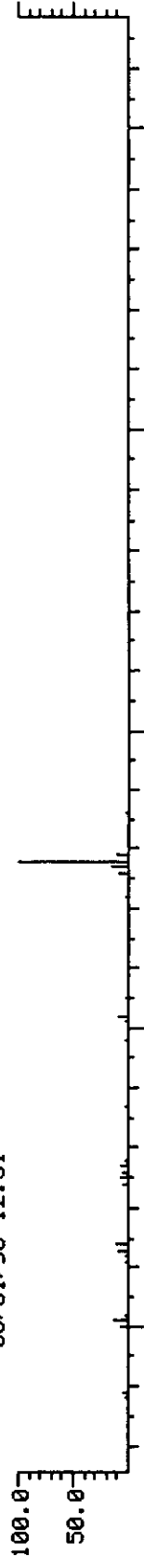
08/31/98 12:31

RIC: 195840.

BASE M/Z: 128

128

79232.



M/Z

100.0

50.0

0.0

-50.0

-100.0

50

100

150

200

250

TARGET COMPOUND COMPARISON

COMPOUND: C470 2-METHYLNAPHTHALENE 142 S2#145 91-57-6

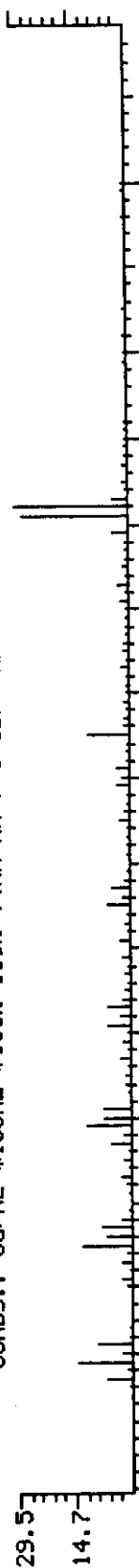
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08/31/98 21:59

SAMPLE: S-MM5-3 1/35A/100M INST. ID: F16

CONDS.: UG/ML *100Z/100Z *KNA/NA >1/35A NA M

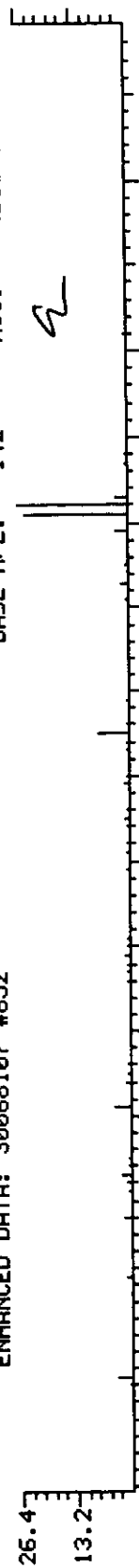
14688.



ENHANCED DATA: 30068107 #652

BASE M/Z: 142 RIC: 49600.

13136.

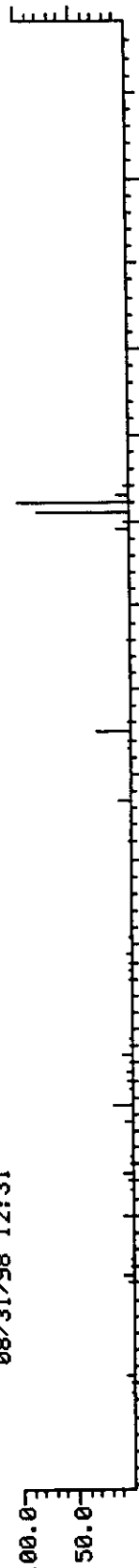


STANDARD FILE: ST16980831 #651

08/31/98 12:31

BASE M/Z: 142 RIC: 198144.

49792.



Data Reduced by: 2 Date: 9/18/98
Data Reviewed by: Date:

Data File: 30068107

QUANTERRA GC/MS TIC REPORT (Part 1)

Sample: S-MM5-3 1/3SA/100M INST. ID: F16
Analyst: DAT Date Analyzed: 08/31/98 21:59
Run Factor: 300.

Concentration in Sample			
# SCAN	(UG/A)	CAS #	
1 444	10000.	1120-21-4	
UNDECANE or isomer			
2 556	12000.	62108-23-0	
DECANE, 2,5,6-TRIMETHYL- or isomer			
3 569	9900.	17301-23-4	
UNDECANE, 2,6-DIMETHYL- or isomer			
4 616	11000.	00-00-0	5700
UNKNOWN			
X 621	7500.	2809-64-5	
NAPHTHALENE, 1,2,3,4-TETRAHYDRO-5-METHYL- or isomer			
6 632	19000.	54105-67-8	
HEPTADECANE, 2,6-DIMETHYL- or isomer			
7 653	13000.	69225-59-8	
1,3-DIOXASPIRO[5.5]UNDECAN-9-ONE, 3,3-DIMETHYL-			unknown 5700
8 663	19000.	2801-84-5	
DECANE, 2,4-DIMETHYL- or isomer			
X 678	8100.	00-00-0	5700
UNKNOWN			
10 700	9100.	00-00-0	
UNKNOWN			

11 738 22000. 3891-98-3 ~~alkane~~ 5746400
DODECANE, 2,6,10-TRIMETHYL-

12 764 36000. 629-59-4
TETRADECANE

13 776 16000. 575-41-7
NAPHTHALENE, 1,3-DIMETHYL- or isomer

14 820 26000. 18344-37-1
HEPTADECANE, 2,6,10,14-TETRAMETHYL- or isomer

15 859 35000. 17312-57-1
DODECANE, 3-METHYL- or isomer

16 904 11000. 52783-43-4 2911858
~~NONADECANOL~~ alkane

17 948 35000. 629-92-5
NONADECANE or isomer

18 988 22000. 629-92-5
NONADECANE

19 1036 46000. 54105-67-8
HEPTADECANE, 2,6-DIMETHYL

~~20~~ 1054 7400. 34303-81-6
~~3-HEXADECENE, (Z)-~~ alkene 5720000

21 1114 14000. 29812-79-1
~~HYDROXYLAMINE, O-DECYL-~~ alkane

22 1119 17000. 629-92-5
NONADECANE or isomer

~~23~~ 1191 9000. 629-92-5
NONADECANE or isomer

24 1335 10000. 54105-67-8
HEPTADECANE, 2,6-DIMETHYL-

~~25~~ 1402 8000. 629-92-5
NONADECANE

DATA FILE: 30068107

QUANTERRA GC/MS TIC REPORT (Part 2)

CONCENTRATION = AREA(TIC)*CONC(IS)/AREA(IS)

#	FIT	PURITY	INT.		RRT	AREA	HEIGHT	AMOUNT		LIB	LIB #
			STD.	RT				AS ANALYZED			
1	990	746	2	4:56	0.536	283914.	85573.	33.684	NB	11607.	
2	945	884	2	6:11	0.671	328256.	110848.	38.945	NB	19015	
3	979	888	2	6:19	0.686	277376.	87296.	32.908	NB	19054.	
4	819	392	2	6:51	0.743	312896.	46976.	37.122	UK	1.	
5	898	644	2	6:54	0.749	210688.	47232.	24.996	NB	8945.	
6	945	794	2	7:01	0.762	538176.	133888.	63.850	NB	37462.	
7	869	653	2	7:15	0.788	364315.	73034.	43.223	NB	22336.	
8	966	848	2	7:22	0.800	544896.	138240.	64.647	NB	15356.	
9	884	466	2	7:32	0.818	227072.	58419.	26.940	UK	1.	
10	824	399	2	7:47	0.844	256768.	30080.	30.463	UK	1.	
11	975	822	2	8:12	0.890	620928.	165376.	73.667	NB	25991	
12	951	723	2	8:29	0.922	1009280.	270848.	119.742	NB	22530.	
13	958	638	2	8:37	0.936	451456.	64768.	53.561	NB	11625.	
14	980	837	2	9:07	0.989	738048.	155392.	87.563	NB	42196.	
15	960	803	2	9:33	1.036	993536.	249088.	117.874	NB	19007.	
16	923	647	2	10:03	1.090	323072.	72192.	38.330	NB	40233.	
17	948	843	2	10:32	1.144	969600.	241408.	115.034	NB	37465.	
18	945	832	3	10:59	0.908	712448.	182016.	72.262	NB	37465.	
19	976	875	3	11:31	0.952	1498370.	258304.	151.977	NB	37462.	
20	914	722	3	11:43	0.969	243136.	57504.	24.661	NB	28768.	
21	962	804	3	12:23	1.024	445568.	136192.	45.193	NB	15969.	
22	964	839	3	12:26	1.028	572544.	170240.	58.072	NB	37465.	
23	938	769	3	13:14	1.095	294144.	98688.	29.834	NB	37465.	
24	890	720	4	14:50	0.860	194624.	47580.	33.353	NB	37462.	
25	953	662	4	15:35	0.903	154880.	34688.	26.542	NB	37465.	

Library Search Data: 30068107 # 444 Base m/z: 43
 08/31/98 21:59:00 + 4:56 Cali: 30068107 # 3 RIC: 68608.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 736 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 11607 UNDECANE
 2 19015 DECANE, 2,5,6-TRIMETHYL-
 3 8104 OCTANE, 3,5-DIMETHYL-
 4 11602 OCTANE, 2,4,6-TRIMETHYL-
 5 19016 UNDECANE, 4,7-DIMETHYL-
 6 22530 TETRADECANE
 7 14793 1-UNDECENE, 4-METHYL-
 8 19013 DECANE, 2,5,9-TRIMETHYL-
 9 5151 HEPTANE, 3,4-DIMETHYL-

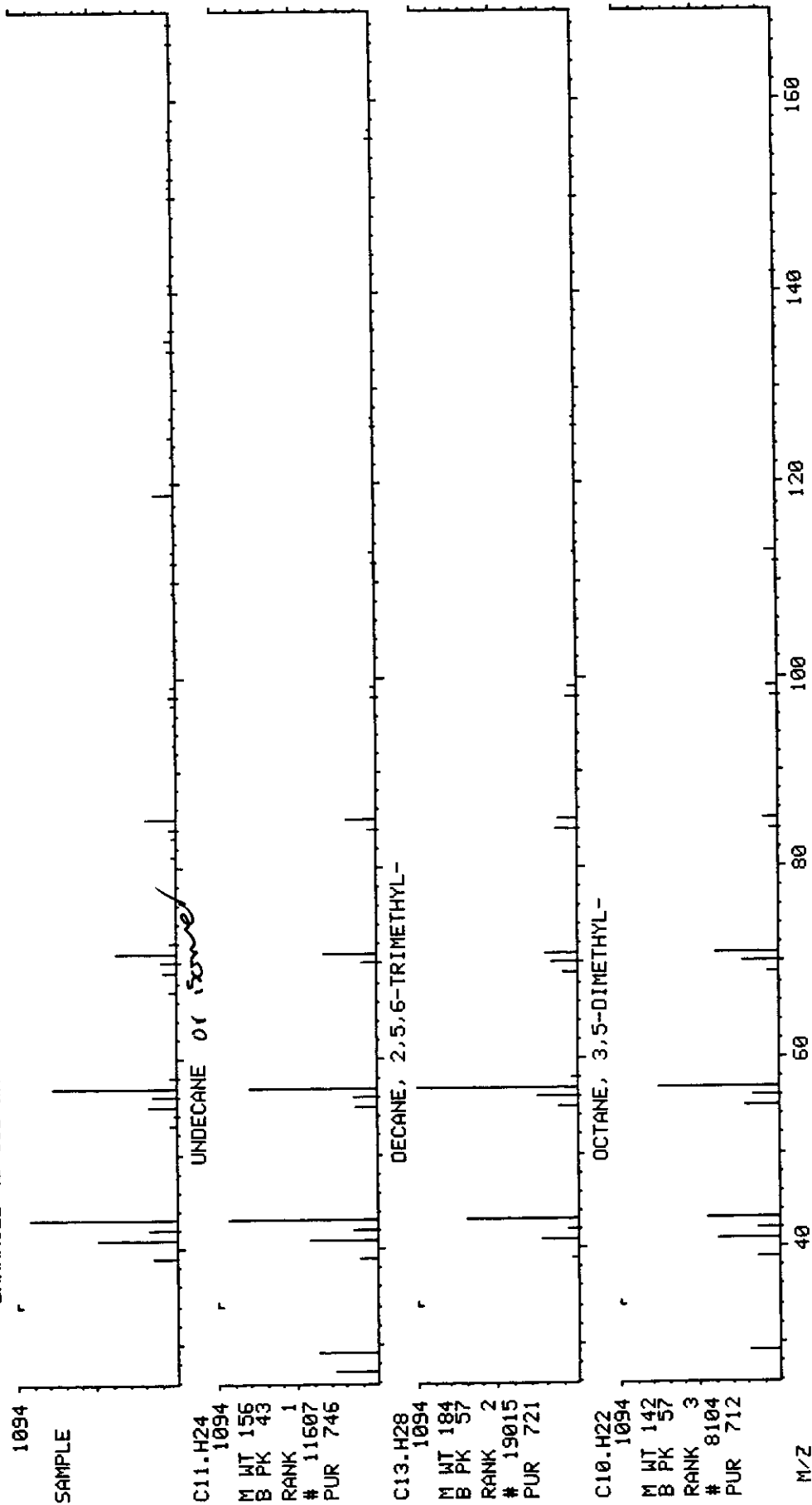
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C11.H24	156	43	746	990	749
2	C13.H28	184	57	721	933	723
3	C10.H22	142	57	712	938	718
4	C11.H24	156	57	708	958	708
5	C13.H28	184	43	704	942	710
6	C14.H30	198	43	702	913	755
7	C12.H24	168	43	699	921	735
8	C13.H28	184	57	697	924	697
9	C9.H20	128	43	695	915	697

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	1120-21-4
2	—	—	—	—	62108-23-0
3	—	—	—	—	15869-93-9
4	—	—	—	—	62016-37-9
5	—	—	—	—	17301-32-5
6	—	—	—	—	629-59-4
7	—	—	—	—	74630-39-0
8	—	—	—	—	62108-22-9
9	—	—	—	—	922-28-1

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:59:00 + 4:56
 SAMPLE: S-MMS-3 1/35A/100M
 CONDS.: UG/ML *100Z/100Z *(NA/NA) /1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068107 # 444
 CALI: 30068107 # 3

BASE M/Z: 43
 RIC: 68608.



Library Search Data: 30068107 # 556 Base m/z: 43
 08/31/98 21:59:00 + 6:11 Cali: 30068107 # 3 RIC: 96896.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 206 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 19015 DECANE, 2,5,6-TRIMETHYL-
 2 11602 OCTANE, 2,4,6-TRIMETHYL-
 3 19016 UNDECANE, 4,7-DIMETHYL-
 4 8104 OCTANE, 3,5-DIMETHYL-
 5 8089 NONANE, 2-METHYL-
 6 5151 HEPTANE, 3,4-DIMETHYL-
 7 19013 DECANE, 2,5,9-TRIMETHYL-
 8 15353 2,6-DIMETHYLDECANE
 9 8074 OCTANE, 2,7-DIMETHYL-

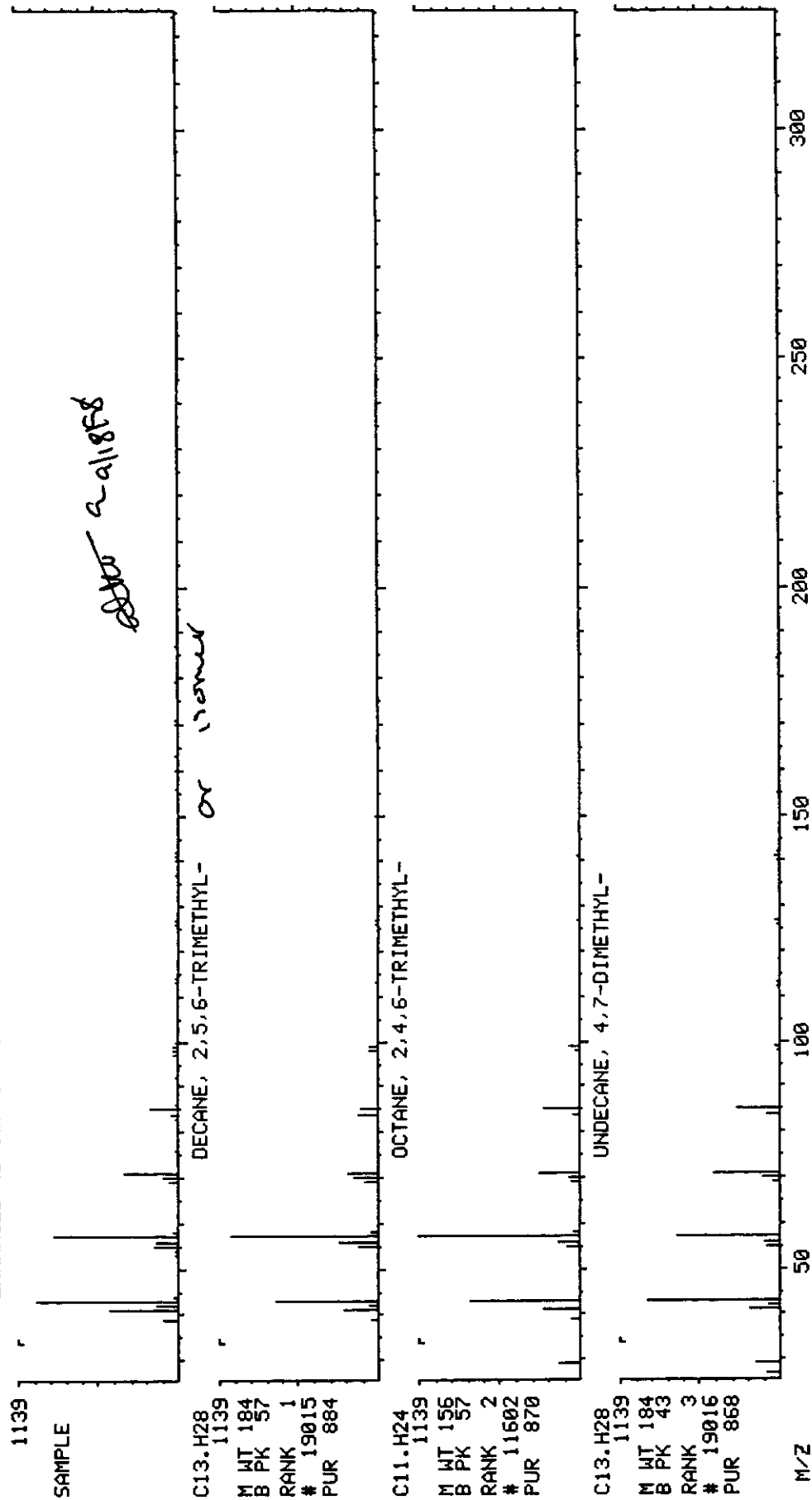
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H28	184	57	884	945	885
2	C11.H24	156	57	870	971	870
3	C13.H28	184	43	868	952	874
4	C10.H22	142	57	858	934	866
5	C10.H22	142	43	853	951	856
6	C9.H20	128	43	845	923	848
7	C13.H28	184	57	840	926	850
8	C12.H26	170	43	837	893	878
9	C10.H22	142	43	836	923	836

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	62108-23-0
2	—	—	—	—	62016-37-9
3	—	—	—	—	17301-32-5
4	—	—	—	—	15869-93-9
5	—	—	—	—	871-83-0
6	—	—	—	—	922-28-1
7	—	—	—	—	62108-22-9
8	—	—	—	—	13150-81-7
9	—	—	—	—	1072-16-8

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:59:00 + 6:11
 SAMPLE: 5-MMS-3 1/35A/100M
 COND5.: UG/ML *100ML *100Z/100Z *(NA/NA) 1/35A NA M
 ENHANCED (5 15B 2N 0T)

DATA: 30068107 # 556
 CALI: 30068107 # 3

BASE M/Z: 43
 RIC: 96896.



Library Search Data: 30068107 # 569 Base m/z: 57
 08/31/98 21:59:00 + 6:20 Cali: 30068107 # 3 RIC: 76800.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N OT)

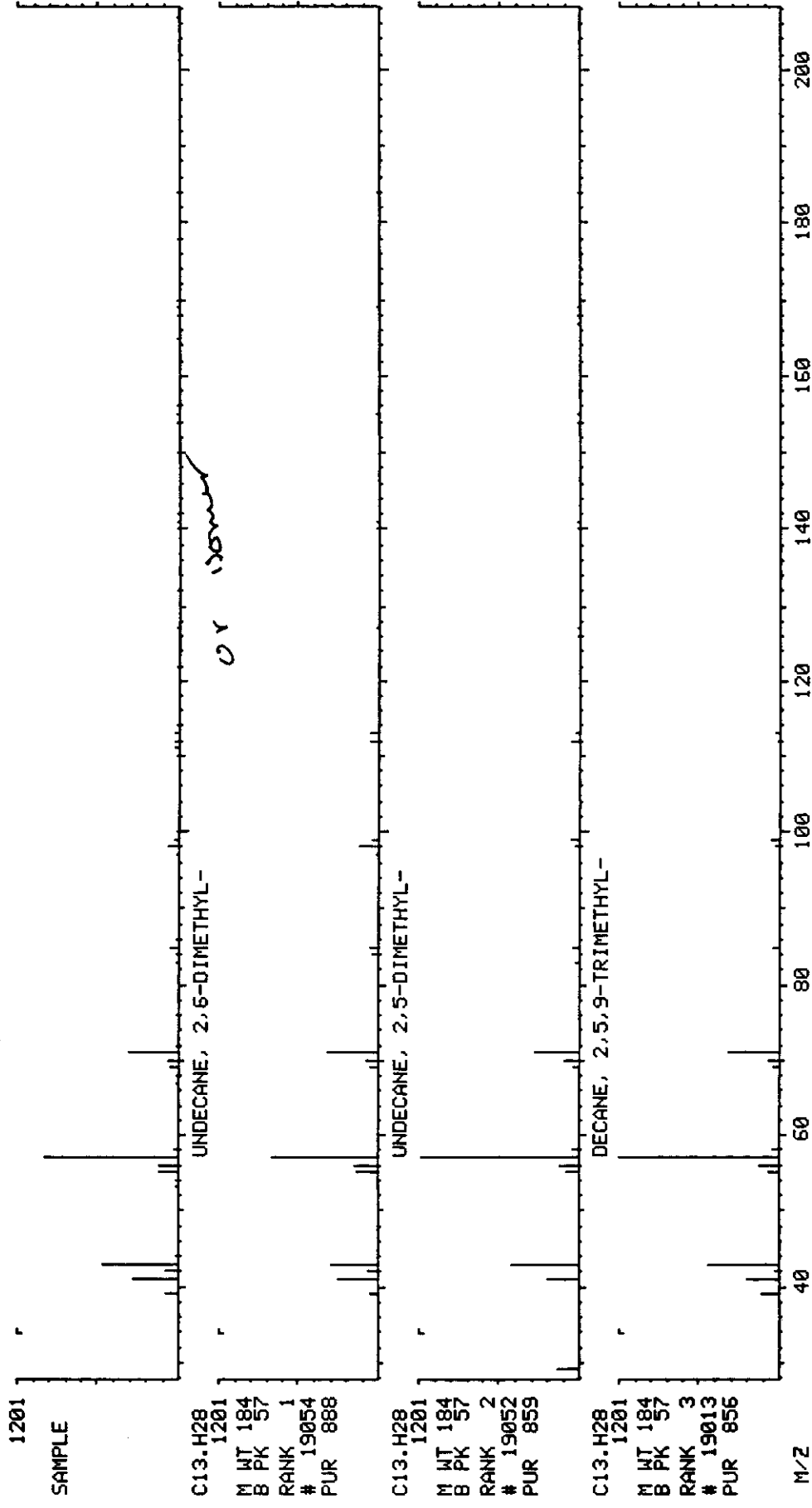
62231 spectra in LIBRARYNB searched for maximum PURITY
 860 matched at least 7 of the 16 largest peaks in the unknown

Rank	In.	Name
1	19054	UNDECANE, 2,6-DIMETHYL-
2	19052	UNDECANE, 2,5-DIMETHYL-
3	19013	DECANE, 2,5,9-TRIMETHYL-
4	19002	DODECANE, 6-METHYL-
5	19026	DECANE, 2,6,8-TRIMETHYL-
6	8104	OCTANE, 3,5-DIMETHYL-
7	18996	UNDECANE, 3,6-DIMETHYL-
8	19004	UNDECANE, 4,6-DIMETHYL-
9	8099	OCTANE, 2,6-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H28	184	57	888	979	888
2	C13.H28	184	57	859	975	862
3	C13.H28	184	57	856	975	856
4	C13.H28	184	57	854	972	857
5	C13.H28	184	57	850	968	855
6	C10.H22	142	57	842	960	842
7	C13.H28	184	57	831	967	834
8	C13.H28	184	57	824	959	824
9	C10.H22	142	57	823	946	829

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	17301-23-4
2	—	—	—	—	17301-22-3
3	—	—	—	—	62108-22-9
4	—	—	—	—	6044-71-9
5	—	—	—	—	62108-26-3
6	—	—	—	—	15869-93-9
7	—	—	—	—	17301-28-9
8	—	—	—	—	17312-82-2
9	—	—	—	—	2051-30-1

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:59:00 + 6:20
 SAMPLE: 5-MM5-3 1/35A/100M
 CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068107 # 569
 CALI: 30068107 # 3
 BASE M/Z: 57
 RIC: 76800.



Library Search Data: 30068107 # 616 Base m/z: 131
 08/31/98 21:59:00 + 6:51 Cali: 30068107 # 3 RIC: 35968.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 789 matched at least 5 of the 16 largest peaks in the unknown

Rank In. Name
 1 8953 7-ETHYLIDENE BICYCLO[4.2.1]NONA-2,4-DIENE
 2 24994 1,1'-BICYCLOHEXYL, 2-(1-METHYLETHYL)-, TRANS-
 3 8915 PROPANEDINITRILE, CYCLOHEXYLIDENE-
 4 24993 1,1'-BICYCLOHEXYL, 2-(1-METHYLETHYL)-, CIS-
 5 47544 BENZENE, 1,1'-[3-(3-CYCLOPENTYLPROPYL)-1,5-PENTANEDIYL]BIS-
 6 8902 7,7-NORCARANEDICARBONITRILE, CIS-
 7 33979 BORINIC ACID, DIETHYL-, 1-CYCLODODECEN-1-YL ESTER
 8 24997 1,1'-BICYCLOHEXYL, 2-PROPYL-, TRANS-
 9 16369 BENZENE, (3,3-DIMETHYL-4-PENTENYL)-

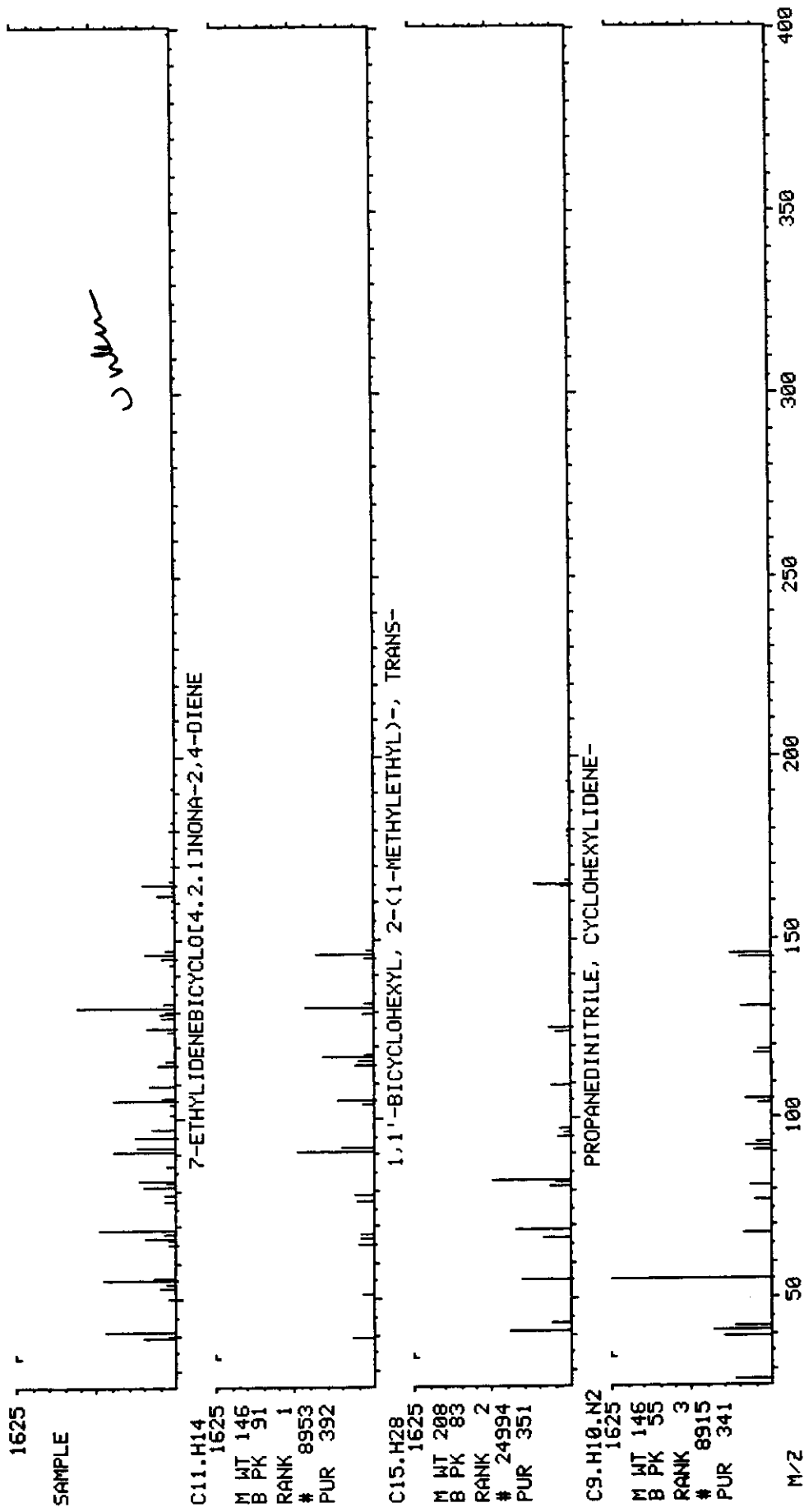
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C11.H14	146	91	392	819	456
2	C15.H28	208	83	351	876	365
3	C9.H10.N2	146	55	341	733	395
4	C15.H28	208	83	339	863	356
5	C25.H34	334	92	339	584	364
6	C9.H10.N2	146	55	335	722	388
7	C16.H31.O.B	250	83	327	845	365
8	C15.H28	208	41	313	765	332
9	C13.H18	174	104	312	647	334

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	94400-10-9
2	—	—	—	—	50991-16-7
3	—	—	—	—	4354-73-8
4	—	—	—	—	50991-15-6
5	—	—	—	—	55191-62-3
6	—	—	—	—	29782-28-3
7	—	—	—	—	61142-73-2
8	—	—	—	—	54934-89-3
9	—	—	—	—	61142-18-5

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:59:00 + 6:51
 SAMPLE: S-MMS-3 1/35A/100M
 COND.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068107 # 616
 CALI: 30068107 # 3

BASE M/Z: 131
 RIC: 35968.



1
 12
 13
 14

Library Search Data: 30068107 # 621 Base m/z: 118
 08/31/98 21:59:00 + 6:54 Cal: 30068107 # 3 RIC: 31552.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

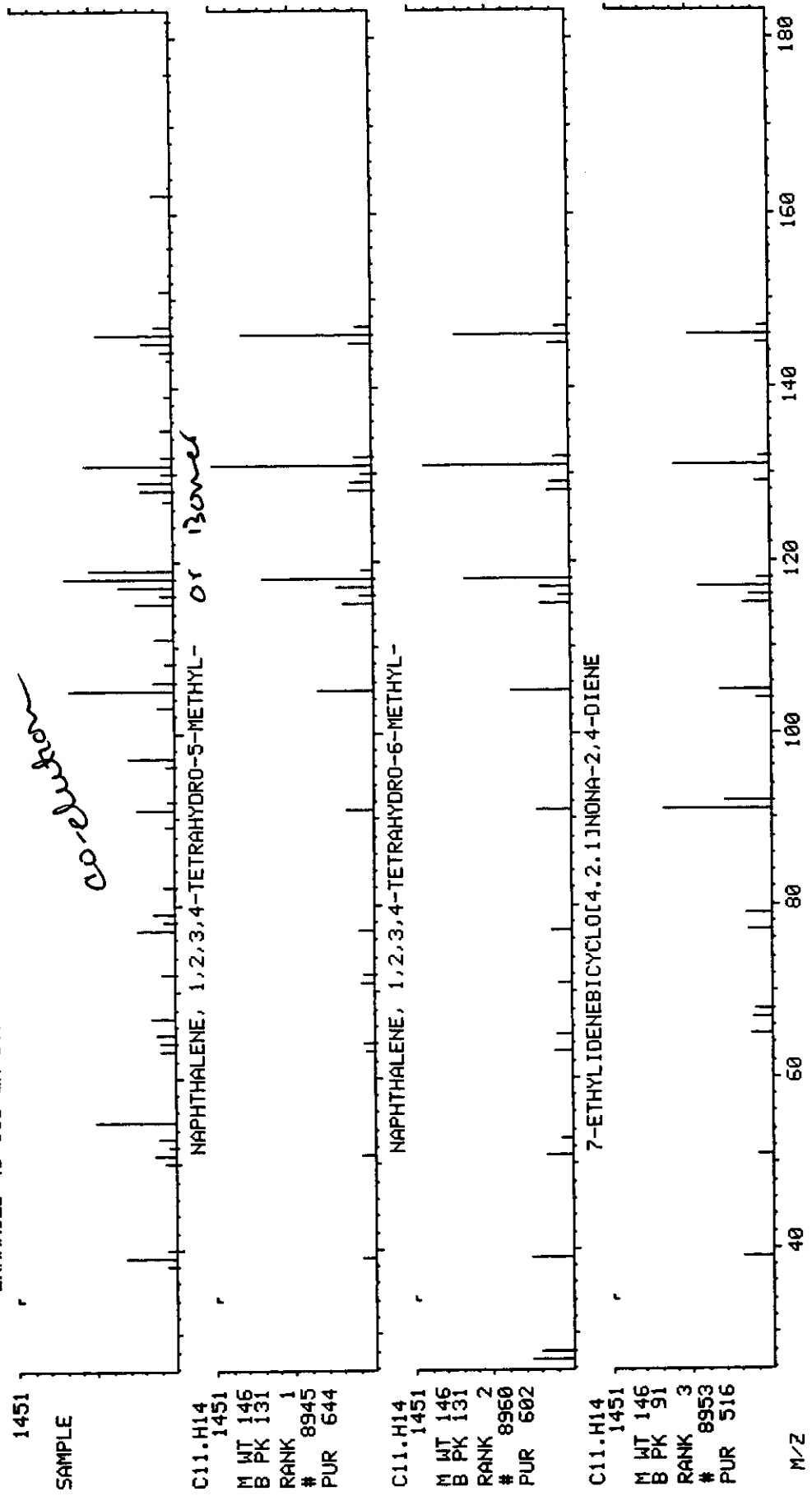
62231 spectra in LIBRARYNB searched for maximum PURITY
 740 matched at least 5 of the 16 largest peaks in the unknown

Rank In. Name
 1 8945 NAPHTHALENE, 1,2,3,4-TETRAHYDRO-5-METHYL-
 2 8960 NAPHTHALENE, 1,2,3,4-TETRAHYDRO-6-METHYL-
 3 8953 7-ETHYLIDENEBICYCLO[4.2.1]NONA-2,4-DIENE
 4 8929 NAPHTH[1,2-B]OXIRENE, 1A,2,3,7B-TETRAHYDRO-
 5 8938 NAPHTHALENE, 1,2,3,4-TETRAHYDRO-1-METHYL-
 6 8944 BENZENE, (1-METHYLENEBUTYL)-
 7 8932 1(2H)-NAPHTHALENONE, 3,4-DIHYDRO-
 8 8973 BENZENE, 2-ETHENYL-1,3,5-TRIMETHYL-
 9 32808 BENZENE, (1-PROPYL-1-NONENYL)-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C11.H14 ✓	146	131	644	898	650
2	C11.H14 ✓	146	131	602	947	609
3	C11.H14 ✓	146	91	516	833	535
4	C10.H10.O	146	146	504	799	580
5	C11.H14	146	131	502	861	516
6	C11.H14	146	118	499	846	520
7	C10.H10.O	146	118	489	767	547
8	C11.H14	146	131	488	871	492
9	C18.H28	244	118	484	696	605

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	2809-64-5
2	—	—	—	—	1680-51-9
3	—	—	—	—	94400-10-9
4	—	—	—	—	2461-34-9
5	—	—	—	—	1559-81-5
6	—	—	—	—	5676-32-4
7	—	—	—	—	529-34-0
8	—	—	—	—	769-25-5
9	—	—	—	—	54986-39-9

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:59:00 + 6:54
 SAMPLE: 5-MM5-3 1/35A/100M
 COND5.: UG/ML *100ML *100Z/100Z *(NA/NA >1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30058107 # 621
 CALI: 30058107 # 3
 BASE M/Z: 118
 RIC: 31552.



Library Search Data: 30068107 # 632 Base m/z: 57
 08/31/98 21:59:00 + 7:02 Cali: 30068107 # 3 RIC: 116992.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

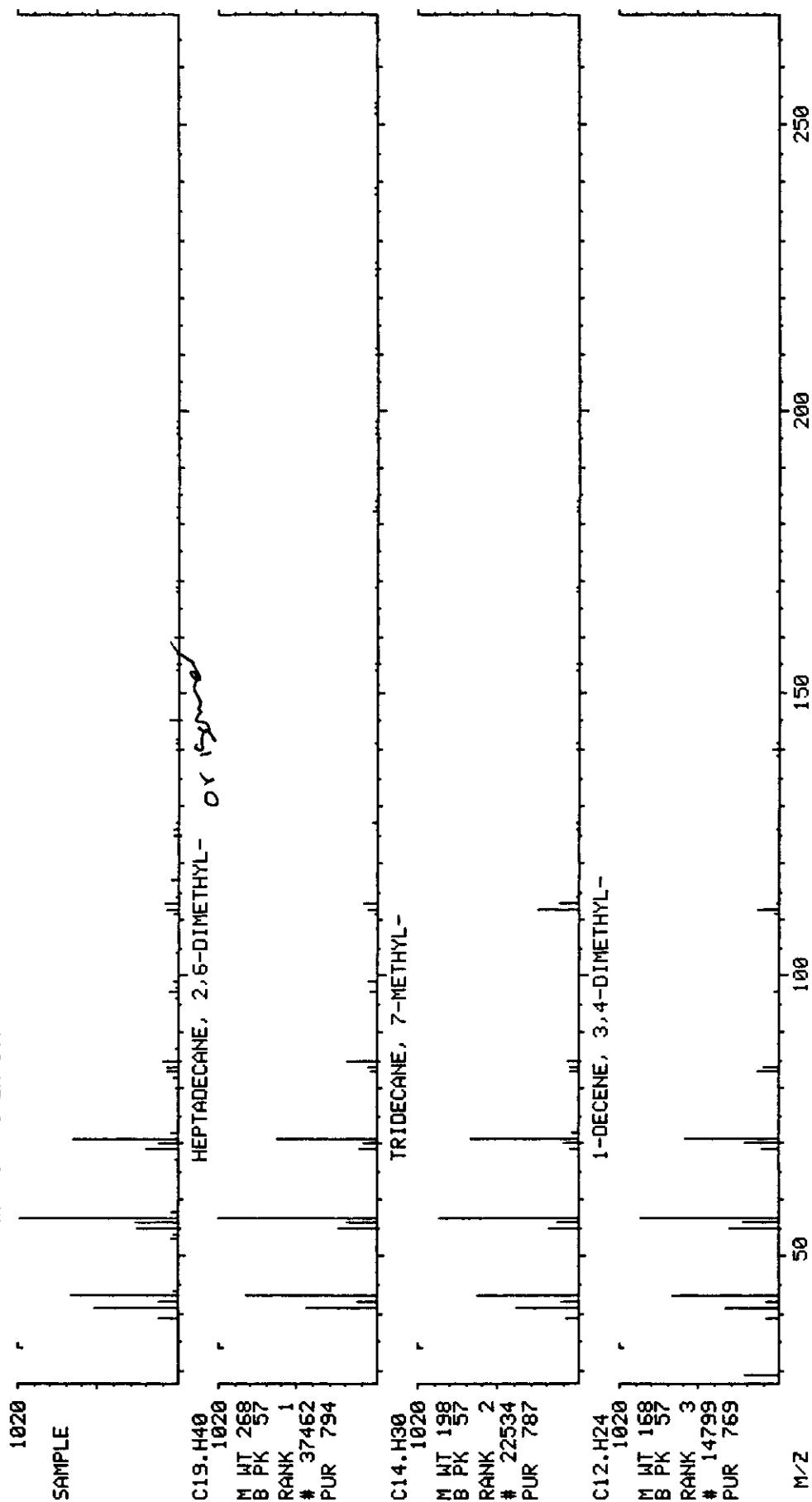
62231 spectra in LIBRARYNB searched for maximum PURITY
 323 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 22534 TRIDECANE, 7-METHYL-
 3 14799 1-DECENE, 3,4-DIMETHYL-
 4 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 5 22535 DODECANE, 4,6-DIMETHYL-
 6 15353 2,6-DIMETHYLDECANE
 7 19023 DECANE, 2,6,7-TRIMETHYL-
 8 19523 1-DECANOL, 2-ETHYL-
 9 25994 DODECANE, 2,6,11-TRIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	794	945	819
2	C14.H30	198	57	787	937	790
3	C12.H24	168	57	769	948	770
4	C21.H44	296	57	766	928	787
5	C14.H30	198	57	763	945	765
6	C12.H26	170	43	762	927	762
7	C13.H28	184	57	762	969	762
8	C12.H26.O	186	57	760	935	764
9	C15.H32	212	57	758	956	760

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	54105-67-8
2	---	---	---	---	26730-14-3
3	---	---	---	---	50871-03-9
4	---	---	---	---	18344-37-1
5	---	---	---	---	61141-72-8
6	---	---	---	---	13150-81-7
7	---	---	---	---	62108-25-2
8	---	---	---	---	21078-65-9
9	---	---	---	---	31295-56-4

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:59:00 + 7:02
 SAMPLE: S-MM5-3 1/35A/100M
 COND.S.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068107 # 632
 CALI: 30068107 # 3
 BASE M/Z: 57
 RIC: 116992.



Library Search Data: 30068107 # 653 Base m/z: 142
 08/31/98 21:59:00 + 7:16 Cali: 30068107 # 3 RIC: 73600.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARY8 searched for maximum PURITY
 439 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 22336 1,5-DIOXASPIRO[5.5]UNDECAN-9-ONE, 3,3-DIMETHYL-
 2 7849 3(2H)-FURANONE, 4-METHOXY-2,5-DIMETHYL-
 3 15148 1,3,2-DIOXABORINANE, 2,4-DIETHYL-5,6-DIMETHYL-
 4 27506 DECANE, 3-BROMO-
 5 22434 CYCLOHEXANECARBOXYLIC ACID, 1-BUTYL-, METHYL ESTER
 6 11522 HEPTANE, 3-[(ETHENYLOXY)METHYL]-
 7 32423 HEPTANE, 3,3'-[OXYBIS(METHYLENE)]BIS-
 8 14759 CYCLOPROPANE, 1-(2-METHYLBUTYL)-1-(1-METHYLPROPYL)-
 9 42522 DECANE, 1,1'-OXYBIS-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C11.H18.O3	198	141	653	869	657
2	C7.H10.O3	142	43	510	706	544
3	C9.H19.O2.B	170	56	449	623	528
4	C10.H21.BR	220	57	434	600	444
5	C12.H22.O2	198	83	406	634	494
6	C10.H20.O	156	57	400	556	421
7	C16.H34.O	242	57	387	537	407
8	C12.H24	168	57	364	491	394
9	C20.H42.O	298	43	358	643	406

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	69225-59-8
2	—	—	—	—	4077-47-8
3	—	—	—	—	74744-56-2
4	—	—	—	—	30571-71-2
5	—	—	—	—	7362-81-4
6	—	—	—	—	103-44-6
7	—	—	—	—	10143-60-9
8	—	—	—	—	64723-36-0
9	—	—	—	—	2456-28-2

BASE M/Z: 142
RIC: 73600.

MID LIBRARY SEARCH (LIBRARYNB)

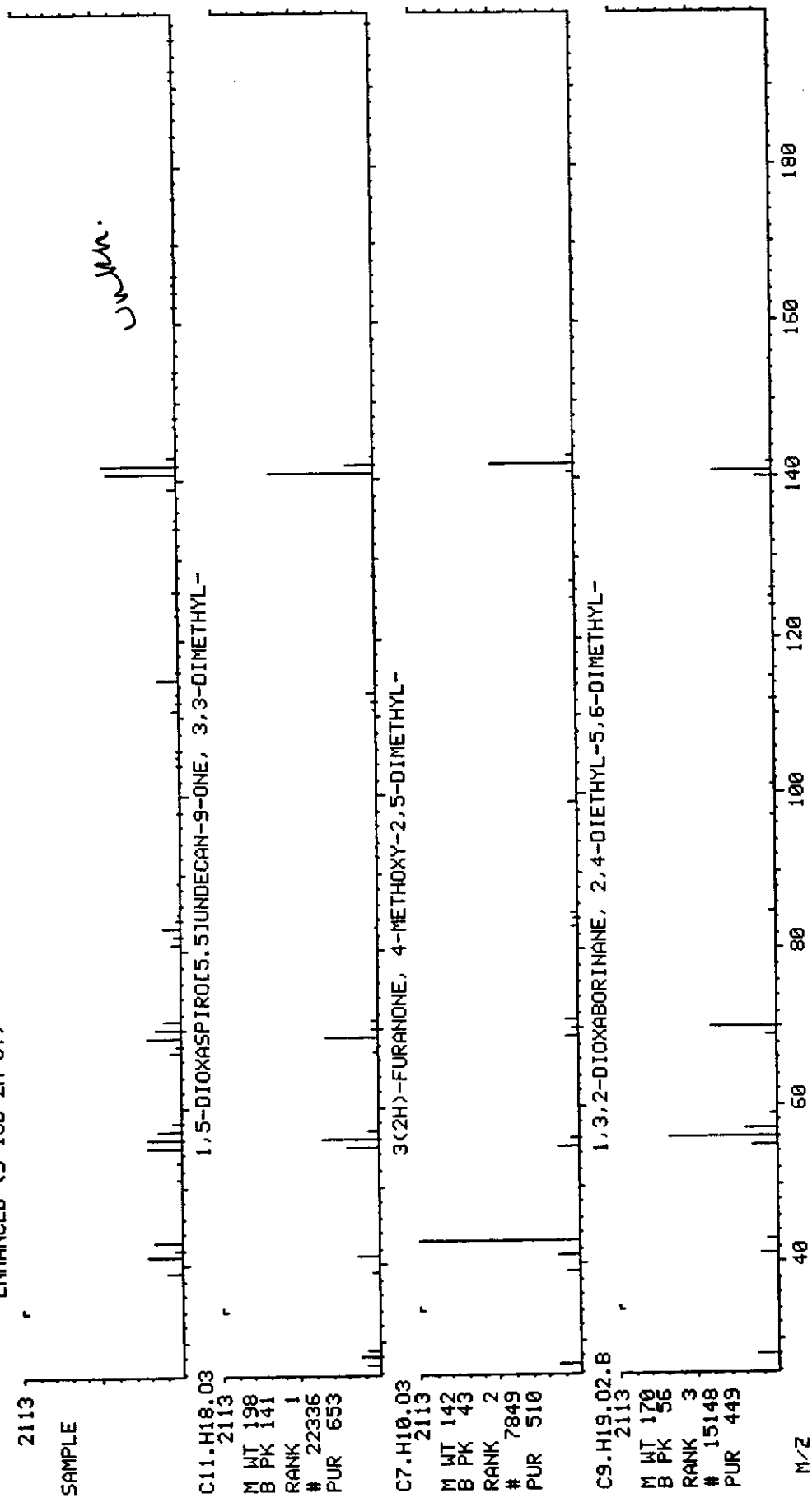
08/31/98 21:59:00 + 7:16

SAMPLE: S-MM5-3 1/35A/100M

INST. ID: F16

COND5.: UG/ML *100ML *100%/100% *(NA/NA)/1/35A NA M

ENHANCED (5 15B 2N 0T) (11 N 5B 2N 0T)



1141

Library Search Data: 30068107 # 663 Base m/z: 57
 08/31/98 21:59:00 + 7:22 Cali: 30068107 # 3 RIC: 119168.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

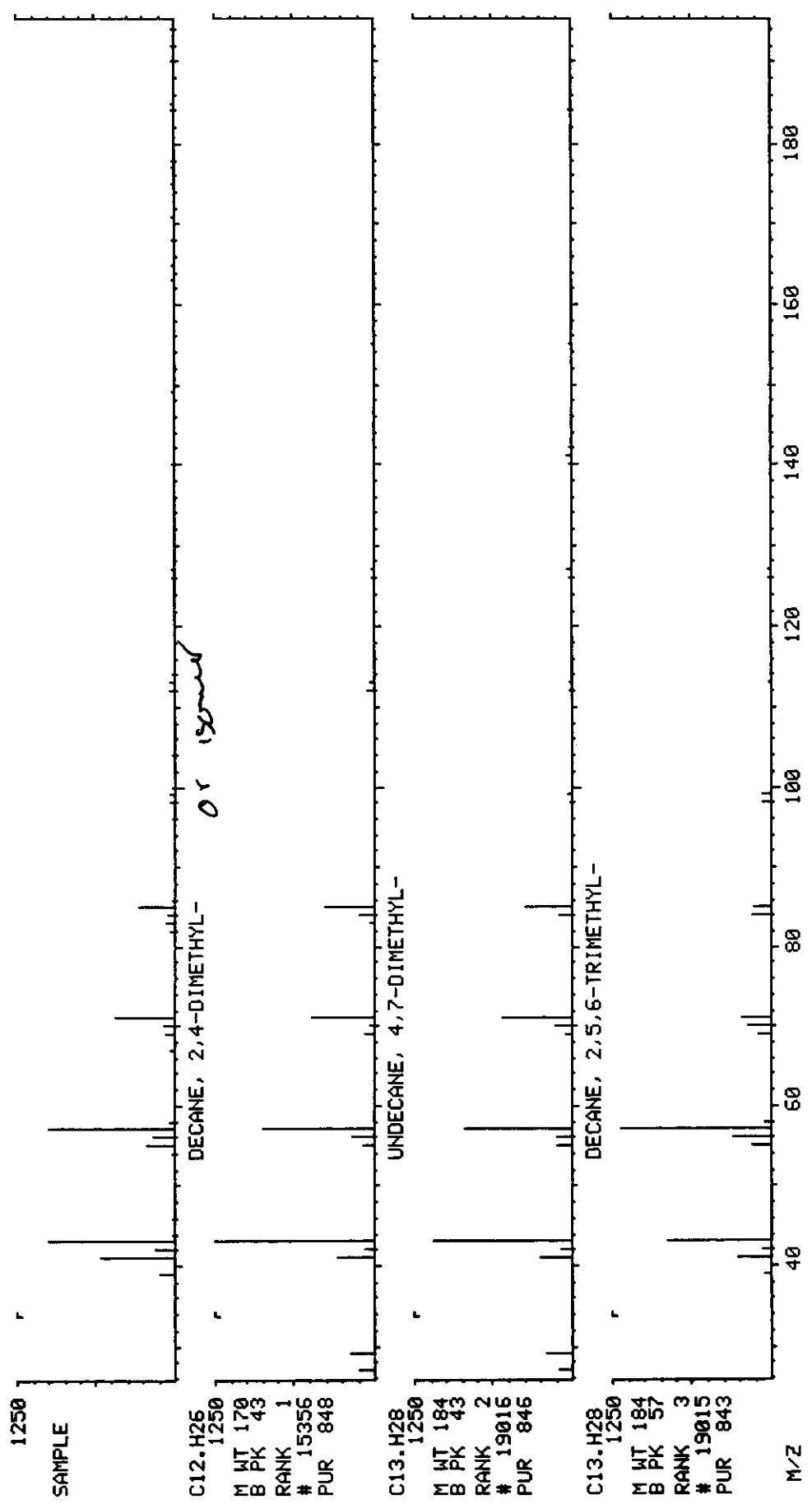
62231 spectra in LIBRARYNB searched for maximum PURITY
 372 matched at least 8 of the 16 largest peaks in the unknown

Rank	In.	Name
1	15356	DECANE, 2,4-DIMETHYL-
2	19016	UNDECANE, 4,7-DIMETHYL-
3	19015	DECANE, 2,5,6-TRIMETHYL-
4	11602	OCTANE, 2,4,6-TRIMETHYL-
5	8104	OCTANE, 3,5-DIMETHYL-
6	15359	UNDECANE, 5-METHYL-
7	18985	TRIDECANE
8	19013	DECANE, 2,5,9-TRIMETHYL-
9	19026	DECANE, 2,6,8-TRIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H26	170	43	848	966	849
2	C13.H28	184	43	846	941	861
3	C13.H28	184	57	843	937	843
4	C11.H24	156	57	836	971	837
5	C10.H22	142	57	836	938	844
6	C12.H26	170	43	835	953	840
7	C13.H28	184	57	831	950	846
8	C13.H28	184	57	826	935	826
9	C13.H28	184	57	824	930	841

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	2801-84-5
2	—	—	—	—	17301-32-5
3	—	—	—	—	62108-23-0
4	—	—	—	—	62016-37-9
5	—	—	—	—	15869-93-9
6	—	—	—	—	1632-70-8
7	—	—	—	—	629-50-5
8	—	—	—	—	62108-22-9
9	—	—	—	—	62108-26-3

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:59:00 + 7:22
 SAMPLE: S-MM5-3 1/35A/100M
 CONDS.: UG/ML #100ML #100% /100% *(NA/NA) /1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30068107 # 653
 CALI: 30068107 # 3
 BASE M/Z: 57
 RIC: 119168.



Library Search Data: 30068107 # 678 Base m/z: 57
 08/31/98 21:59:00 + 7:32 Cali: 30068107 # 3 RIC: 56960.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

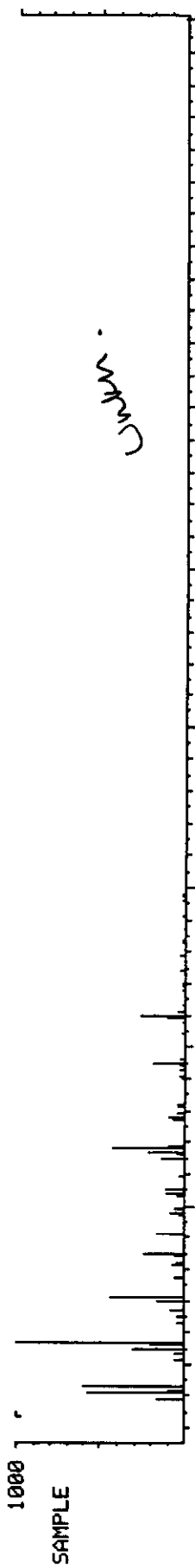
62231 spectra in LIBRARYNB searched for maximum PURITY
 427 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 12534 NAPHTHALENE, 1,2,3,4-TETRAHYDRO-2,6-DIMETHYL-
 2 12552 NAPHTHALENE, 1,2,3,4-TETRAHYDRO-2,7-DIMETHYL-
 3 32792 FALCARINOL (2)-(-)-1,9-HEPTADECADIENE-4,6-DIYNE-3-OL
 4 12526 1,4-DIMETHYL-1,2,3,4-TETRAHYDRONAPHTHALENE
 5 30843 4,4,6-TRIMETHYL-6-PHENYLTETRAHYDRO-1,3-OXAZINE-2-THIONE
 6 5159 NONANE
 7 30039 BENZENE, (4-METHYL-1-DECENYL)-
 8 12576 1H-INDENE, 1-ETHYL-2,3-DIHYDRO-1-METHYL-
 9 37251 1-iodo-2-methylnonane

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H16	160	118	466	884	468
2	C12.H16	160	118	447	959	450
3	C17.H24.O	244	55	414	733	449
4	C12.H16	160	118	400	914	418
5	C13.H17.O.N.S	235	118	398	744	492
6	C9.H20	128	43	394	945	394
7	C17.H26	230	104	389	615	444
8	C12.H16	160	131	387	669	387
9	C10.H21.I	268	43	385	923	401

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	7524-63-2
2	---	---	---	---	13065-07-1
3	---	---	---	---	- -
4	---	---	---	---	- -
5	---	---	---	---	- -
6	---	---	---	---	111-84-2
7	---	---	---	---	62337-88-6
8	---	---	---	---	56298-75-0
9	---	---	---	---	- -

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:59:00 + 7:32
 SAMPLE: S-MM5-3 1/35A/100M
 COND.S.: UG/ML *100ML *100Z/100Z *(NA/NA) /1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068107 # 678
 CALI: 30068107 # 3
 BASE M/Z: 57
 RIC: 56950.



Library Search Data: 30068107 # 700 Base m/z: 81
 08/31/98 21:59:00 + 7:47 Cali: 30068107 # 3 RIC: 28224.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

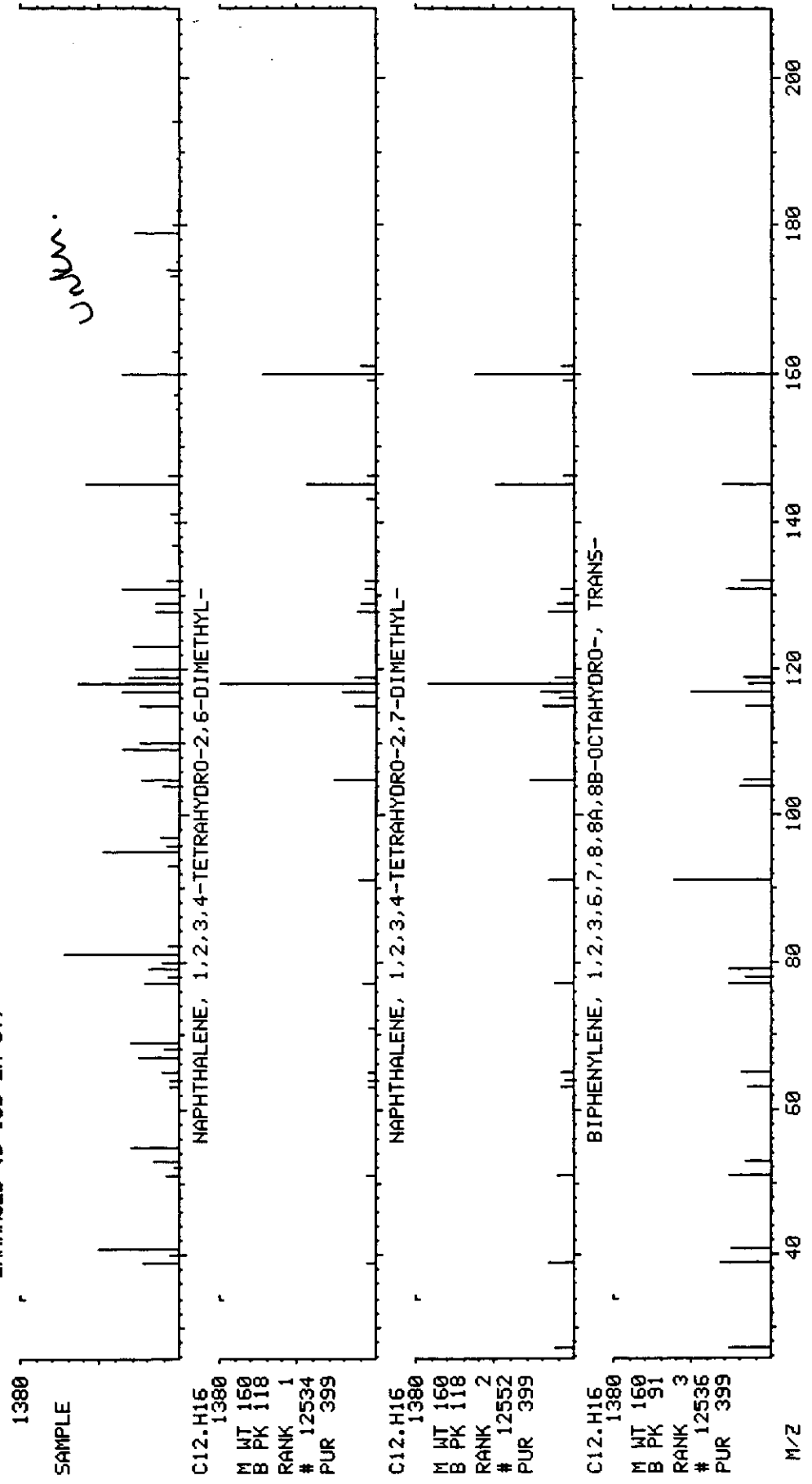
62231 spectra in LIBRARYNB searched for maximum PURITY
 897 matched at least 4 of the 16 largest peaks in the unknown

Rank In. Name
 1 12534 NAPHTHALENE, 1,2,3,4-TETRAHYDRO-2,6-DIMETHYL-
 2 12552 NAPHTHALENE, 1,2,3,4-TETRAHYDRO-2,7-DIMETHYL-
 3 12536 BIPHENYLENE, 1,2,3,6,7,8,8A,8B-OCTAHYDRO-, TRANS-
 4 41258 6,9-OCTADECADIYNOIC ACID, METHYL ESTER
 5 12526 1,4-DIMETHYL-1,2,3,4-TETRAHYDRONAPHTHALENE
 6 12577 BENZENE, 4-(2-BUTENYL)-1,2-DIMETHYL-, (E)-
 7 12573 NAPHTHALENE, 1,2,3,4-TETRAHYDRO-1,5-DIMETHYL-
 8 16368 BENZENE, (2,2-DIMETHYLCYCLOPENTYL)-
 9 12557 NAPHTHALENE, 1,2,3,4-TETRAHYDRO-6,7-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H16	160	118	399	824	438
2	C12.H16	160	118	399	841	442
3	C12.H16	160	91	399	804	447
4	C19.H30.O2	290	91	371	638	516
5	C12.H16	160	118	363	839	413
6	C12.H16	160	145	352	703	417
7	C12.H16	160	145	349	734	383
8	C13.H18	174	104	335	688	399
9	C12.H16	160	145	329	738	369

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	---	---	---	---	7524-63-2
2	---	---	---	---	13065-07-1
3	---	---	---	---	28229-15-4
4	---	---	---	---	56847-03-1
5	---	---	---	---	- -
6	---	---	---	---	54340-86-2
7	---	---	---	---	21564-91-0
8	---	---	---	---	19960-99-7
9	---	---	---	---	1076-61-5

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:59:00 + 7:47
 SAMPLE: S-MMS-3 1/35A/100M
 COND.S.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)
 DATA: 30068107 # 700
 CALLI: 30068107 # 3
 BASE M/Z: 81
 RIC: 28224.



1127

Library Search Data: 30068107 # 738 Base m/z: 57
 08/31/98 21:59:00 + 8:13 Cali: 30068107 # 3 RIC: 147968.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 712 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 25991 DODECANE, 2,6,10-TRIMETHYL-
 2 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 3 26001 DODECANE, 2,7,10-TRIMETHYL-
 4 37462 HEPTADECANE, 2,6-DIMETHYL-
 5 22535 DODECANE, 4,6-DIMETHYL-
 6 25997 PENTADECANE
 7 25994 DODECANE, 2,6,11-TRIMETHYL-
 8 22530 TETRADECANE
 9 15352 UNDECANE, 2-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C15.H32	212	57	822	975	823
2	C21.H44	296	57	821	940	842
3	C15.H32	212	57	816	988	816
4	C19.H40	268	57	814	941	838
5	C14.H30	198	57	812	955	812
6	C15.H32	212	57	806	942	821
7	C15.H32	212	57	804	977	807
8	C14.H30	198	43	791	922	803
9	C12.H26	170	43	790	942	796

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	3891-98-3
2	—	—	—	—	18344-37-1
3	—	—	—	—	74645-98-0
4	—	—	—	—	54105-67-8
5	—	—	—	—	61141-72-8
6	—	—	—	—	629-62-9
7	—	—	—	—	31295-56-4
8	—	—	—	—	629-59-4
9	—	—	—	—	7045-71-8

BASE M/Z: 57
RIC: 147958.

DATA: 30068107 # 738
CALI: 30068107 # 3

MID LIBRARY SEARCH (LIBRARYNB)

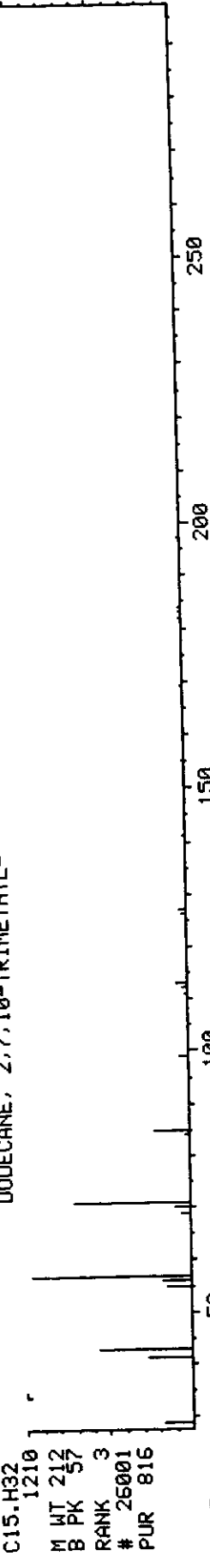
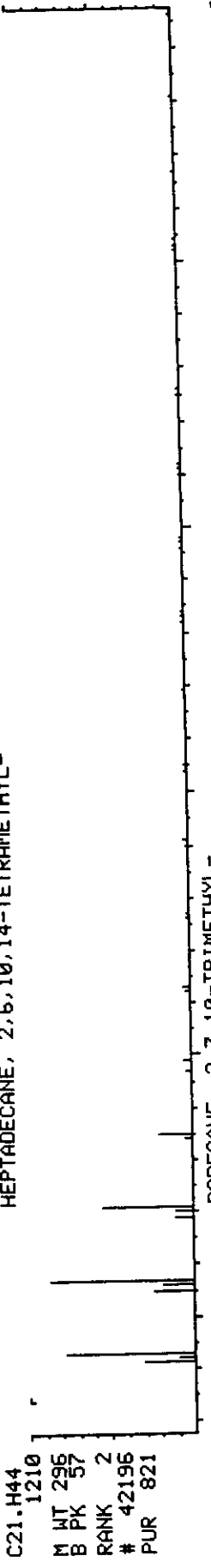
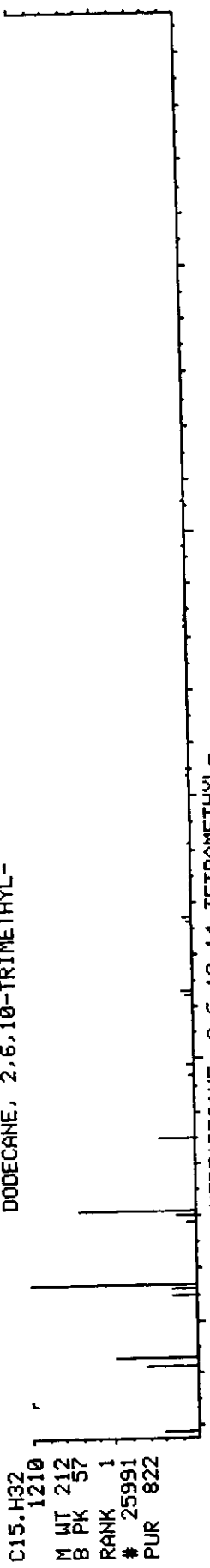
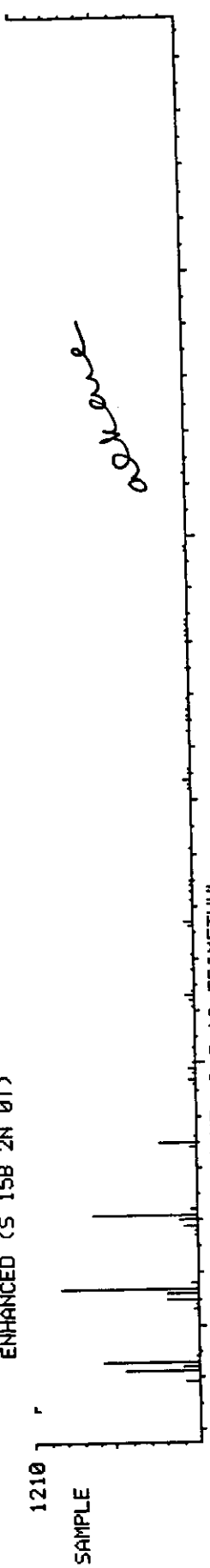
08/31/98 21:59:00 + 8:13

SAMPLE: S-MMS-3 1/35A/100M

CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA) /1/35A NA M

ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068107 # 764 Base m/z: 43
 08/31/98 21:59:00 + 8:30 Cali: 30068107 # 3 RIC: 227584.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

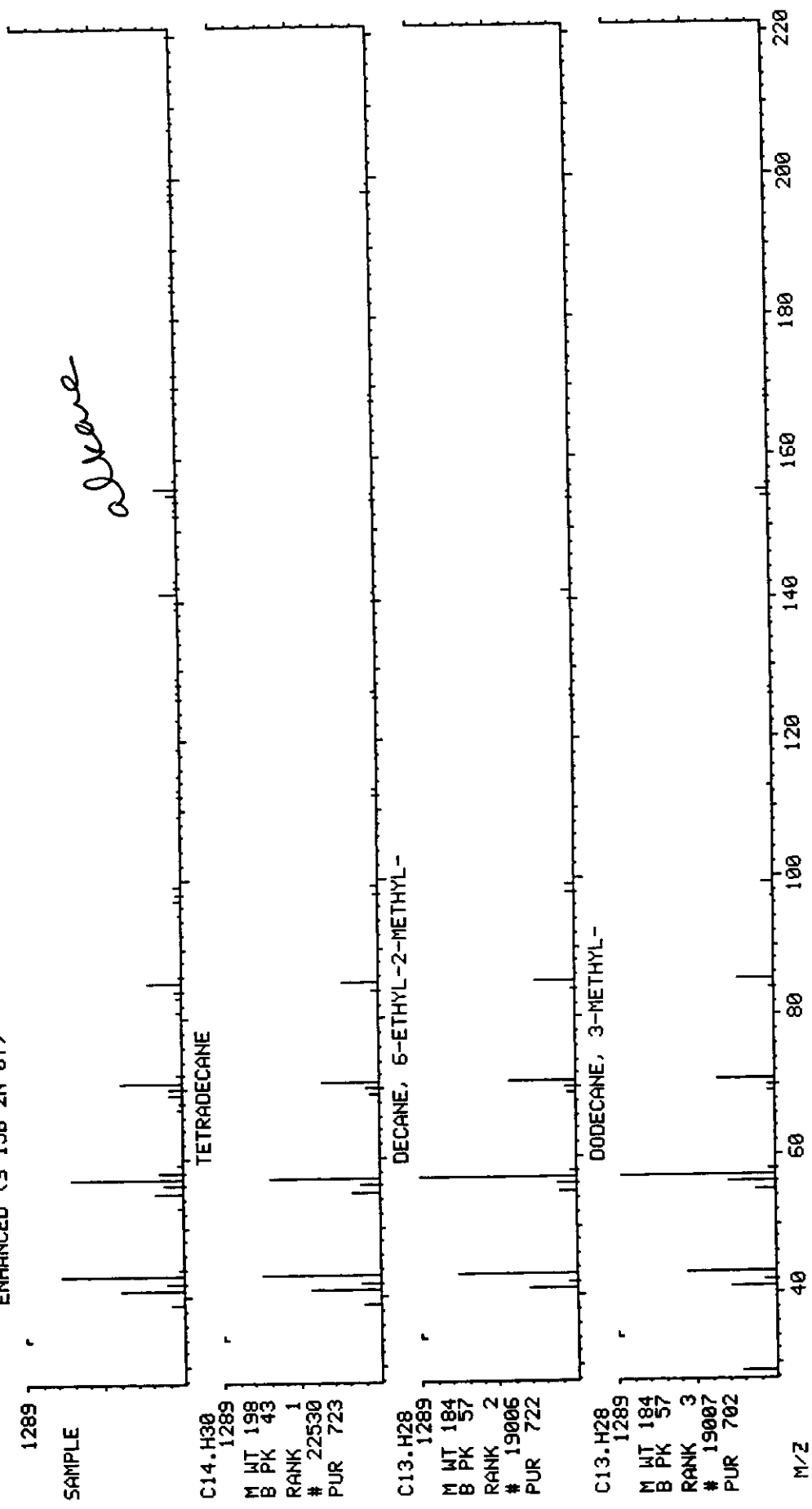
62231 spectra in LIBRARYNB searched for maximum PURITY
 672 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 22530 TETRADECANE
 2 19006 DECANE, 6-ETHYL-2-METHYL-
 3 19007 DODECANE, 3-METHYL-
 4 11612 NONANE, 2,5-DIMETHYL-
 5 11607 UNDECANE
 6 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 7 39681 1-iodoundecane
 8 18990 UNDECANE, 2,9-DIMETHYL-
 9 37252 DECANE, 1-iodo-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C14.H30	198	43	723	951	726
2	C13.H28	184	57	722	956	722
3	C13.H28	184	57	702	896	702
4	C11.H24	156	57	701	861	701
5	C11.H24	156	43	698	966	698
6	C21.H44	296	57	688	915	715
7	C11.H23.I	282	57	688	860	694
8	C13.H28	184	57	688	867	688
9	C10.H21.I	268	57	686	883	691

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-59-4
2	—	—	—	—	62108-21-8
3	—	—	—	—	17312-57-1
4	—	—	—	—	17302-27-1
5	—	—	—	—	1120-21-4
6	—	—	—	—	18344-37-1
7	—	—	—	—	4282-44-4
8	—	—	—	—	17301-26-7
9	—	—	—	—	2050-77-3

MID LIBRARY SEARCH <LIBRARY.NB>
 08/31/98 21:59:00 + 8:30
 SAMPLE: 5-MM5-3 1/35A/100M
 COND.S.: UG/ML *100%/100% *(NA/NA)/1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068107 # 764
 CALI: 30068107 # 3
 BASE M/Z: 43
 RIC: 227584.



1151

Library Search Data: 30068107 # 776 Base m/z: 156
 08/31/98 21:59:00 + 8:38 Cali: 30068107 # 3 RIC: 56512.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 277 matched at least 5 of the 16 largest peaks in the unknown

Rank	In.	Name
1	11625	NAPHTHALENE, 1,3-DIMETHYL-
2	11618	NAPHTHALENE, 2,3-DIMETHYL-
3	11629	NAPHTHALENE, 1,5-DIMETHYL-
4	11619	NAPHTHALENE, 1,2-DIMETHYL-
5	11615	NAPHTHALENE, 2,7-DIMETHYL-
6	11628	NAPHTHALENE, 1,6-DIMETHYL-
7	11621	NAPHTHALENE, 1,8-DIMETHYL-
8	11620	NAPHTHALENE, 2,6-DIMETHYL-
9	11624	NAPHTHALENE, 1,7-DIMETHYL-

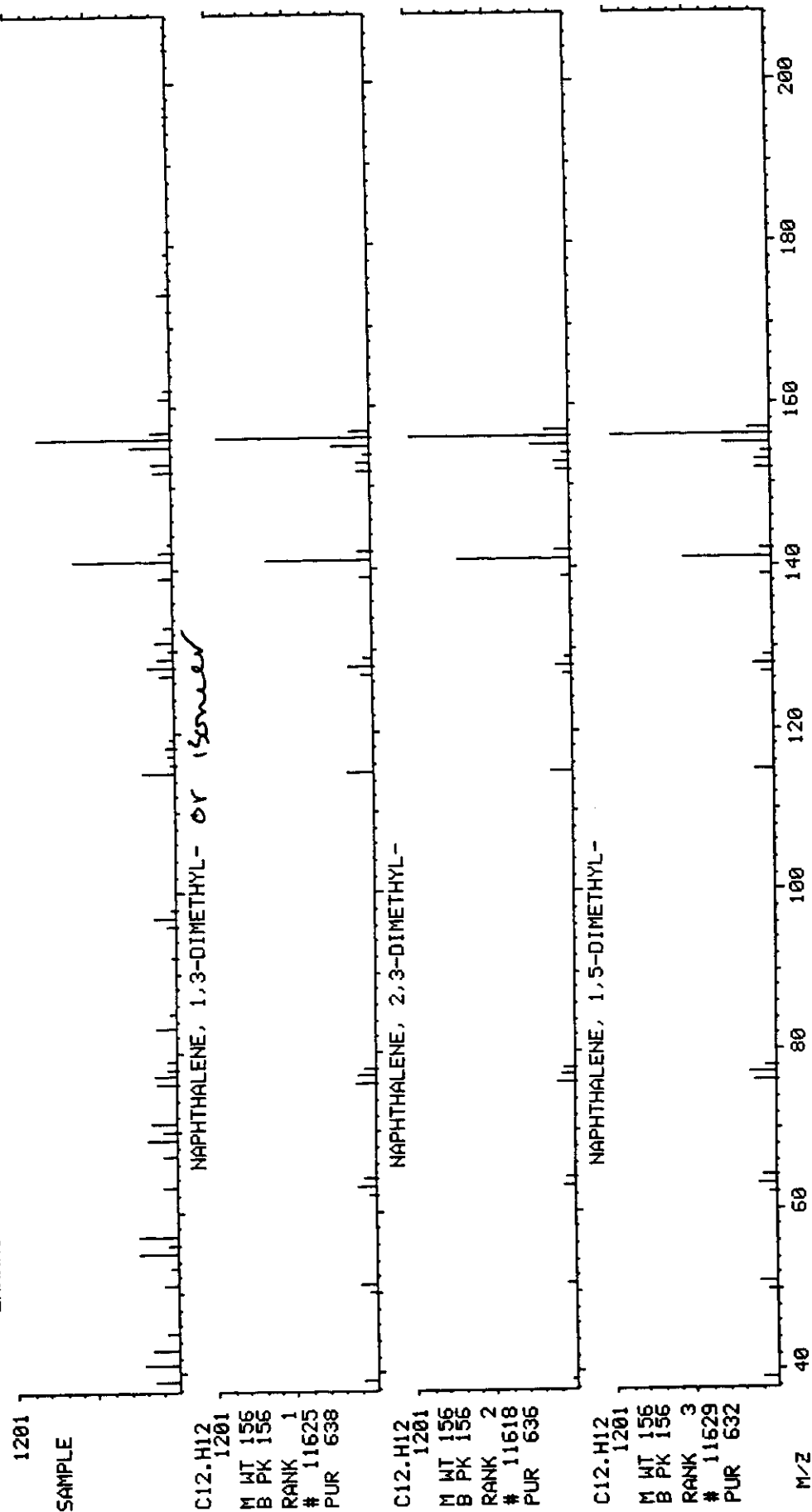
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C12.H12	156	156	638	958	660
2	C12.H12	156	156	636	957	651
3	C12.H12	156	156	632	950	658
4	C12.H12	156	156	630	948	646
5	C12.H12	156	156	627	943	653
6	C12.H12	156	156	624	939	649
7	C12.H12	156	156	619	943	647
8	C12.H12	156	156	617	949	648
9	C12.H12	156	156	616	936	652

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	575-41-7
2	—	—	—	—	581-40-8
3	—	—	—	—	571-61-9
4	—	—	—	—	573-98-8
5	—	—	—	—	582-16-1
6	—	—	—	—	575-43-9
7	—	—	—	—	569-41-5
8	—	—	—	—	581-42-0
9	—	—	—	—	575-37-1

DATA: 30068107 # 776
CALI: 30068107 # 3
BASE M/Z: 156
RIC: 56512.

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 21:59:00 + 8:38
SAMPLE: S-MM5-3 1/35A/100M
CONDOS.: UG/ML *100ML *100Z/100Z *(NA/NA) /1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



1-1-03

Library Search Data: 30068107 # 820 Base m/z: 57
 08/31/98 21:59:00 + 9:07 Cali: 30068107 # 3 RIC: 145408.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 855 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 25997 PENTADECANE
 4 25991 DODECANE, 2,6,10-TRIMETHYL-
 5 22530 TETRADECANE
 6 18987 UNDECANE, 2,8-DIMETHYL-
 7 19016 UNDECANE, 4,7-DIMETHYL-
 8 15353 2,6-DIMETHYLDECANE
 9 37465 NONADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C21.H44	296	57	837	980	844
2	C19.H40	268	57	808	952	824
3	C15.H32	212	57	787	946	799
4	C15.H32	212	57	782	964	783
5	C14.H30	198	43	782	938	793
6	C13.H28	184	43	782	964	785
7	C13.H28	184	43	780	950	781
8	C12.H26	170	43	779	925	788
9	C19.H40	268	57	778	928	810

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	18344-37-1
2	—	—	—	—	54105-67-8
3	—	—	—	—	629-62-9
4	—	—	—	—	3891-98-3
5	—	—	—	—	629-59-4
6	—	—	—	—	17301-25-6
7	—	—	—	—	17301-32-5
8	—	—	—	—	13150-81-7
9	—	—	—	—	629-92-5

DATA: 30068107 # 820
CALI: 30068107 # 3
BASE M/Z: 57
RIC: 145408.

MID LIBRARY SEARCH (LIBRARYNB)

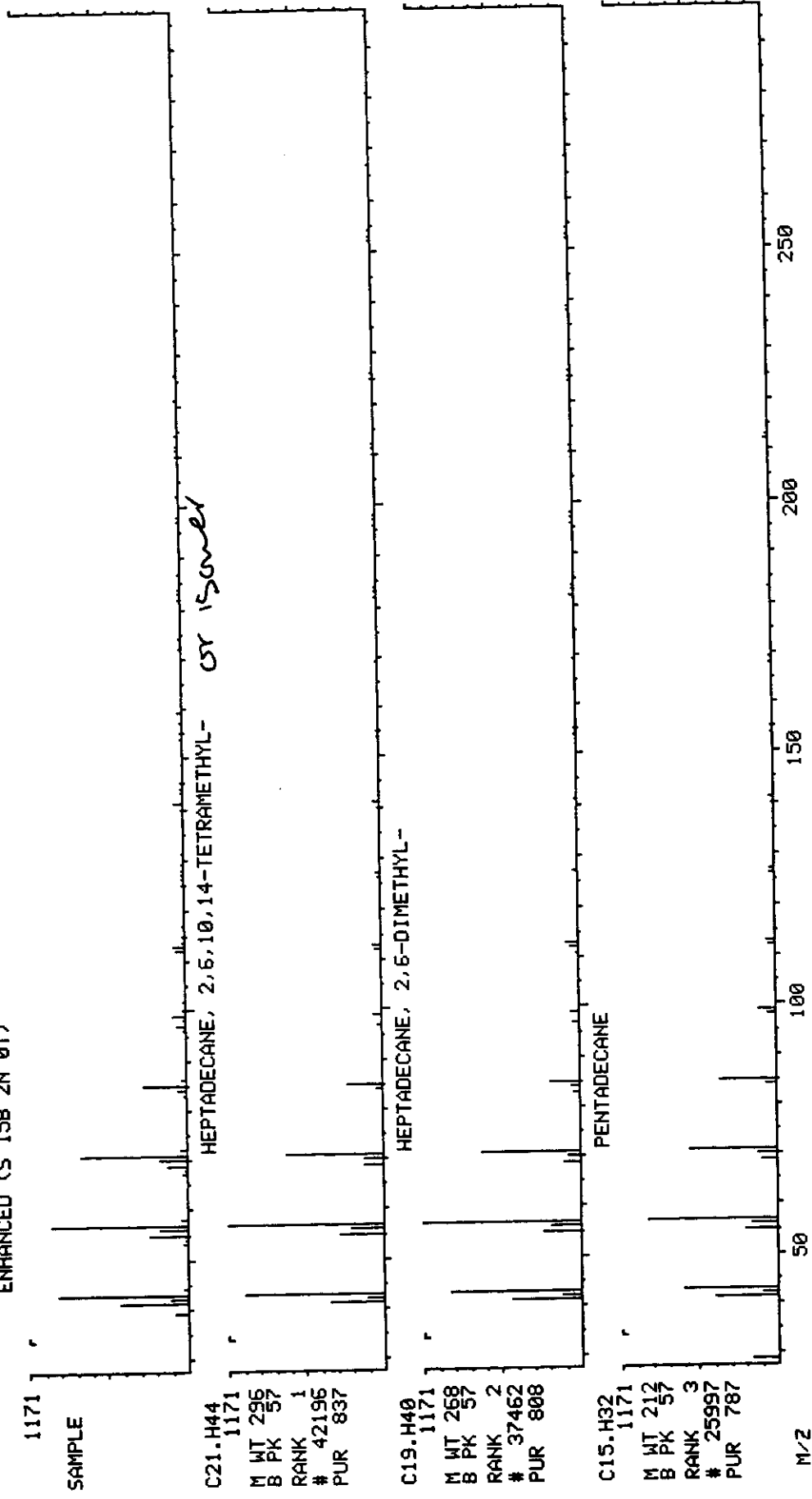
08/31/98 21:59:00 + 9:07

SAMPLE: S-MM5-3 1/35A/100M

INST. ID: F16

CONDOS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M

ENHANCED (S 15B 2N 0T)



1171
57
837

Library Search Data: 30068107 # 859 Base m/z: 57
 08/31/98 21:59:00 + 9:33 Cali: 30068107 # 3 RIC: 195840.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

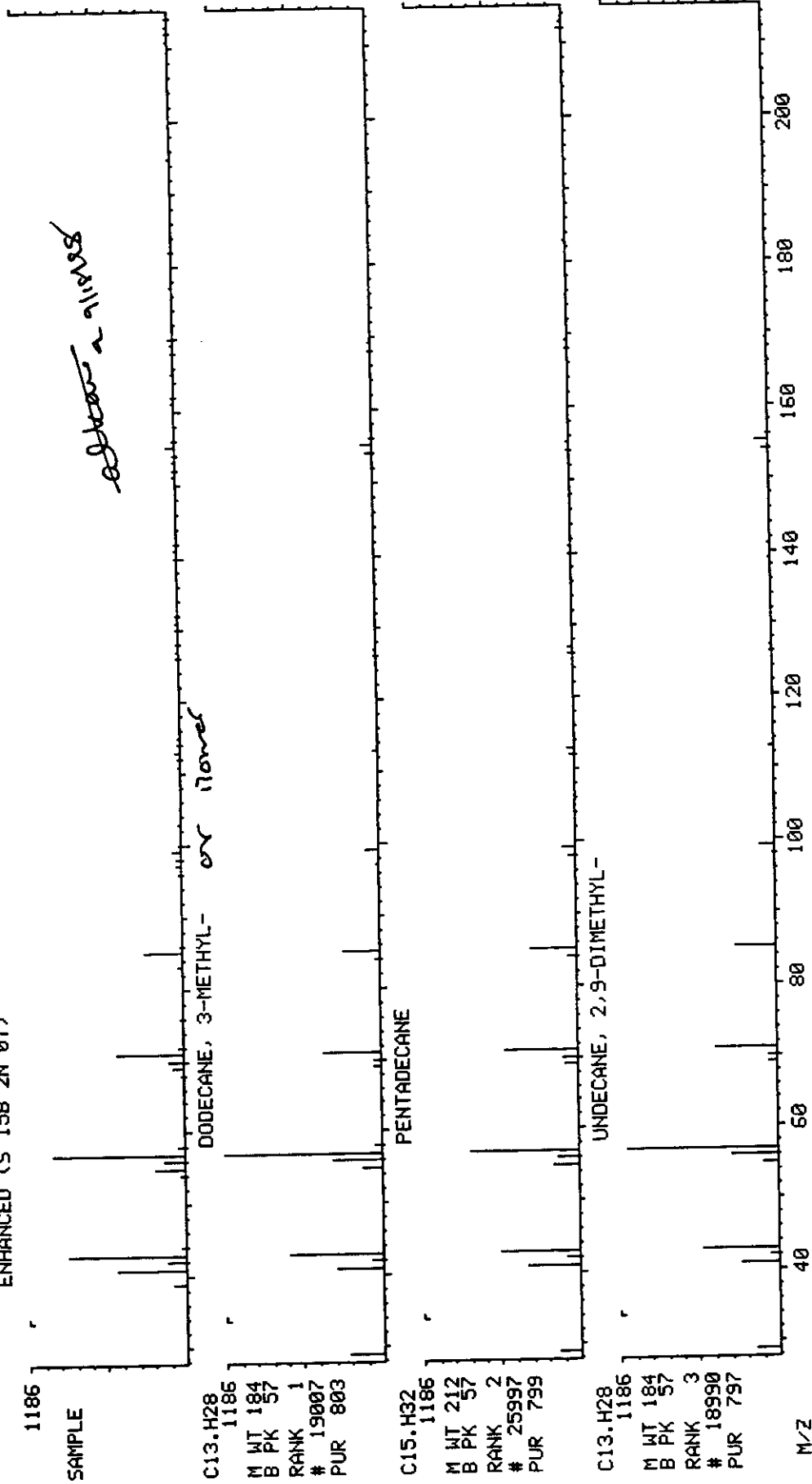
62231 spectra in LIBRARYNB searched for maximum PURITY
 559 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 19007 DODECANE, 3-METHYL-
 2 25997 PENTADECANE
 3 18990 UNDECANE, 2,9-DIMETHYL-
 4 19026 DECANE, 2,6,8-TRIMETHYL-
 5 29249 TETRADECANE, 6,9-DIMETHYL-
 6 39681 1-iodoundecane
 7 15352 UNDECANE, 2-METHYL-
 8 19006 DECANE, 6-ETHYL-2-METHYL-
 9 15357 DECANE, 2,9-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H28	184	57	803	960	810
2	C15.H32	212	57	799	941	818
3	C13.H28	184	57	797	957	799
4	C13.H28	184	57	796	946	799
5	C16.H34	226	57	792	934	810
6	C11.H23.I	282	57	792	968	800
7	C12.H26	170	43	792	936	803
8	C13.H28	184	57	792	958	800
9	C12.H26	170	43	790	962	790

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	17312-57-1
2	—	—	—	—	629-62-9
3	—	—	—	—	17301-26-7
4	—	—	—	—	62108-26-3
5	—	—	—	—	55045-13-1
6	—	—	—	—	4282-44-4
7	—	—	—	—	7045-71-8
8	—	—	—	—	62108-21-8
9	—	—	—	—	1002-17-1

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:59:00 + 9:33
 SAMPLE: 5-MM5-3 1/35A/100M
 COND.S.: UG/ML *100ML *100Z/100Z *(NA/NA) /1/35A NA M
 ENHANCED (S 158 2N 0T)
 DATA: 30068107 # 859
 CALI: 30068107 # 3
 BASE M/Z: 57
 RIC: 195840.



Library Search Data: 30068107 # 904 Base m/z: 57
 08/31/98 21:59:00 + 10:03 Cali: 30068107 # 3 RIC: 73344.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 782 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 40233 NONADECANOL
 2 42521 1-EICOSANOL
 3 37444 OXIRANE, HEXADECYL-
 4 37449 OCTADECANAL
 5 22482 2-TRIDECEN-1-OL, (E)-
 6 40193 CIS-9,10-EPOXYOCTADECAN-1-OL
 7 25971 OXIRANE, DODECYL-
 8 32052 OXIRANE, TETRADECYL-
 9 42180 OCTADECANE, 1-(ETHENYLOXY)-

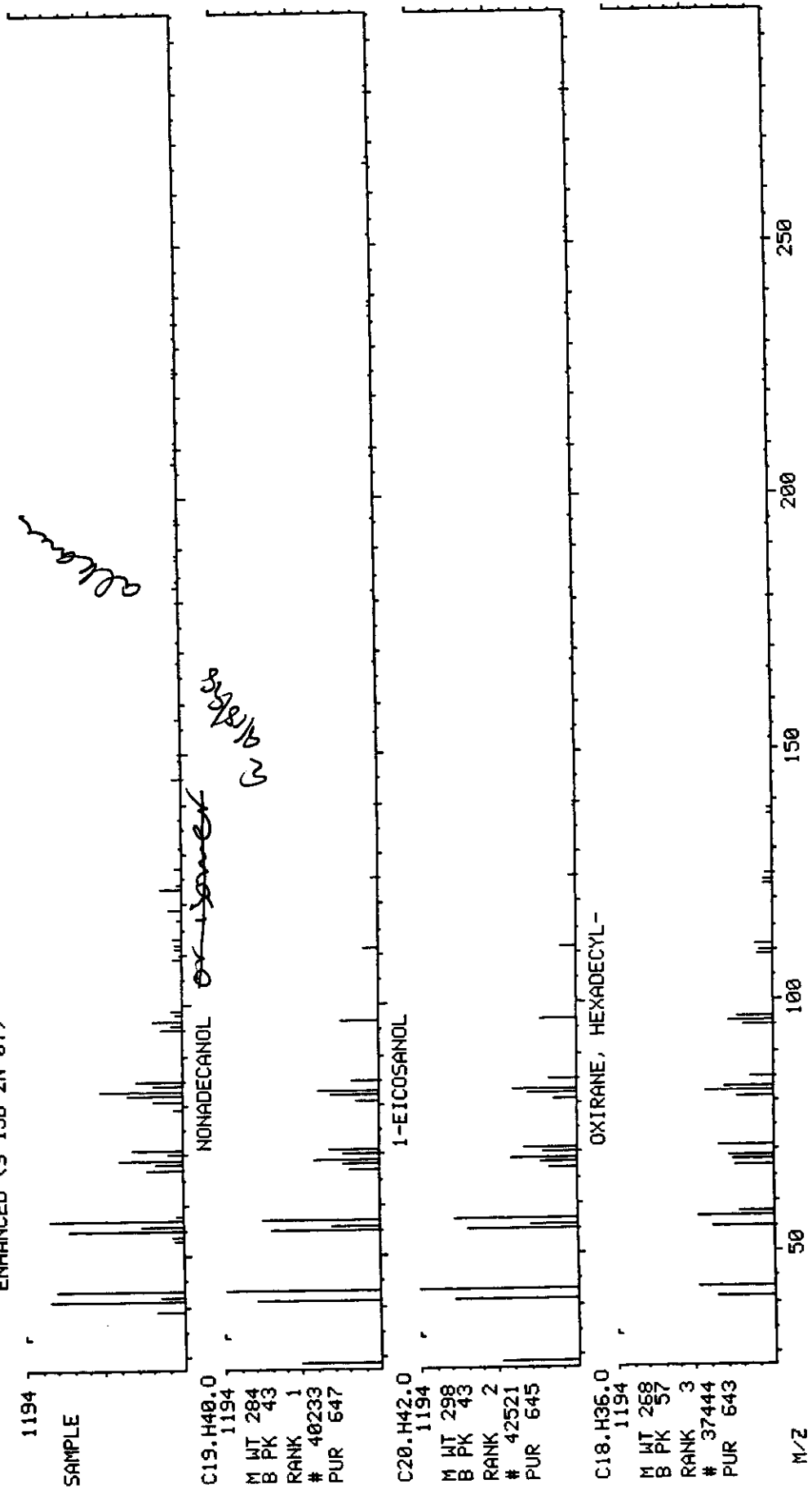
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40.O	284	43	647	923	686
2	C20.H42.O	298	43	645	919	688
3	C18.H36.O	268	57	643	867	672
4	C18.H36.O	268	43	640	873	683
5	C13.H26.O	198	57	637	923	648
6	C18.H36.O2	284	55	636	906	666
7	C14.H28.O	212	41	636	911	642
8	C16.H32.O	240	41	630	942	630
9	C20.H40.O	296	43	621	930	656

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	52783-43-4
2	—	—	—	—	629-96-9
3	—	—	—	—	7390-81-0
4	—	—	—	—	638-66-4
5	—	—	—	—	74962-98-4
6	—	—	—	—	13980-12-6
7	—	—	—	—	3234-28-4
8	—	—	—	—	7320-37-8
9	—	—	—	—	930-02-9

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 21:59:00 + 10:03
 SAMPLE: S-MM5-3 1/35A/100M
 COND.: UG/ML *100ML *100Z/100Z *KNA/NA >1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068107 # 904
 CALI: 30068107 # 3

BASE M/Z: 57
 RIC: 73344.



1150

Library Search Data: 30068107 # 948 Base m/z: 57
 08/31/98 21:59:00 + 10:33 Cali: 30068107 # 3 RIC: 229888.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 360 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 15969 HYDROXYLAMINE, O-DECYL-
 5 19016 UNDECANE, 4,7-DIMETHYL-
 6 25997 PENTADECANE
 7 25994 DODECANE, 2,6,11-TRIMETHYL-
 8 26001 DODECANE, 2,7,10-TRIMETHYL-
 9 19054 UNDECANE, 2,6-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	843	948	870
2	C19.H40	268	57	841	941	868
3	C21.H44	296	57	832	943	856
4	C10.H23.O.N	173	43	832	964	835
5	C13.H28	184	43	826	965	826
6	C15.H32	212	57	820	945	854
7	C15.H32	212	57	807	964	811
8	C15.H32	212	57	804	960	809
9	C13.H28	184	57	800	908	810

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	18344-37-1
4	—	—	—	—	29812-79-1
5	—	—	—	—	17301-32-5
6	—	—	—	—	629-62-9
7	—	—	—	—	31295-56-4
8	—	—	—	—	74645-98-0
9	—	—	—	—	17301-23-4

MID LIBRARY SEARCH (LIBRARYNE)
 08/31/98 21:59:00 + 10:33
 SAMPLE: S-MMS-3 1/35A/100M
 COND.: UG/ML *100ML *100% /100% *(NA/NA) /1/35A NA M
 ENHANCED (S 158 2N 0T)
 INST. ID: F16
 DATA: 30068107 # 948
 CALL: 30068107 # 3
 BASE M/Z: 57
 RIC: 229888.

1032
 SAMPLE

C19.H40
 1032

M WT 258
 B PK 57
 RANK 1
 # 37465
 PUR 843

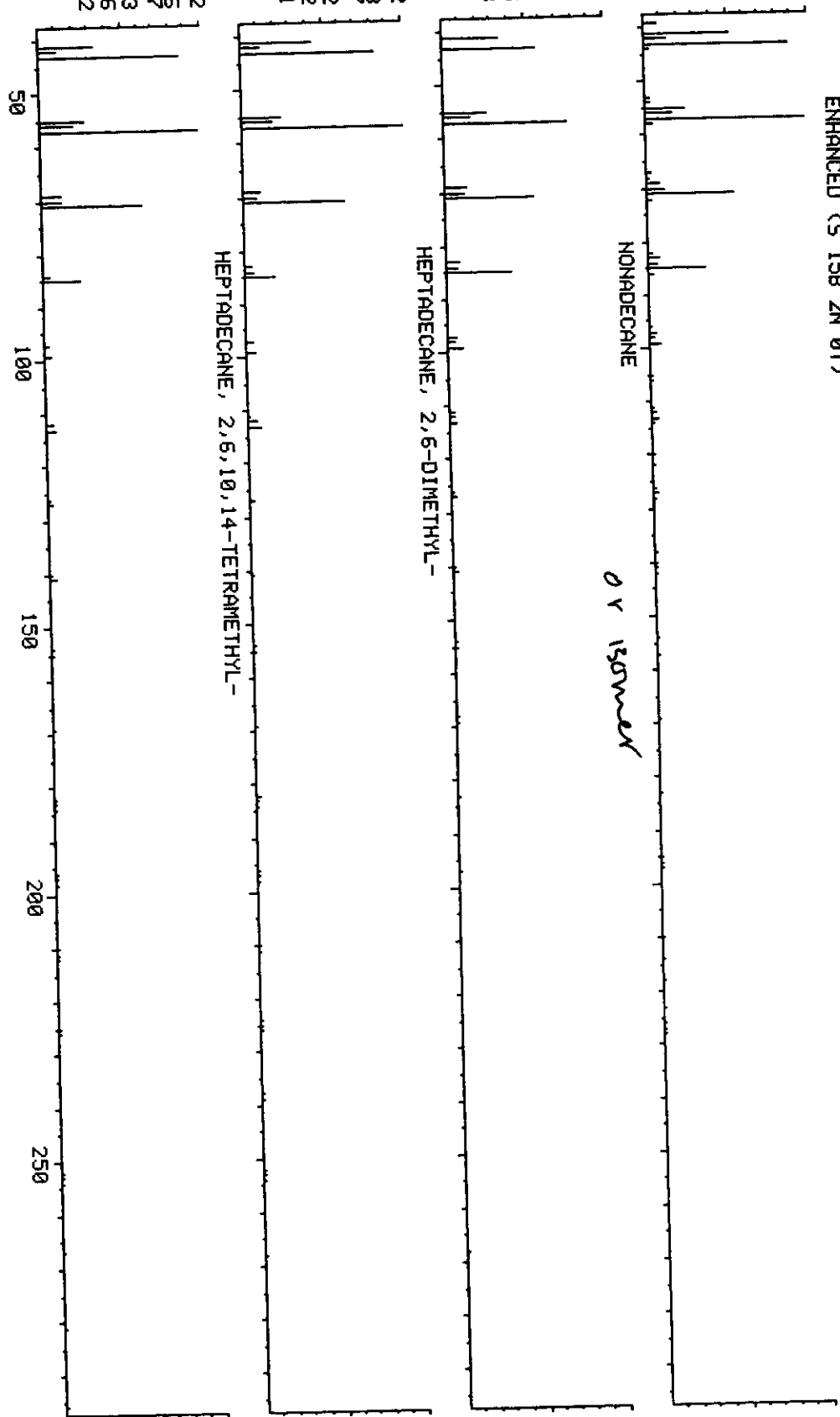
C19.H40
 1032

M WT 258
 B PK 57
 RANK 2
 # 37462
 PUR 841

C21.H44
 1032

M WT 296
 B PK 57
 RANK 3
 # 42196
 PUR 832

M/Z



Library Search Data: 30068107 # 988 Base m/z: 57
 08/31/98 21:59:00 + 10:59 Cali: 30068107 # 3 RIC: 164608.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 644 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 32059 HEPTADECANE
 5 25997 PENTADECANE
 6 19054 UNDECANE, 2,6-DIMETHYL-
 7 22530 TETRADECANE
 8 25996 TRIDECANE, 4,8-DIMETHYL-
 9 29263 HEXADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	832	945	852
2	C19.H40	268	57	832	956	843
3	C21.H44	296	57	826	957	838
4	C17.H36	240	57	824	941	836
5	C15.H32	212	57	824	953	846
6	C13.H28	184	57	822	951	829
7	C14.H30	198	43	820	947	824
8	C15.H32	212	57	814	935	830
9	C16.H34	226	57	810	928	849

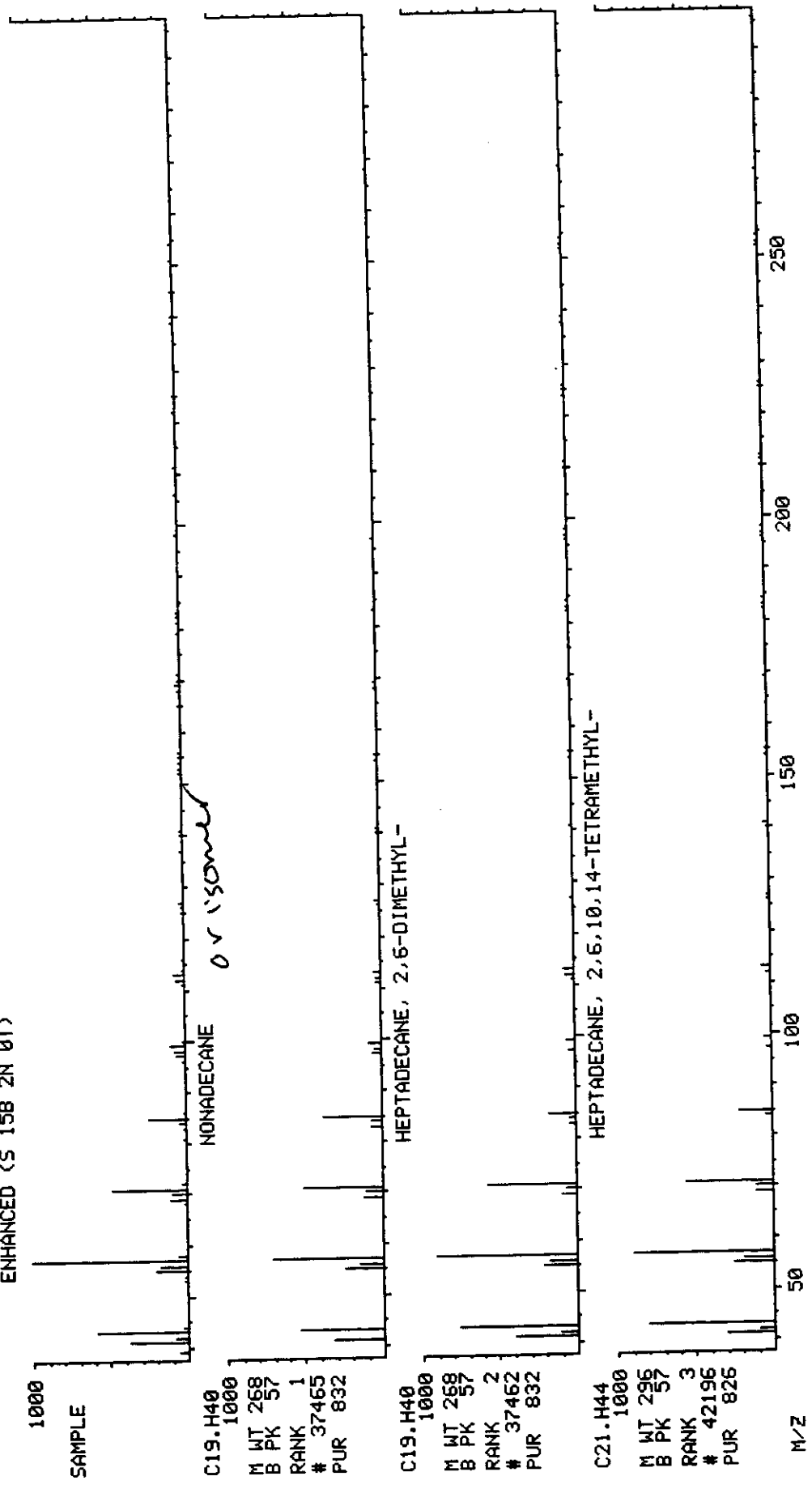
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	18344-37-1
4	—	—	—	—	629-78-7
5	—	—	—	—	629-62-9
6	—	—	—	—	17301-23-4
7	—	—	—	—	629-59-4
8	—	—	—	—	55030-62-1
9	—	—	—	—	544-76-3

BASE M/Z: 57
RIC: 164608.

DATA: 30068107 # 988
CALI: 30068107 # 3

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 21:59:00 + 10:59
SAMPLE: 5-MM5-3 1/35A/100M
CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068107 #1036 Base m/z: 57
 08/31/98 21:59:00 + 11:31 Cali: 30068107 # 3 RIC: 240640.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 713 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 3 37465 NONADECANE
 4 26001 DODECANE, 2,7,10-TRIMETHYL-
 5 25991 DODECANE, 2,6,10-TRIMETHYL-
 6 37466 PENTADECANE, 2,6,10,14-TETRAMETHYL-
 7 22535 DODECANE, 4,6-DIMETHYL-
 8 25994 DODECANE, 2,6,11-TRIMETHYL-
 9 29264 TRIDECANE, 5-PROPYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	875	976	882
2	C21.H44	296	57	833	949	857
3	C19.H40	268	57	831	940	844
4	C15.H32	212	57	826	970	830
5	C15.H32	212	57	825	973	827
6	C19.H40	268	71	820	946	834
7	C14.H30	198	57	818	961	820
8	C15.H32	212	57	809	949	816
9	C16.H34	226	57	806	947	827

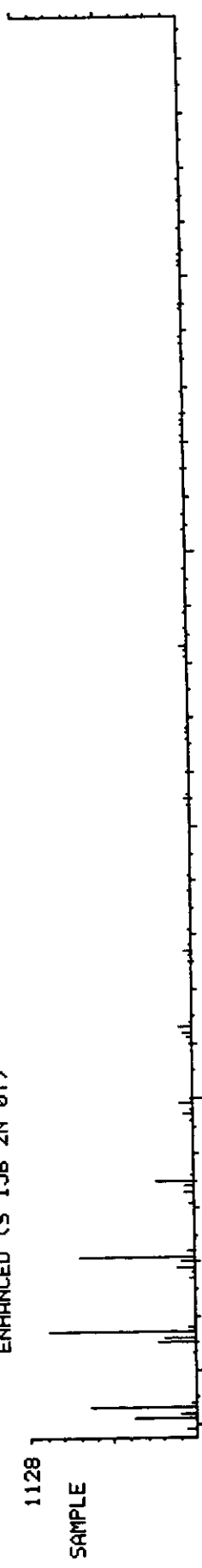
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	18344-37-1
3	—	—	—	—	629-92-5
4	—	—	—	—	74645-98-0
5	—	—	—	—	3891-98-3
6	—	—	—	—	1921-70-6
7	—	—	—	—	61141-72-8
8	—	—	—	—	31295-56-4
9	—	—	—	—	55045-11-9

DATA: 30068107 #1036
CALI: 30068107 # 3

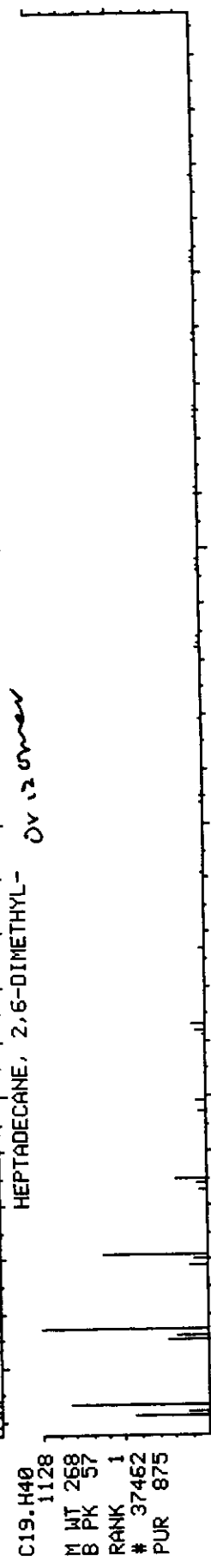
BASE M/Z: 57
RIC: 240640.

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 21:59:00 + 11:31
SAMPLE: S-MM5-3 1/35A/100M
COND5.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)

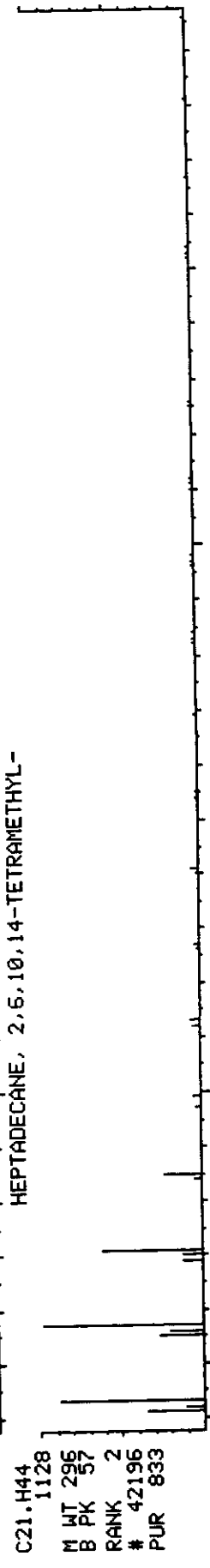
INST. ID: F16



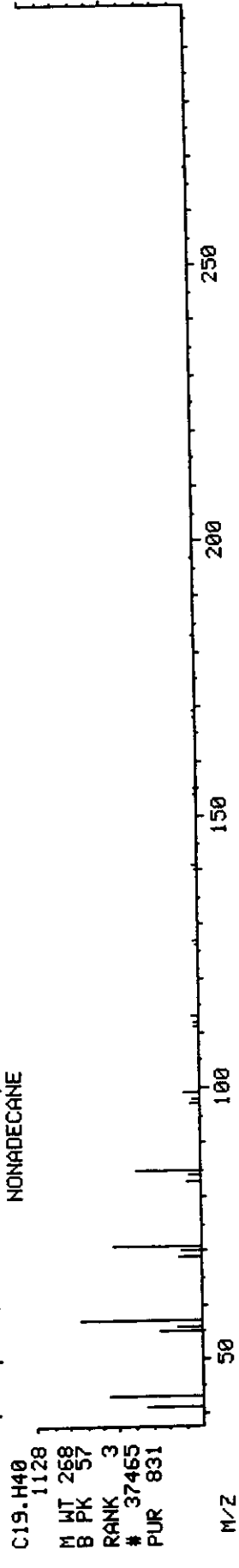
over



HEPTADECANE, 2,6,10,14-TETRAMETHYL-



NONADECANE



Library Search Data: 30068107 #1054 Base m/z: 55
 08/31/98 21:59:00 + 11:43 Cali: 30068107 # 3 RIC: 63616.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 796 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 28768 3-HEXADECENE, (Z)-
 2 21963 3-TETRADECENE, (E)-
 3 28772 7-HEXADECENE, (Z)-
 4 34411 5-OCTADECENE, (E)-
 5 34410 3-OCTADECENE, (E)-
 6 21962 3-TETRADECENE, (Z)-
 7 26409 2-HEXYL-1-OCTANOL
 8 34418 9-OCTADECENE, (E)-
 9 21976 7-TETRADECENE, (E)-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C16.H32	224	55	722	914	749
2	C14.H28	196	41	718	936	723
3	C16.H32	224	55	716	909	739
4	C18.H36	252	55	716	908	741
5	C18.H36	252	69	716	909	744
6	C14.H28	196	41	716	933	719
7	C14.H30.O	214	57	714	928	726
8	C18.H36	252	55	713	905	738
9	C14.H28	196	55	709	929	712

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	34303-81-6
2	—	—	—	—	41446-68-8
3	—	—	—	—	35507-09-6
4	—	—	—	—	7206-21-5
5	—	—	—	—	7206-19-1
6	—	—	—	—	41446-67-7
7	—	—	—	—	- -
8	—	—	—	—	7206-25-9
9	—	—	—	—	41446-63-3

BASE M/Z: 55
RIC: 63616.

DATA: 30068107 #1054
CALI: 30068107 # 3

MID LIBRARY SEARCH (LIBRARYNB)

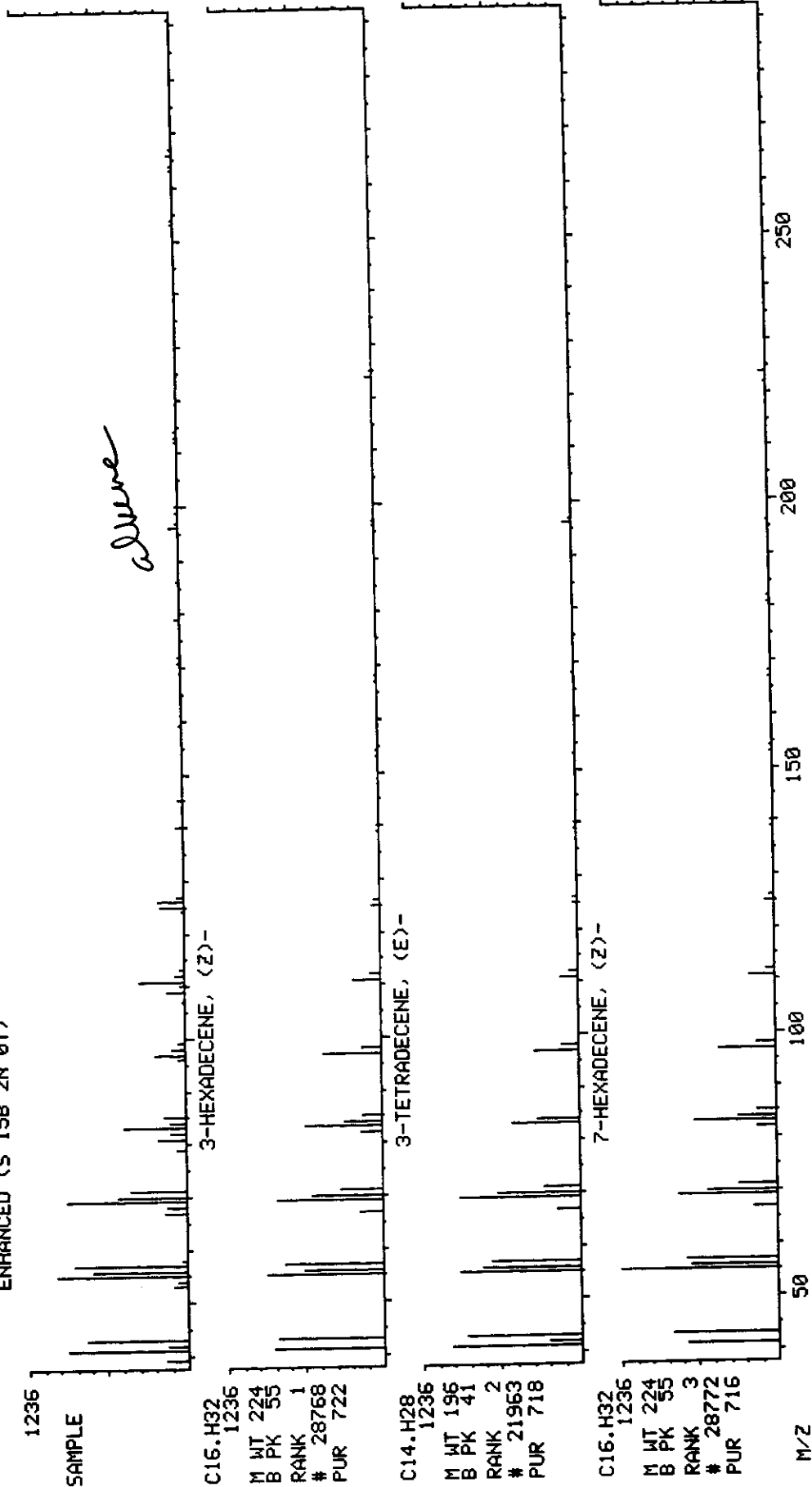
08/31/98 21:59:00 + 11:43

SAMPLE: S-MM5-3 1/35A/100M

INST. ID: F16

CONDS.: UG/ML *100ML *100% /100% *(NA/NA) /1/35A NA M

ENHANCED (S 15B 2N 0T)



Library Search Data: 30068107 #1114 Base m/z: 57
 08/31/98 21:59:00 + 12:23 Cali: 30068107 # 3 RIC: 105344.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 360 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name
 1 15969 HYDROXYLAMINE, O-DECYL-
 2 37465 NONADECANE
 3 37462 HEPTADECANE, 2,6-DIMETHYL-
 4 19016 UNDECANE, 4,7-DIMETHYL-
 5 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 6 25997 PENTADECANE
 7 19523 1-DECANOL, 2-ETHYL-
 8 11602 OCTANE, 2,4,6-TRIMETHYL-
 9 46161 TRICOSANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C10.H23.O.N	173	43	804	962	806
2	C19.H40	268	57	797	940	820
3	C19.H40	268	57	794	930	828
4	C13.H28	184	43	781	963	782
5	C21.H44	296	57	780	939	811
6	C15.H32	212	57	767	937	805
7	C12.H26.O	186	57	763	921	776
8	C11.H24	156	57	760	968	760
9	C23.H48	324	43	758	886	842

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	29812-79-1
2	—	—	—	—	629-92-5
3	—	—	—	—	54105-67-8
4	—	—	—	—	17301-32-5
5	—	—	—	—	18344-37-1
6	—	—	—	—	629-62-9
7	—	—	—	—	21078-65-9
8	—	—	—	—	62016-37-9
9	—	—	—	—	638-67-5

DATA: 30058107 #1114
CALI: 30058107 # 3

BASE M/Z: 57
RIC: 105344.

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 21:59:00 + 12:23

SAMPLE: 5-MM5-3 1/35A/100M

COND.: UG/ML *100ML *100Z/100Z *(NA/NA) 1/35A NA M

ENHANCED (S 15B 2N 0T)

INST. ID: F16

1047

SAMPLE

2-2-18/1998
all none

HYDROXYLAMINE, O-DECYL-

C10.H23.O.N

1047

M WT 173

B PK 43

RANK 1

15969

PUR 804

C19.H40

1047

M WT 268

B PK 57

RANK 2

37465

PUR 797

HEPTADECANE, 2,6-DIMETHYL-

C19.H40

1047

M WT 268

B PK 57

RANK 3

37462

PUR 794

M/Z

1100

Library Search Data: 30068107 #1119 Base m/z: 57
 08/31/98 21:59:00 + 12:27 Cali: 30068107 # 3 RIC: 136704.
 Sample: S-MMS-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 840 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 46161 TRICOSANE
 5 37466 PENTADECANE, 2,6,10,14-TETRAMETHYL-
 6 25991 DODECANE, 2,6,10-TRIMETHYL-
 7 32059 HEPTADECANE
 8 39859 OCTADECANE, 2,6-DIMETHYL-
 9 26001 DODECANE, 2,7,10-TRIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	839	964	839
2	C19.H40	268	57	835	948	839
3	C21.H44	296	57	816	937	830
4	C23.H48	324	43	815	920	856
5	C19.H40	268	71	802	937	806
6	C15.H32	212	57	799	968	800
7	C17.H36	240	57	796	929	822
8	C20.H42	282	57	794	931	799
9	C15.H32	212	57	794	967	794

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	18344-37-1
4	—	—	—	—	638-67-5
5	—	—	—	—	1921-70-6
6	—	—	—	—	3891-98-3
7	—	—	—	—	629-78-7
8	—	—	—	—	75163-97-2
9	—	—	—	—	74645-98-0

BASE M/Z: 57
RIC: 136704.

DATA: 30068107 #1119
CALI: 30068107 # 3

MID LIBRARY SEARCH <LIBRARYNB>

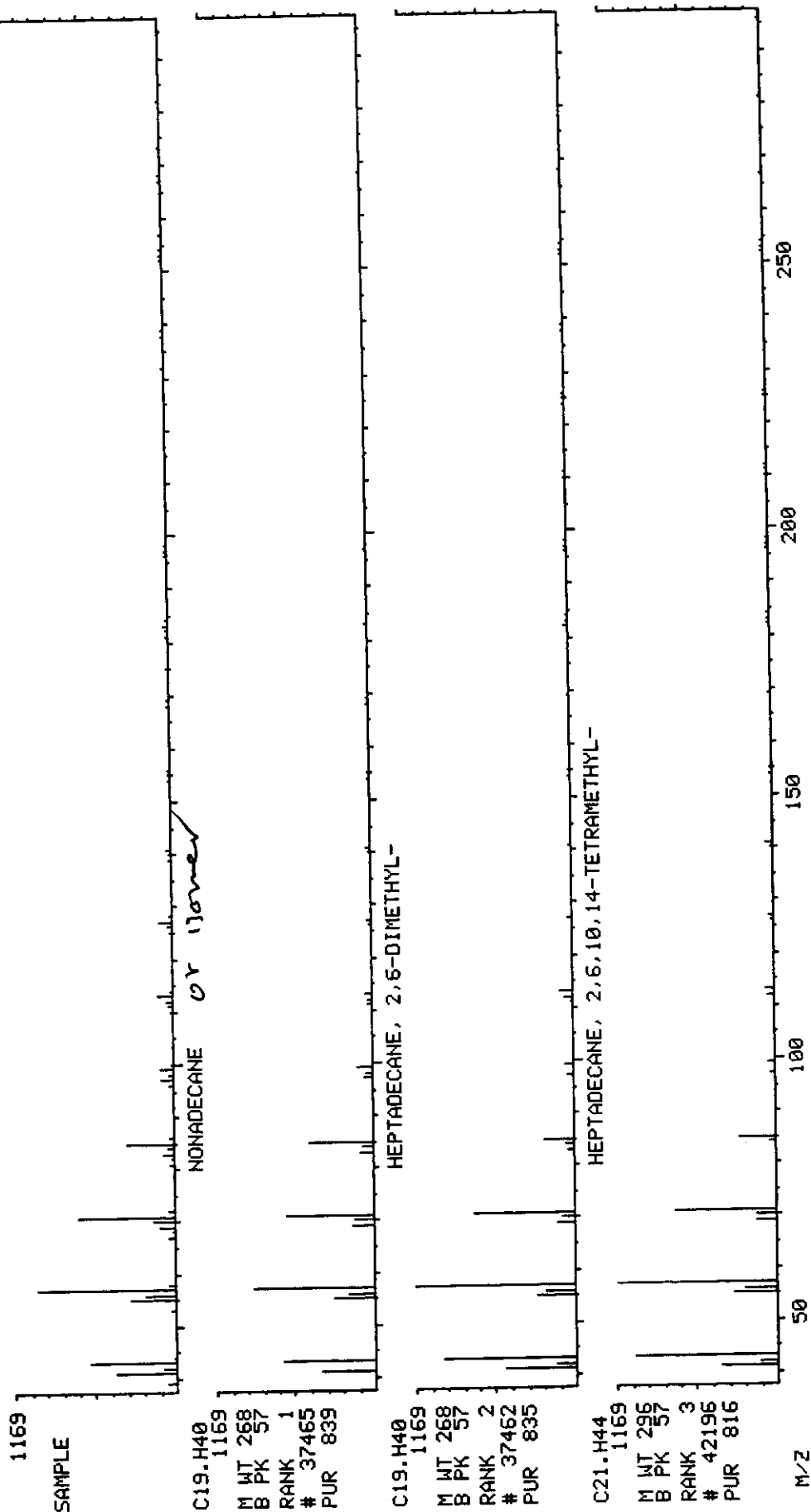
08/31/98 21:59:00 + 12:27

SAMPLE: S-MM5-3 1/35A/100M

INST. ID: F16

CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M

ENHANCED (S 15B 2N 0T)



1171

Library Search Data: 30068107 #1191 Base m/z: 57
 08/31/98 21:59:00 + 13:15 Cali: 30068107 # 3 RIC: 91264.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 294 matched at least 8 of the 16 largest peaks in the unknown

Rank In. Name

- 1 37465 NONADECANE
- 2 15969 HYDROXYLAMINE, O-DECYL-
- 3 37462 HEPTADECANE, 2,6-DIMETHYL-
- 4 19016 UNDECANE, 4,7-DIMETHYL-
- 5 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
- 6 11602 OCTANE, 2,4,6-TRIMETHYL-
- 7 19015 DECANE, 2,5,6-TRIMETHYL-
- 8 12074 1-HEPTANOL, 2-PROPYL-
- 9 25997 PENTADECANE

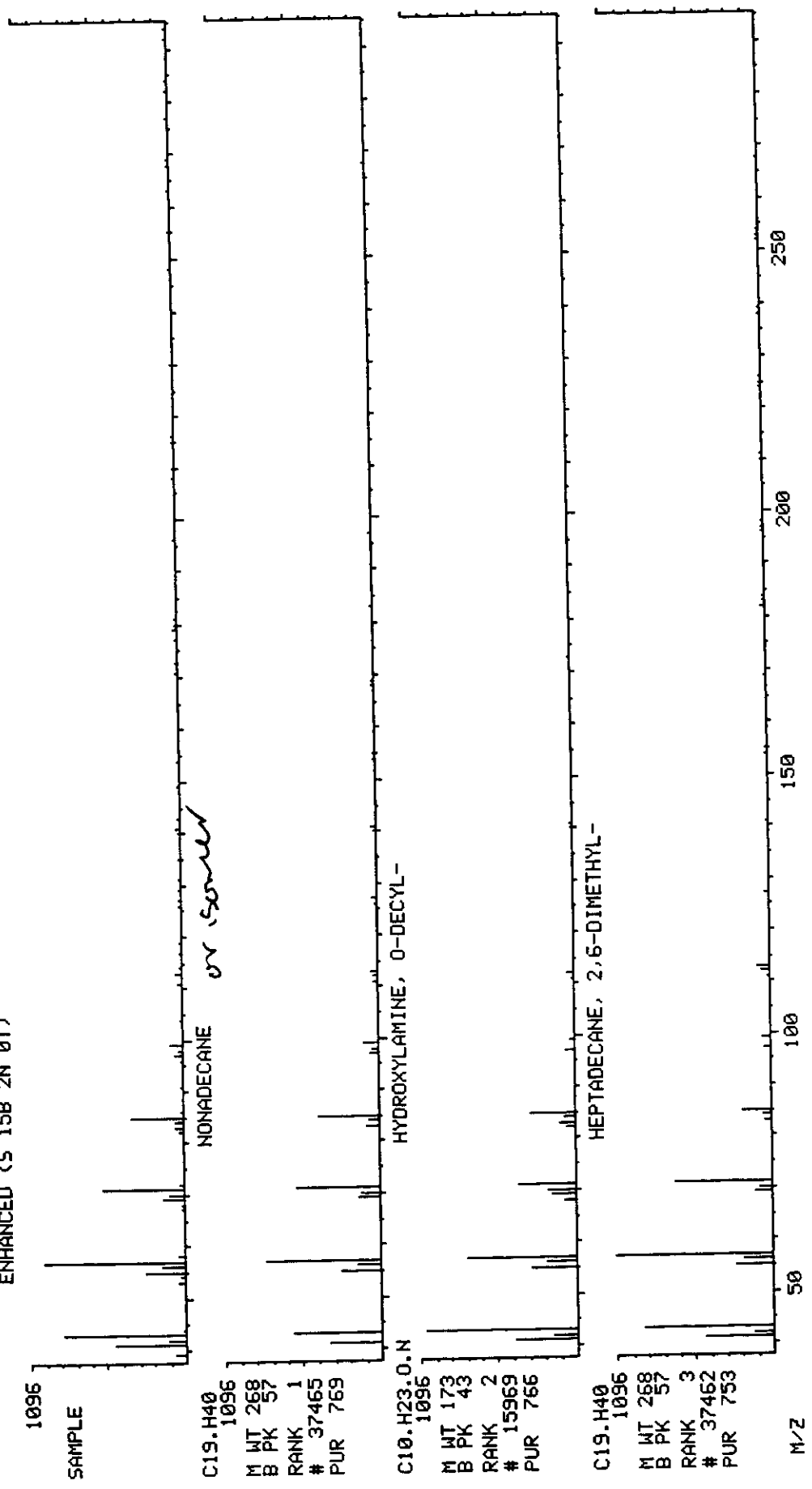
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	769	938	806
2	C10.H23.O.N	173	43	766	949	782
3	C19.H40	268	57	753	907	809
4	C13.H28	184	43	750	955	755
5	C21.H44	296	57	747	922	796
6	C11.H24	156	57	737	965	737
7	C13.H28	184	57	731	928	731
8	C10.H22.O	158	43	731	950	731
9	C15.H32	212	57	728	922	779

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	29812-79-1
3	—	—	—	—	54105-67-8
4	—	—	—	—	17301-32-5
5	—	—	—	—	18344-37-1
6	—	—	—	—	62016-37-9
7	—	—	—	—	62108-23-0
8	—	—	—	—	10042-59-8
9	—	—	—	—	629-62-9

DATA: 30068107 #1191
CALI: 30068107 # 3
BASE M/Z: 57
RIC: 91264.

MTD LIBRARY SEARCH (LIBRARYNB)
08/31/98 21:59:00 + 13:15
SAMPLE: S-MM5-3 1/35A/100M
COND5.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068107 #1335 Base m/z: 57
 08/31/98 21:59:00 + 14:51 Cali: 30068107 # 3 RIC: 45504.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Cords.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 715 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 37456 2-METHYLOCTADECANE
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 29263 HEXADECANE
 5 32058 HEXADECANE, 3-METHYL-
 6 37465 NONADECANE
 7 46161 TRICOSANE
 8 42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-
 9 34816 HEPTADECANE, 4-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	720	890	739
2	C19.H40	268	43	713	917	741
3	C21.H44	296	57	713	908	730
4	C16.H34	226	57	707	945	728
5	C17.H36	240	57	705	940	716
6	C19.H40	268	57	702	948	733
7	C23.H48	324	43	696	902	764
8	C21.H44	296	57	694	930	720
9	C18.H38	254	43	690	881	708

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	- -
3	—	—	—	—	18344-37-1
4	—	—	—	—	544-76-3
5	—	—	—	—	6418-43-5
6	—	—	—	—	629-92-5
7	—	—	—	—	638-67-5
8	—	—	—	—	54833-48-6
9	—	—	—	—	26429-11-8

BASE M/Z: 57
RIC: 45504.

DATA: 30068107 #1335
CALI: 30068107 # 3

MID LIBRARY SEARCH (LIBRARYNB)

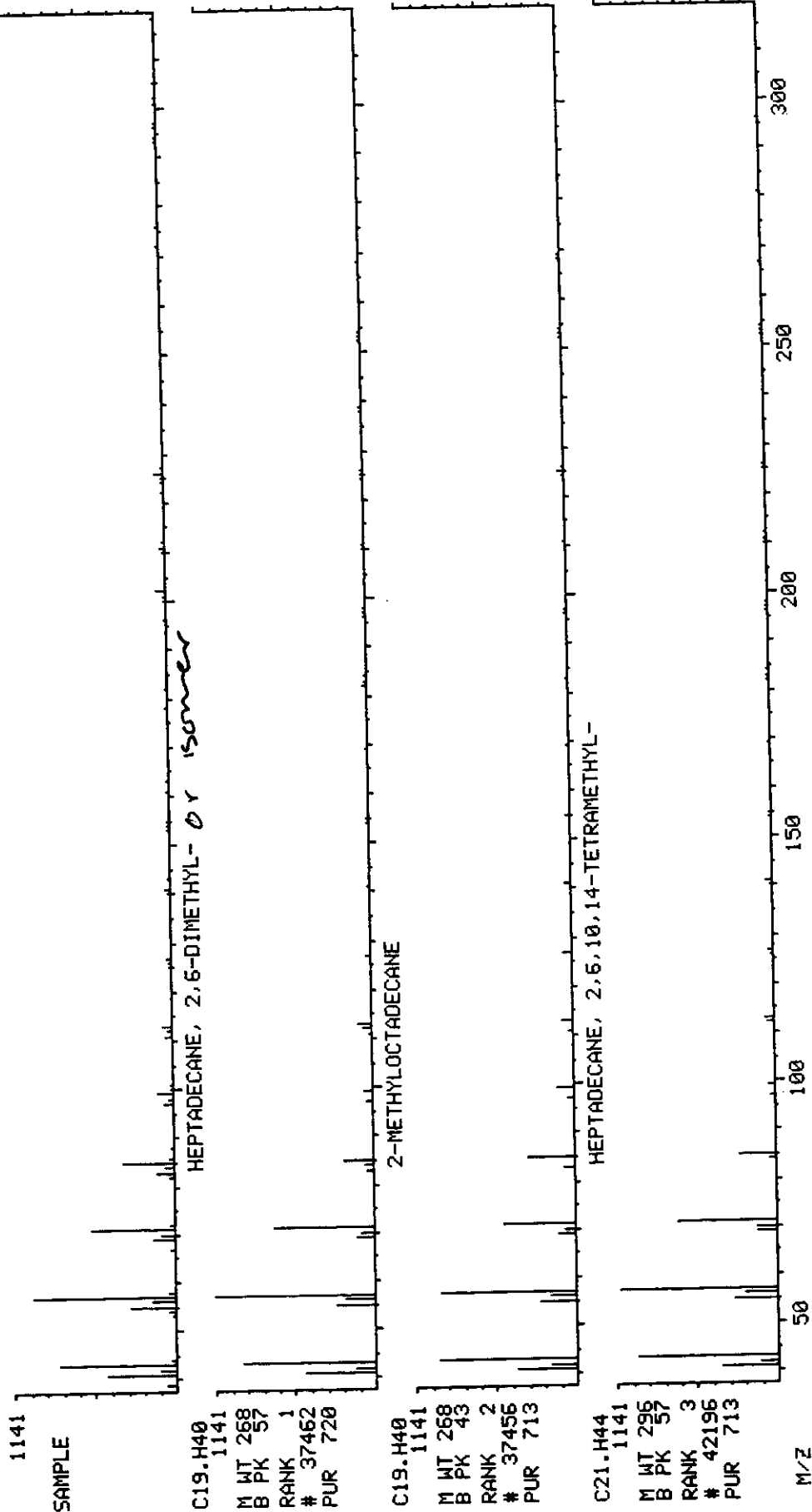
08/31/98 21:59:00 + 14:51

SAMPLE: S-MMS-3 1/35A/100M

CONDOS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M

ENHANCED (S 158 2N 0T)

INST. ID: F16



11.1
CM

Library Search Data: 30068107 #1402 Base m/z: 57
 08/31/98 21:59:00 + 15:36 Cali: 30068107 # 3 RIC: 34624.
 Sample: S-MM5-3 1/3SA/100M INST. ID: F16
 Conds.: UG/ML *100ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 714 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 37462 HEPTADECANE, 2,6-DIMETHYL-
 3 46161 TRICOSANE
 4 15969 HYDROXYLAMINE, O-DECYL-
 5 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 6 22536 TRIDECANE, 3-METHYL-
 7 32059 HEPTADECANE
 8 19523 1-DECANOL, 2-ETHYL-
 9 37456 2-METHYLOCTADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	662	953	688
2	C19.H40	268	57	644	926	667
3	C23.H48	324	43	635	902	700
4	C10.H23.O.N	173	43	633	966	635
5	C21.H44	296	57	625	932	648
6	C14.H30	198	57	624	924	629
7	C17.H36	240	57	624	929	650
8	C12.H26.O	186	57	623	948	627
9	C19.H40	268	43	621	926	657

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	638-67-5
4	—	—	—	—	29812-79-1
5	—	—	—	—	18344-37-1
6	—	—	—	—	6418-41-3
7	—	—	—	—	629-78-7
8	—	—	—	—	21078-65-9
9	—	—	—	—	- -

BASE M/Z: 57
RIC: 34624.

DATA: 30068107 #1402
CALI: 30068107 # 3

MID LIBRARY SEARCH (LIBRARYNB)

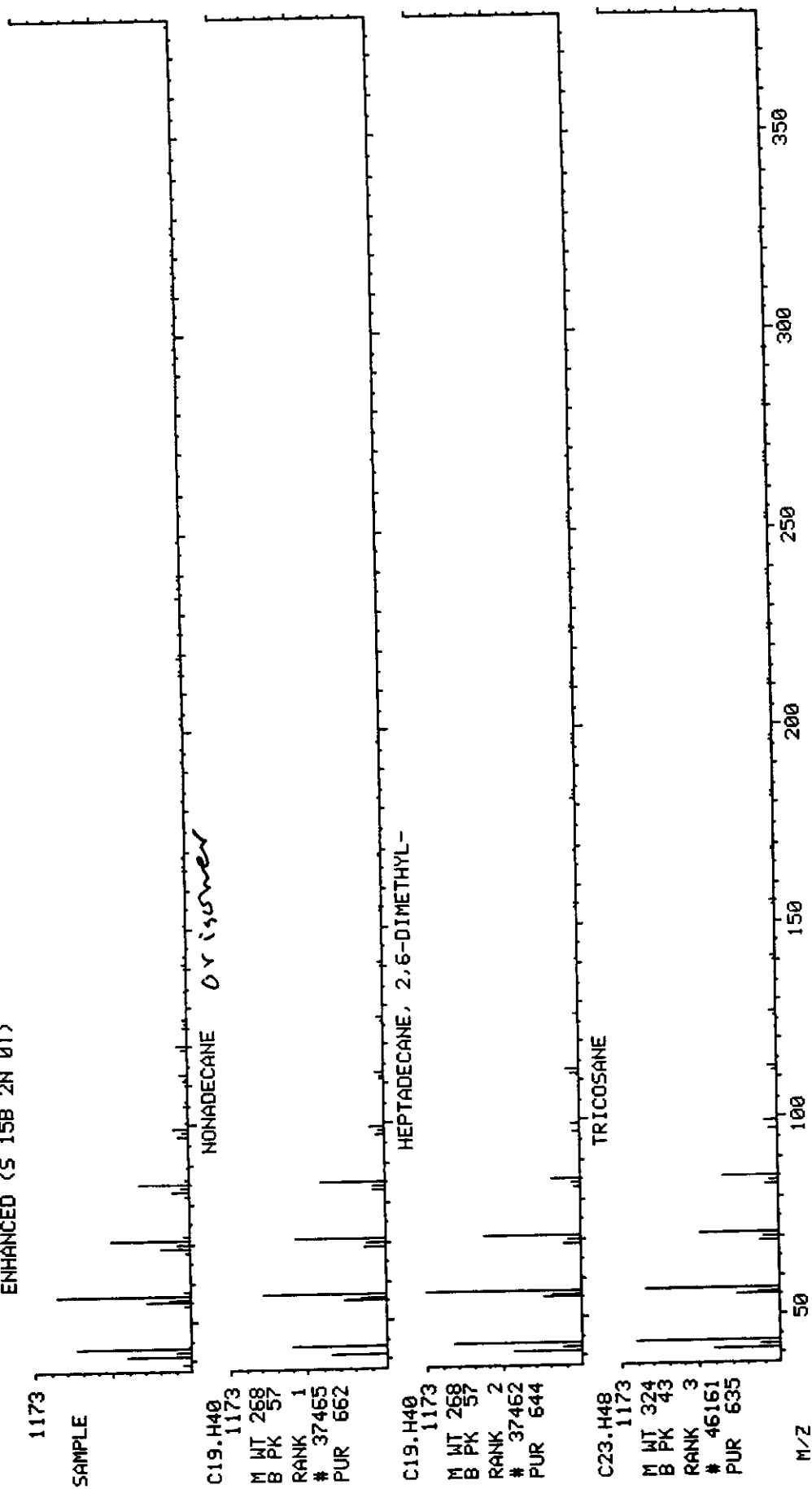
08/31/98 21:59:00 + 15:36

SAMPLE: 5-MIN5-3 1/35A/100M

INST. ID: F16

CONDS.: UG/ML *100ML *100Z/100Z *(NA/NA)/1/35A NA M

ENHANCED (5 15B 2N 0T)



1177

TIC SELECTION REPORT

DATA FILE: 30068107

THE FOLLOWING PEAKS WERE REJECTED BECAUSE
AT LEAST 40 % OF THEIR SIZE WAS ACCOUNTED FOR BY
TARGET COMPOUNDS ELUTING WITHIN 4 SCANS OF THE
PEAK TOP.

SCAN	SIZE	AMOUNT
340	178208.	21.143
530	370816.	43.994
829	337152.	40.000
1088	394368.	40.000
1552	233408.	40.000
1783	216272.	40.000

TOTAL NUMBER OF UNIDENTIFIED PEAKS WITH SIZE
GREATER THAN 10 % OF THE CLOSEST INTERNAL STANDARD
THAT DOES NOT HAVE INTERFERENCES = 103

INTERNAL STANDARDS THAT HAVE RIC SIZE LESS THAN
50 % OR GREATER THAN 200 % OF THE ESTIMATED RIC SIZE
ARE CONSIDERED TO HAVE INTERFERENCES AND WILL NOT BE USED
FOR QUANTITATION.

#	INTERNAL STANDARD	PERCENT OF ESTIMATED		SCAN
		RIC SIZE	RIC SIZE	
1	C140 NAPHTHALENE-D8	370816.	211 *	530
2	C150 ACENAPHTHENE-D10	337152.	156	829
3	C160 PHENANTHRENE-D10	394368.	163	1088
4	C170 CHRYSENE-D12	233408.	140	1552
5	C175 PERYLENE-D12	216272.	158	1783

* INDICATES INTERFERENCE

SIZE = AREA

Semivolatile Organics
Method 0010/8270

Client Name: Pacific Environmental Services
Client ID: S-MM5-FB-F, FH, XAD, COND, BH
LAB ID: 300681-0008-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 26 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
Phenol	ND	ug/Sample	30	R
bis(2-Chloroethyl)ether	ND	ug/Sample	30	
2-Chlorophenol	ND	ug/Sample	30	
1,3-Dichlorobenzene	ND	ug/Sample	30	
1,4-Dichlorobenzene	ND	ug/Sample	30	
Benzyl alcohol	ND	ug/Sample	30	
1,2-Dichlorobenzene	ND	ug/Sample	30	
2-Methylphenol	ND	ug/Sample	30	
2,2'-Oxybis(1-chloropropane)	ND	ug/Sample	30	
3/4-Methylphenol	ND	ug/Sample	30	
N-Nitroso-di-n-propylamine	ND	ug/Sample	30	
Hexachloroethane	ND	ug/Sample	30	
Nitrobenzene	ND	ug/Sample	30	
Isophorone	ND	ug/Sample	30	
2-Nitrophenol	ND	ug/Sample	30	
2,4-Dimethylphenol	ND	ug/Sample	30	
Benzoic acid	ND	ug/Sample	150	
bis(2-Chloroethoxy)-methane	ND	ug/Sample	30	
2,4-Dichlorophenol	ND	ug/Sample	30	
1,2,4-Trichlorobenzene	ND	ug/Sample	30	
Naphthalene	ND	ug/Sample	30	
4-Chloroaniline	ND	ug/Sample	30	
Hexachlorobutadiene	ND	ug/Sample	30	
4-Chloro-3-methylphenol	ND	ug/Sample	30	
2-Methylnaphthalene	ND	ug/Sample	30	
Hexachlorocyclopentadiene	ND	ug/Sample	30	
2,4,6-Trichlorophenol	ND	ug/Sample	30	
2,4,5-Trichlorophenol	ND	ug/Sample	150	
2-Chloronaphthalene	ND	ug/Sample	30	
2-Nitroaniline	ND	ug/Sample	30	
Dimethyl phthalate	ND	ug/Sample	30	
Acenaphthylene	ND	ug/Sample	30	
3-Nitroaniline	ND	ug/Sample	150	
Acenaphthene	ND	ug/Sample	30	
2,4-Dinitrophenol	ND	ug/Sample	150	
4-Nitrophenol	ND	ug/Sample	150	
Dibenzofuran	ND	ug/Sample	30	

Note R = Reporting limit(s) raised due to sample volume limitations.
ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.
Rev 230787

1178

~~1186~~
HHS

Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: S-MM5-FB-F, FH, XAD, COND, BH
LAB ID: 300681-0008-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 26 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
2,4-Dinitrotoluene	ND	ug/Sample	30	
2,6-Dinitrotoluene	ND	ug/Sample	30	
Diethyl phthalate	ND	ug/Sample	30	
4-Chlorophenyl phenyl ether	ND	ug/Sample	30	
Fluorene	ND	ug/Sample	30	
4-Nitroaniline	ND	ug/Sample	150	
4,6-Dinitro-2-methylphenol	ND	ug/Sample	150	
N-Nitrosodiphenylamine	ND	ug/Sample	30	
4-Bromophenyl phenyl ether	ND	ug/Sample	30	
Hexachlorobenzene	ND	ug/Sample	30	
Pentachlorophenol	ND	ug/Sample	150	
Phenanthrene	ND	ug/Sample	30	
Anthracene	ND	ug/Sample	30	
Di-n-butyl phthalate	ND	ug/Sample	30	
Fluoranthene	ND	ug/Sample	30	
Pyrene	ND	ug/Sample	30	
Butyl benzyl phthalate	ND	ug/Sample	30	
3,3'-Dichlorobenzidine	ND	ug/Sample	60	
Benzo(a)anthracene	ND	ug/Sample	30	
bis(2-Ethylhexyl)-phthalate	ND	ug/Sample	30	
Chrysene	ND	ug/Sample	30	
Di-n-octyl phthalate	ND	ug/Sample	30	
Benzo(b)fluoranthene	ND	ug/Sample	30	
Benzo(k)fluoranthene	ND	ug/Sample	30	
Benzo(a)pyrene	ND	ug/Sample	30	
Indeno(1,2,3-cd)pyrene	ND	ug/Sample	30	
Dibenz(a,h)anthracene	ND	ug/Sample	30	
Benzo(g,h,i)perylene	ND	ug/Sample	30	
Acetophenone	ND	ug/Sample	30	
4-Aminobiphenyl	ND	ug/Sample	150	
Aniline	ND	ug/Sample	30	
Benzidine	ND	ug/Sample	300	
3,3'-Dimethylbenzidine	ND	ug/Sample	60	
N-Nitrosodimethylamine	ND	ug/Sample	30	
N-Nitrosomorpholine	ND	ug/Sample	30	
Pentachloronitrobenzene (PCNB)	ND	ug/Sample	150	
o-Toluidine	ND	ug/Sample	60	

ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.
Rev 230787

1180

Semivolatile Organics
Method 0010/8270

(cont.)

Client Name: Pacific Environmental Services
Client ID: S-MM5-FB-F, FH, XAD, COND, BH
LAB ID: 300681-0008-SA
Matrix: AIRTRAIN
Authorized: 30 JUL 98

Sampled: 26 JUL 98
Prepared: 31 JUL 98

Received: 30 JUL 98
Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Wet wt. Result	Units	Reporting Limit	Qualifier
2-Methoxybenzenamine	ND	ug/Sample	--	K
Biphenyl	ND	ug/Sample	--	K
Chloroacetophenone	ND	ug/Sample	--	K
Cumene	ND	ug/Sample	--	K
DBCP (1,2-Dibromo-3-chloropropane)	ND	ug/Sample	--	K
Benzo(e)pyrene	ND	ug/Sample	--	K
N-N-Diethylaniline	ND	ug/Sample	--	K
Dimethylaniline	ND	ug/Sample	--	K
3,3'-Dimethoxybenzidine	ND	ug/Sample	--	K
Hydroquinone	ND	ug/Sample	--	K
4,4'-Methyl-bis(2-chloroaniline)	ND	ug/Sample	--	K
4-Nitrodiphenyl	ND	ug/Sample	--	K
Trifluralin	ND	ug/Sample	--	K

Surrogate	Recovery	Acceptable Range
Nitrobenzene-d5	58 %	45 - 107
2-Fluorobiphenyl	79 %	62 - 110
Terphenyl-d14	101 %	58 - 135
Phenol-d5	59 %	43 - 130
2-Fluorophenol	54 %	36 - 111
2,4,6-Tribromophenol	74 %	58 - 131

Note K = Identified by mass spectrum only; quantitation based on 1:1 response with internal standard.

ND = Not Detected

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.
Rev 230787

1181

Semivolatiles Library Search (20 Compound TID)
Method 8270

Client Name: Pacific Environmental Services

Client ID: S-MM5-FB-F, FH, XAD, COND, BH

LAB ID: 300681-0008-SA

Matrix: AIRTRAIN

Authorized: 30 JUL 98

Sampled: 26 JUL 98

Prepared: NA

Received: 30 JUL 98

Analyzed: 31 AUG 98

Dilution Factor: 1.0

Parameter	Result	Units	Reporting Limit	Qualifier
n-Nonane	360	ug/Sample	--	0
Benzaldehyde	29	ug/Sample	--	0
Heptadecane, 2,6-dimethyl-	52	ug/Sample	--	0
4-Hydroxy-4-methyl-2-pentanone	380	ug/Sample	--	0
Hexadecanoic acid	33	ug/Sample	--	0
Unknown alkane	51	ug/Sample	--	0
Tricosane	31	ug/Sample	--	0
Unknown alkane	30	ug/Sample	--	0
Unknown	36	ug/Sample	--	
Unknown alkane	45	ug/Sample	--	
5-Eiconsene, (E) -	150	ug/Sample	--	0
Unknown	49	ug/Sample	--	
Unknown	34	ug/Sample	--	
Unknown alkane	26	ug/Sample	--	
Unknown	57	ug/Sample	--	
Unknown	77	ug/Sample	--	
Unknown	58	ug/Sample	--	
Unknown	42	ug/Sample	--	
Unknown alkene	30	ug/Sample	--	
Unknown	27	ug/Sample	--	

Note 0 = Or structurally similar compound (isomer).
NA = Not Applicable

Reported By: Emily Uebelhoer

Approved By: Karin Yee

The cover letter is an integral part of this report.
Rev 230787

1182

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068108
Std Id: ST16980831

Sample: S-MMS-FB 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 22:29
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Surrogate Spike Recoveries
8270-G, LIMS 10/10/96

Surrogate	Surrogate Amount (ug)		% Recovery	
	Spiked	Measured	Measured	QC limits
CS20 NITROBENZENE-D5	50.00	29.00	58.0	45 107
CS25 2-FLUOROBIPHENY	50.00	39.30	78.6	62 110
CS30 TERPHENYL-D14	50.00	50.70	101.	58 135
CS45 PHENOL-D5	100.0	58.80	58.8	43 130
CS50 2-FLUOROPHENOL	100.0	53.80	53.8	36 111
CS55 2,4,6-TRIBROMOP	100.0	74.20	74.2	58 131

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting
				Limit
HEXACHLOROBENZENE-C13		ND	UG/A	30.0
C310 N-NITROSODIMETHYLAMINE		ND	UG/A	30.0
PYRIDINE		ND	UG/A	60.0
2-PICOLINE		ND	UG/A	30.0
N-NITROSOMETHYLETHYLAMINE		ND	UG/A	30.0
METHYLMETHANESULFONATE		ND	UG/A	30.0
N-NITROSODIETHYLAMINE		ND	UG/A	30.0
ETHYLMETHANE SULFONATE		ND	UG/A	30.0
PENTACHLOROETHANE		ND	UG/A	30.0
C320 ANILINE		ND	UG/A	30.0
C315 PHENOL		ND	UG/A	30.0
C325 BIS(2-CHLOROETHYL)ETHER		ND	UG/A	30.0
C330 2-CHLOROPHENOL		ND	UG/A	30.0
C335 1,3-DICHLOROBENZENE		ND	UG/A	30.0
C340 1,4-DICHLOROBENZENE		ND	UG/A	30.0
C345 BENZYL ALCOHOL		ND	UG/A	30.0
C350 1,2-DICHLOROBENZENE		ND	UG/A	30.0
C355 2-METHYLPHENOL		ND	UG/A	30.0
C360 2,2'-OXYBIS(1-CLPROPAN)		ND	UG/A	30.0
C361 ACETOPHENONE		ND	UG/A	30.0
N-NITROSPYRROLIDINE		ND	UG/A	30.0
N-NITROSOMORPHOLINE		ND	UG/A	30.0
3-METHYL PHENOL		ND	UG/A	30.0
C365 4-METHYLPHENOL		ND	UG/A	30.0
C370 N-NITROSO-DI-N-PROPYLAM		ND	UG/A	30.0
O-TOLUIDINE		ND	UG/A	30.0
C375 HEXACHLOROETHANE		ND	UG/A	30.0
C410 NITROBENZENE		ND	UG/A	30.0
N-NITROSOPIPERIDINE		ND	UG/A	30.0

Reviewed by:

149/17/96

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068108
Std Id: ST16980831

Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 22:29
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C415 ISOPHORONE		ND	UG/A	30.0
C420 2-NITROPHENOL		ND	UG/A	30.0
C425 2,4-DIMETHYLPHENOL		ND	UG/A	30.0
C435 BIS(2-CHLOROETHOXY)METH		ND	UG/A	30.0
C440 2,4-DICHLOROPHENOL		ND	UG/A	30.0
C430 BENZOIC ACID		ND	UG/A	150.0
AA-DIMETHYLPHENETHYLAMINE		ND	UG/A	30.0
C445 1,2,4-TRICHLOROBENZENE		ND	UG/A	30.0
C450 NAPHTHALENE		ND	UG/A	30.0
C455 4-CHLOROANILINE		ND	UG/A	30.0
2,6-DICHLOROPHENOL		ND	UG/A	30.0
HEXACHLOROPROPENE		ND	UG/A	30.0
C460 HEXACHLOROBUTADIENE		ND	UG/A	30.0
P-PHENYLENE DIAMINE		ND	UG/A	30.0
N-NITROSODI-N-BUTYLAMINE		ND	UG/A	30.0
C465 4-CHLORO-3-METHYLPHENO		ND	UG/A	30.0
SAFROLE		ND	UG/A	30.0
C470 2-METHYLNAPHTHALENE		ND	UG/A	30.0
1,2,4,5-TETRACHLOROBENZENE		ND	UG/A	30.0
ISOSAFROLE (#1)		ND	UG/A	60.0
C510 HEXACHLOROCYCLOPENTADI		ND	UG/A	30.0
C515 2,4,6-TRICHLOROPHENOL		ND	UG/A	30.0
C520 2,4,5-TRICHLOROPHENOL		ND	UG/A	30.0
ISOSAFROLE (#2)		ND	UG/A	60.0
C525 2-CHLORONAPHTHALENE		ND	UG/A	30.0
1-CHLORONAPHTHALENE		ND	UG/A	15.0
C530 2-NITROANILINE		ND	UG/A	150.0
1,4-NAPHTHOQUINONE		ND	UG/A	30.0
C535 DIMETHYLPHTHALATE		ND	UG/A	30.0
1,3-DINITROBENZENE		ND	UG/A	30.0
C540 ACENAPHTHYLENE		ND	UG/A	30.0
C543 2,6-DINITROTOLUENE		ND	UG/A	30.0
C545 3-NITROANILINE		ND	UG/A	150.0
C550 ACENAPHTHENE		ND	UG/A	30.0
C555 2,4-DINITROPHENOL		ND	UG/A	150.0
C565 DIBENZOFURAN		ND	UG/A	30.0
C560 4-NITROPHENOL		ND	UG/A	150.0
PENTACHLOROBENZENE		ND	UG/A	30.0
C570 2,4-DINITROTOLUENE		ND	UG/A	30.0
1-NAPHTHYLAMINE		ND	UG/A	30.0
2-NAPHTHYLAMINE		ND	UG/A	30.0
2,3,4,6-TETRACHLOROPHENOL		ND	UG/A	60.0
C580 DIETHYLPHTHALATE		ND	UG/A	30.0
C590 FLUORENE		ND	UG/A	30.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

Data File: 30068108
Std Id: ST16980831

Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 22:29
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
=====				
C585 4-CHLOROPHENYL-PHENYLE		ND	UG/A	30.0
5-NITRO-O-TOLUIDINE		ND	UG/A	30.0
C595 4-NITROANALINE		ND	UG/A	150.0
C610 4,6-DINITRO-2-METHYLPH		ND	UG/A	150.0
C615 N-NITROSODIPHENYLAMINE		ND	UG/A	30.0
C620 AZOBENZENE		ND	UG/A	30.0
SYM-TRINITROBENZENE		ND	UG/A	30.0
C625 4-BROMOPHENYL-PHENYLET		ND	UG/A	30.0
PHENACETIN		ND	UG/A	30.0
DIALATE	234	ND	UG/A	30.0
C630 HEXACHLOROBENZENE		ND	UG/A	30.0
4-AMINOBIIPHENYL		ND	UG/A	30.0
C635 PENTACHLOROPHENOL		ND	UG/A	150.0
PRONAMIDE		ND	UG/A	30.0
PENTACHLORONITROBENZENE		ND	UG/A	150.0
C640 PHENANTHRENE		ND	UG/A	30.0
C645 ANTHRACENE		ND	UG/A	30.0
2SECBUTYL-4,6-DINITROPHENOL		ND	UG/A	30.0
C647 CARBAZOLE		ND	UG/A	30.0
C650 DI-N-BUTYLPHTHALATE		ND	UG/A	30.0
4-NITROQUINOLINE-1-OXIDE		ND	UG/A	30.0
METHAPYRILENE		ND	UG/A	30.0
ISODRIN		ND	UG/A	30.0
C655 FLUORANTHENE		ND	UG/A	30.0
CHLOROBENZILATE		ND	UG/A	30.0
C710 BENZIDINE		ND	UG/A	300.0
C715 PYRENE		ND	UG/A	30.0
ARAMITE (#1)		ND	UG/A	30.0
ARAMITE (#2)		ND	UG/A	30.0
P-DIMETHYLAMINOAZOBENZENE		ND	UG/A	30.0
3,3'-DIMETHYLBENZIDINE		ND	UG/A	30.0
KEPONE		ND	UG/A	150.0
C720 BUTYLBENZYLPHTHALATE		ND	UG/A	30.0
2-ACETYLAMINOFLUORENE		ND	UG/A	30.0
C730 BENZO(A)ANTHRACENE		ND	UG/A	30.0
C725 3,3'-DICHLOROBENZIDINE		ND	UG/A	60.0
C740 CHRYSENE		ND	UG/A	30.0
C745 BIS(2-ETHYLHEXYL)PHTHA		ND	UG/A	30.0
3-METHYLCHOLANTHRENE		ND	UG/A	30.0
C760 DI-N-OCTYL PHTHALATE		ND	UG/A	30.0
C765 BENZO(B)FLUORANTHENE		ND	UG/A	30.0
7,12-DIMETHYLBENZANTHRACENE		ND	UG/A	30.0
C770 BENZO(K)FLUORANTHENE		ND	UG/A	30.0
HEXACHLOROPHENE		ND	UG/A	30.0

QUANTERRA GC/MS
Target Compound Data Summary Sheet

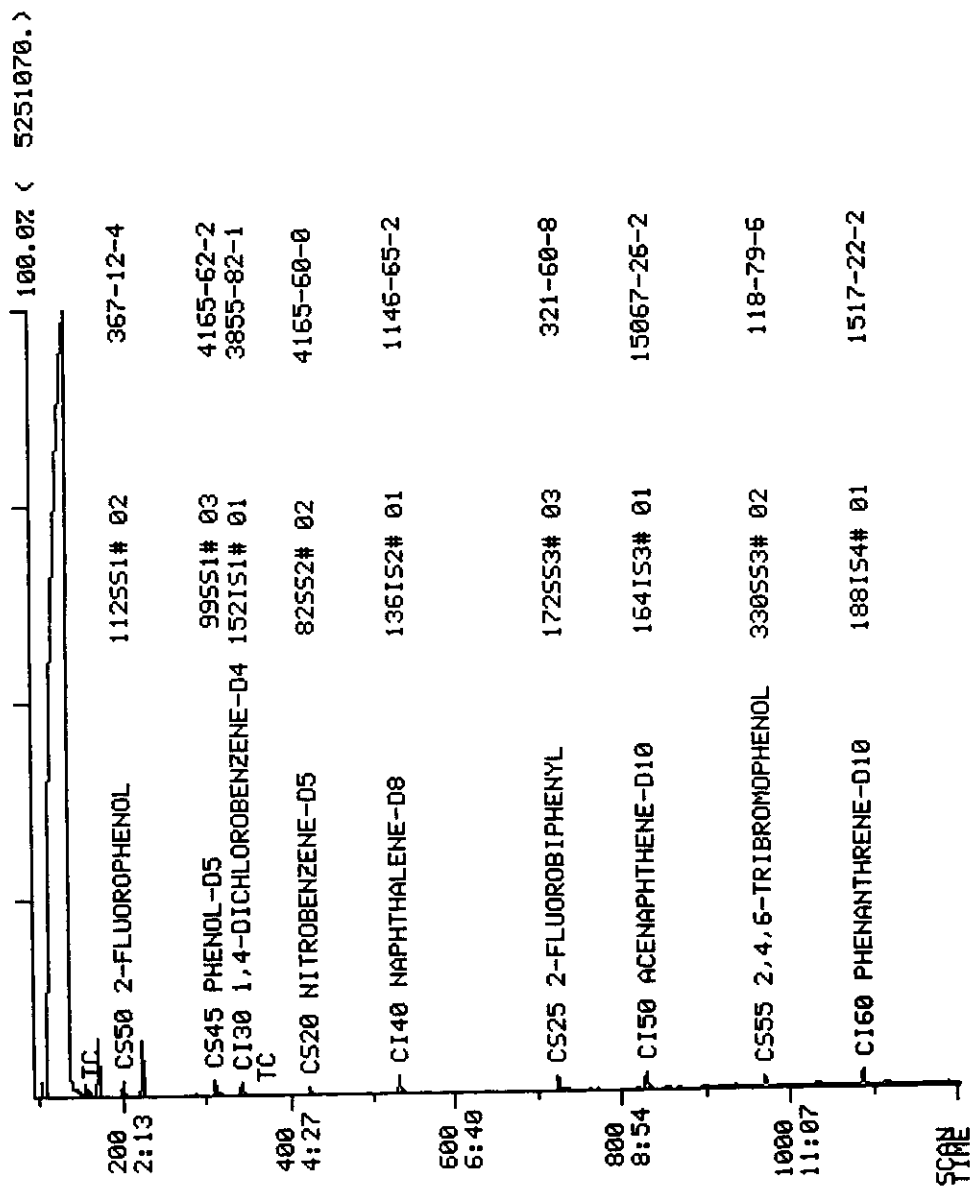
Data File: 30068108
Std Id: ST16980831

Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
Client: PACIFI Date Analyzed: 08/31/98 22:29
Analyst: DAT Instrument ID: F16 Run Factor: 3.00
Quan List Threshold: 0.95 Surrogate Vol.: 1.00

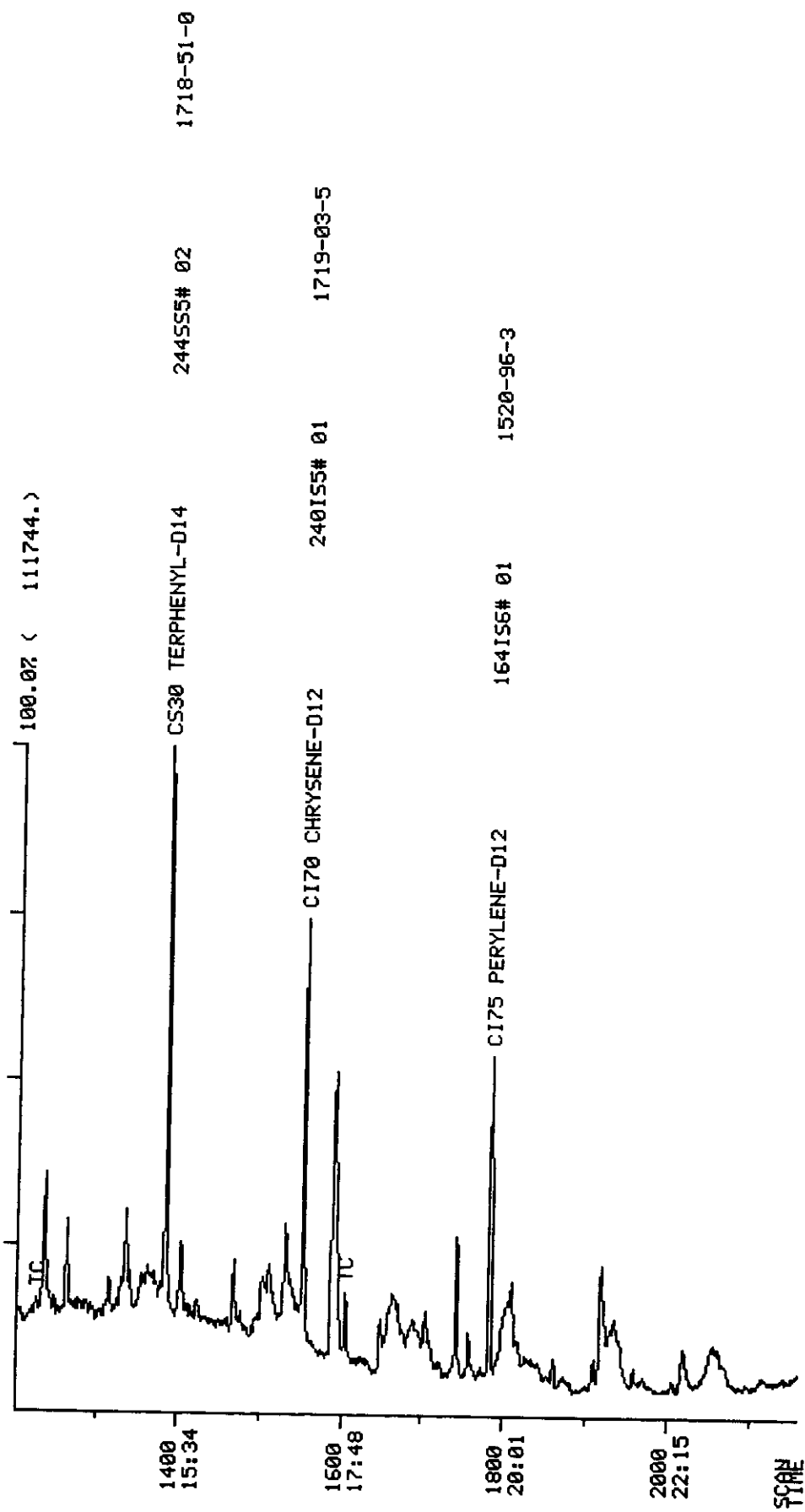
Target Compounds: SAP9

Parameter	Scan	Result	Units	Reporting Limit
C775 BENZO(A)PYRENE		ND	UG/A	30.0
C780 INDENO(1,2,3-CD)PYRENE		ND	UG/A	30.0
C785 DIBENZ(A,H)ANTHRACENE		ND	UG/A	30.0
C790 BENZO(G,H,I)PERYLENE		ND	UG/A	30.0

DATA FROM FILE: 30068108 SCANS 93 TO 1203 ACQUIRED: 08/31/98 22:29:00
 SAMPLE: S-MMS-FB 1/35A/1ML CALI: 30068108 #3
 COND.: UG/ML *1002/1002 *(NA/NA) 1/35A NA M INST. ID: F16



DATA FROM FILE: 30058108 SCANS 1203 TO 2158 ACQUIRED: 08/31/98 22:29:00
CALI: 30058108 #3
SAMPLE: S-MMS-F8 1/35A/1ML INST. ID: F16
CONDS.: UG/ML *100%/100% *(NA/NA >1/35A NA M



ADDITIONAL STANDARD USED FOR THIS REPORT IS: ST16980831A

QUANTERRA QUANTITATION SUMMARY

File: 30068108

Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
Analyst: DAT Instrument Id: F16 Analyzed: 08/31/98 22:29

Compounds with amounts less than 0.20 reported as NOT FOUND

Library		Units: UG/ML							
No	Entry	Name	Mass	Meth	Scan	Ref	Area	RRF(L)	Amount
1	S1#	1 C130 1,4-DICHLOROBENZENE	152	A	BB	340	1	36214. 1.000	40.000
2	S2#	1 C140 NAPHTHALENE-DB	136	A	BB	529	2	139245. 1.000	40.000
3	S3#	1 C150 ACENAPHTHENE-D10	164	A	BB	827	3	74547. 1.000	40.000
4	S4#	1 C160 PHENANTHRENE-D10	188	A	BB	1086	4	121683. 1.000	40.000
5	S5#	1 C170 CHRYSENE-D12	240	A	BB	1551	5	65622. 1.000	40.000
6	S6#	1 C175 PERYLENE-D12	264	A	BV	1782	6	53016. 1.000	40.000
7	S2#	2 CS20 NITROBENZENE-D5	82	A	BB	421	2	47363. 0.470	28.975
8	S3#	3 CS25 2-FLUOROBIPHENYL	172	A	BB	722	3	90468. 1.234	39.328
9	S5#	2 CS30 TERPHENYL-D14	244	A	BB	1383	5	82729. 0.995	50.670
10	S1#	3 CS45 PHENOL-D5	99	A	BB	308	1	107781. 2.025	58.774
11	S1#	2 CS50 2-FLUOROPHENOL	112	A	BB	198	1	66619. 1.368	53.779
12	S3#	2 CS55 2,4,6-TRIBROMOPHENO	330	A	BB	968	3	32579. 0.236	74.180
13	S4#	4 HEXACHLOROBENZENE-C13	294			4	NOT FOUND		
14	S1#	15 C310 N-NITROSODIMETHYLAM	74			1	NOT FOUND		
15	S1#	20 PYRIDINE	79			1	NOT FOUND		
16	S1#	30 2-PICOLINE	93			1	NOT FOUND		
17	S1#	40 N-NITROSOMETHYLETHYLAMIN	42	A	VV	159	1	5080. 0.811	6.918
18	S1#	60 METHYLMETHANESULFONATE	80			1	NOT FOUND		
19	S1#	70 N-NITROSODIETHYLAMINE	102			1	NOT FOUND		
20	S1#	85 ETHYLMETHANE SULFONATE	79			1	NOT FOUND		
21	S1#	95 PENTACHLOROETHANE	117			1	NOT FOUND		
22	S1#	100 C320 ANILINE	93			1	NOT FOUND		
23	S1#	105 C315 PHENOL	94	A	BB	309	1	604. 1.968	0.339
24	S1#	110 C325 BIS(2-CHLOROETHYL)E	93			1	NOT FOUND		
25	S1#	115 C330 2-CHLOROPHENOL	128			1	NOT FOUND		
26	S1#	125 C335 1,3-DICHLOROBENZENE	146			1	NOT FOUND		
27	S1#	130 C340 1,4-DICHLOROBENZENE	146			1	NOT FOUND		
28	S1#	145 C345 BENZYL ALCOHOL	108	A	BB	370	1	2842. 0.854	3.674
29	S1#	150 C350 1,2-DICHLOROBENZENE	146			1	NOT FOUND		
30	S1#	160 C355 2-METHYLPHENOL	108			1	NOT FOUND		
31	S1#	165 C360 2,2'-OXYBIS(1-CLPRO	45			1	NOT FOUND		
32	S1#	170 C361 ACETOPHENONE	105			1	NOT FOUND		
33	S1#	175 N-NITROSOPYRROLIDINE	100			1	NOT FOUND		
34	S1#	180 N-NITROSOMORPHOLINE	56			1	NOT FOUND		
35	S1#	182 3-METHYL PHENOL	108			1	NOT FOUND		
36	S1#	185 C365 4-METHYLPHENOL	108			1	NOT FOUND		
37	S1#	190 C370 N-NITROSO-DI-N-PROP	70			1	NOT FOUND		
38	S1#	195 O-TOLUIDINE	106			1	NOT FOUND		
39	S1#	200 C375 HEXACHLOROETHANE	117			1	NOT FOUND		
40	S2#	10 C410 NITROBENZENE	77			2	NOT FOUND		
41	S2#	15 N-NITROSOPIPERIDINE	42			2	NOT FOUND		
42	S2#	20 C415 ISOPHORONE	82			2	NOT FOUND		
43	S2#	25 C420 2-NITROPHENOL	139			2	NOT FOUND		
44	S2#	30 C425 2,4-DIMETHYLPHENOL	107			2	NOT FOUND		
45	S2#	35 C435 BIS(2-CHLOROETHOXY)	93			2	NOT FOUND		

46	S2# 40	C440 2,4-DICHLOROPHENOL	162	2	NOT FOUND
47	S2# 45	C430 BENZOIC ACID	122	2	NOT FOUND
48	S2# 50	AA-DIMETHYLPHENETHYLAMIN	58	2	NOT FOUND
49	S2# 55	C445 1,2,4-TRICHLOROBENZ	180	2	NOT FOUND

50	S2# 60	C450 NAPHTHALENE	128	2	NOT FOUND			
51	S2# 80	C455 4-CHLOROANILINE	127	2	NOT FOUND			
52	S2# 85	2,6-DICHLOROPHENOL	162	2	NOT FOUND			
53	S2# 90	HEXACHLOROPROPENE	213	2	NOT FOUND			
54	S2# 95	C460 HEXACHLOROBUTADIENE	225	2	NOT FOUND			
55	S2#115	P-PHENYLENE DIAMINE	108	2	NOT FOUND			
56	S2#120	N-NITROSODI-N-BUTYLAMINE	84	2	NOT FOUND			
57	S2#130	C465 4-CHLORO-3-METHYLPH	107	2	NOT FOUND			
58	S2#140	SAFROLE	162	2	NOT FOUND			
59	S2#145	C470 2-METHYLNAPHTHALENE	142	2	NOT FOUND			
60	S3# 10	1,2,4,5-TETRACHLOROBENZE	214	3	NOT FOUND			
61	S3# 15	ISOSAFROLE (#1)	162	3	NOT FOUND			
62	S3# 20	C510 HEXACHLOROCYCLOPENT	237	3	NOT FOUND			
63	S3# 25	C515 2,4,6-TRICHLOROPHEN	196	3	NOT FOUND			
64	S3# 30	C520 2,4,5-TRICHLOROPHEN	196	3	NOT FOUND			
65	S3# 35	ISOSAFROLE (#2)	104	3	NOT FOUND			
66	S3# 40	C525 2-CHLORONAPHTHALENE	162	3	NOT FOUND			
67	S3# 42	1-CHLORONAPHTHALENE	162	3	NOT FOUND			
68	S3# 45	C530 2-NITROANILINE	65 A VB	761	3	298. 0.502	0.319	
69	S3# 50	1,4-NAPHTHOQUINONE	158	3	NOT FOUND			
70	S3# 55	C535 DIMETHYLPHTHALATE	163	3	NOT FOUND			
71	S3# 60	1,3-DINITROBENZENE	168	3	NOT FOUND			
72	S3# 65	C540 ACENAPHTHYLENE	152	3	NOT FOUND			
73	S3# 70	C543 2,6-DINITROTOLUENE	165 A BB	814	3	460. 0.324	0.762	
74	S3# 75	C545 3-NITROANILINE	138	3	NOT FOUND			
75	S3# 80	C550 ACENAPHTHENE	153	3	NOT FOUND			
76	S3# 85	C555 2,4-DINITROPHENOL	184	3	NOT FOUND			
77	S3# 90	C565 DIBENZOFURAN	168	3	NOT FOUND			
78	S3# 95	C560 4-NITROPHENOL	109	3	NOT FOUND			
79	S3#100	PENTACHLOROBENZENE	250	3	NOT FOUND			
80	S3#105	C570 2,4-DINITROTOLUENE	165 A BB	879	3	222. 0.397	0.300	
81	S3#110	1-NAPHTHYLAMINE	143	3	NOT FOUND			
82	S3#115	2-NAPHTHYLAMINE	143	3	NOT FOUND			
83	S3#120	2,3,4,6-TETRACHLOROPHENO	232	3	NOT FOUND			
84	S3#130	C580 DIETHYLPHTHALATE	149 A VB	931	3	1395. 1.369	0.547	
85	S3#135	C590 FLUORENE	166	3	NOT FOUND			
86	S3#140	C585 4-CHLOROPHENYL-PHEN	204	3	NOT FOUND			
87	S3#145	5-NITRO-O-TOLUIDINE	152 A VB	939	3	208. 0.356	0.314	
88	S3#150	C595 4-NITROANILINE	138 A BB	948	3	118. 0.307	0.206	
89	S4# 10	C610 4,6-DINITRO-2-METHY	198	4	NOT FOUND			
90	S4# 15	C615 N-NITROSODIPHENYLAM	169	4	NOT FOUND			
91	S4# 20	C620 AZOBENZENE	77 A BB	966	4	1071. 1.171	0.301	
92	S4# 25	SYM-TRINITROBENZENE	75	4	NOT FOUND			
93	S4# 30	C625 4-BROMOPHENYL-PHENY	248	4	NOT FOUND			
94	S4# 35	PHENACETIN	108 A VV	1035	4	320. 0.317	0.332	
95	S4# 37	DIALATE	234	4	NOT FOUND			
96	S4# 40	C630 HEXACHLOROBENZENE	284	4	NOT FOUND			
97	S4# 45	4-AMINOBI-PHENYL	169	4	NOT FOUND			
98	S4# 50	C635 PENTACHLOROPHENOL	266	4	NOT FOUND			
99	S4# 55	PRONAMIDE	173	4	NOT FOUND			
100	S4# 60	PENTACHLORONITROBENZENE	237	4	NOT FOUND			
101	S4# 65	C640 PHENANTHRENE	178 A BB	1089	4	1753. 1.033	0.558	
102	S4# 70	C645 ANTHRACENE	178	4	NOT FOUND			
103	S4# 75	2SECBUTYL-4,6-DINITROPHE	211	4	NOT FOUND			
104	S4# 80	C647 CARBAZOLE	167	4	NOT FOUND			
105	S4# 85	C650 DI-N-BUTYLPHTHALATE	149 A BB	1226	4	4322. 1.247	1.140	
106	S4#100	4-NITROQUINOLINE-1-OXIDE	190 A BB	1237	4	272. 0.073	1.233	
107	S4#105	METHAPYRILENE	58 A BB	1263	4	803. 0.305	0.867	
108	S4#106	ISODIRIN	193	4	NOT FOUND			
109	S4#110	C655 FLUORANTHENE	202	4	NOT FOUND			
110	S4#120	CHLOROBENZILATE	139	4	NOT FOUND			

111	S5# 10	C710 BENZIDINE	184	5	NOT FOUND		
112	S5# 15	C715 PYRENE	202	5	NOT FOUND		
113	S5# 20	ARAMITE (#1)	185	5	NOT FOUND		
114	S5# 25	ARAMITE (#2)	185	5	NOT FOUND		
115	S5# 30	P-DIMETHYLAMINOAZOBENZEN	120	5	NOT FOUND		
116	S5# 35	3,3'-DIMETHYLBENZIDINE	212	5	NOT FOUND		
117	S5# 37	KEPONE	272	5	NOT FOUND		
118	S5# 40	C720 BUTYLBENZYLPHTHALAT	149	5	NOT FOUND		
119	S5# 45	2-ACETYLAMINOFLUORENE	181	A VV 1516	5	330.	0.432 0.465
120	S5# 50	C730 BENZO(A)ANTHRACENE	228	A BB 1551	5	574.	1.125 0.311
121	S5# 55	C725 3,3'-DICHLOROBENZID	252	5	NOT FOUND		
122	S5# 60	C740 CHRYSENE	228	5	NOT FOUND		
123	S5# 65	C745 BIS(2-ETHYLHEXYL)PH	149	A BB 1603	5	9579.	0.980 5.959
124	S5# 85	3-METHYLCHOLANTHRENE	268	5	NOT FOUND		
125	S6# 10	C760 DI-N-OCTYL PHTHALAT	149	6	NOT FOUND		
126	S6# 15	C765 BENZO(B)FLUORANTHEN	252	6	NOT FOUND		
127	S6# 20	7,12-DIMETHYLBENZANTHRAC	256	6	NOT FOUND		
128	S6# 25	C770 BENZO(K)FLUORANTHEN	252	6	NOT FOUND		
129	S6# 30	HEXACHLOROPHENE	196	6	NOT FOUND		
130	S6# 35	C775 BENZO(A)PYRENE	252	6	NOT FOUND		
131	S6# 55	C780 INDENO(1,2,3-CD)PYR	276	6	NOT FOUND		
132	S6# 60	C785 DIBENZ(A,H)ANTHRACE	278	6	NOT FOUND		
133	S6# 65	C790 BENZO(G,H,I)PERYLEN	276	6	NOT FOUND		

QUANTERRA Internal Standard Check

Standard Filename: ST16980831

Analyzed: 08/31/98 12:31

Standard	IS# 1			IS# 2			IS# 3		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	28649.		3.78	100432.		5.90	54347.		9.21
Upper Limit	57297.=+200		4.28	200864.=+200		6.40	108695.=+200		9.71
Lower Limit	14324.=+ 50		3.28	50216.=+ 50		5.40	27174.=+ 50		8.71
Filename									
1 30068108	36214.= 126		3.78	139245.= 138		5.88	74547.= 137		9.20

IS# 1 = C130 1,4-DICHLOROBENZENE-D4

IS# 2 = C140 NAPHTHALENE-D8

IS# 3 = C150 ACENAPHTHENE-D10

* - indicates an I.S. is outside QC limit(s)

Standard	IS# 4			IS# 5			IS# 6		
	Area	%	RT	Area	%	RT	Area	%	RT
ST16980831	98520.		12.08	62289.		17.26	50750.		19.83
Upper Limit	197039.=+200		12.58	124578.=+200		17.76	101500.=+200		20.33
Lower Limit	49260.=+ 50		11.58	31145.=+ 50		16.76	25375.=+ 50		19.33
Filename									
1 30068108	121683.= 123		12.08	65622.= 105		17.25	53016.= 104		19.82

IS# 4 = C160 PHENANTHRENE-D10

IS# 5 = C170 CHRYSENE-D12

IS# 6 = C175 PERYLENE-D12

* - indicates an I.S. is outside QC limit(s)

Data Reduced by: 2 Date: 9/17/98

Data File: 30068108

Data Reviewed by: 1 Date: 9/17/98

QUANTERRA GC/MS TIC REPORT (Part 1)

Sample: S-MM5-FB 1/3SA/1ML

INST. ID: F16

Analyst: DAT

Date Analyzed: 08/31/98 22:29

Run Factor: 3.00

Concentration

in Sample

SCAN (UG/A) CAS #
170 380. 123-42-2
2-PENTANONE, 4-HYDROXY-4-METHYL-

2 222 360. 111-84-2
NONANE *or isomer*

3 285 29. 100-52-7
BENZALDEHYDE

X 932 22. ~~74381-40-1~~
~~PROPANOIC ACID, 2-METHYL-, 1-(1,1-DIMETHYLETHYL)-2-METHYL-1,5-PROPANEDI~~

Unknown 5700

X 985 20. 629-92-5
NONADECANE *or isomer*

6 1033 52. 54105-67-8
HEPTADECANE, 2,6-DIMETHYL

7 1116 24. 629-92-5
NONADECANE

8 1235 33. 57-10-3
HEXADECANOIC ACID

9 1334 51. ~~54105-67-8~~
~~HEPTADECANE, 2,6-DIMETHYL~~

*alkane**5746400*

X 1361 20. 00-00-0
UNKNOWN

5700

11 1401 31. 638-67-5
TRICOSANE or isomer

12 1467 30. ~~638-67-5~~
TRICOSANE alkane 5746400

~~13~~ 1501 24. 00-00-0 5700
UNKNOWN

14 1509 36. 00-00-0
UNKNOWN

15 1529 45. ~~629-92-5~~
~~NONADECANE~~ alkane 5740400

16 1588 150. 74685-30-6
5-EICOSENE, (E)- or isomer

17 1661 49. 00-00-0 5700
UNKNOWN

18 1688 34. 00-00-0
UNKNOWN

19 1703 26. ~~629-92-5~~
~~NONADECANE~~ alkane 5746400

20 1741 57. 00-00-0 5700
UNKNOWN

21 1809 77. 00-00-0
UNKNOWN

22 1917 58. ~~655-53-4~~
4-TETRADECANOL unknown 5700

23 1934
UNKNOWN

42.

00-00-0

5700

24 2016

30.

7208-19-1

3-OCTADECENE, (E)-

alkene

5720000

25 2052

27.

00-00-0

UNKNOWN

5700

DATA FILE: 30068108

QUANTERRA GC/MS TIC REPORT (Part 2)

CONCENTRATION = AREA(TIC)*CONC(1S)/AREA(1S)

#	FIT	PURITY	INT.		RT	RRT	AREA	HEIGHT	AMOUNT		LIB	LIB #
			STD.						AS ANALYZED			
1	955	916	1	1:53	0.321	906380.	388608.	126.768	NB	3241.		
2	986	935	1	2:28	0.420	859436.	372736.	120.203	NB	5159.		
3	967	923	1	3:10	0.539	69204.	17724.	9.679	NB	2017.		
4	915	728	2	10:21	1.127	63056.	21101.	7.252	NB	40501.		
5	926	763	3	10:57	0.907	58112.	13152.	6.546	NB	37465.		
6	939	746	3	11:29	0.951	152832.	27072.	17.215	NB	37462.		
7	955	779	3	12:24	1.028	72432.	21408.	8.159	NB	37465.		
8	977	722	3	13:44	1.137	97728.	22656.	11.008	NB	35182.		
9	892	621	4	14:50	0.860	92576.	15968.	17.096	NB	37462.		
10	664	194	4	15:08	0.877	36800.	6272.	6.796	UK	1.		
11	903	624	4	15:34	0.903	56608.	12576.	10.454	NB	46161.		
12	900	695	4	16:19	0.946	54816.	11600.	10.123	NB	46161.		
13	696	200	4	16:41	0.968	43920.	7152.	8.111	UK	1.		
14	816	451	4	16:47	0.973	64912.	9296.	11.987	UK	1.		
15	955	643	4	17:00	0.986	80832.	15072.	14.927	NB	37465.		
16	975	742	4	17:39	1.024	263344.	45632.	48.631	NB	39516.		
17	800	345	4	18:28	1.071	88976.	7792.	16.431	UK	1.		
18	687	232	5	18:46	0.947	49792.	4368.	11.377	UK	1.		
19	941	723	5	18:56	0.956	37936.	6944.	8.668	NB	37465.		
20	895	502	5	19:21	0.977	83508.	23117.	19.081	UK	1.		
21	904	537	5	20:07	1.015	112176.	10848.	25.631	UK	1.		
22	966	729	5	21:19	1.076	85264.	16481.	19.482	NB	26416.		
23	722	399	5	21:30	1.085	61840.	5968.	14.130	UK	1.		
24	927	696	5	22:25	1.131	44000.	6658.	10.053	NB	34410.		
25	839	412	5	22:49	1.152	39200.	3904.	8.957	UK	1.		

Library Search Data: 30068108 # 170 Base m/z: 43
 08/31/98 22:29:00 + 1:53 Cali: 30068108 # 3 RIC: 278528.
 Sample: S-MMS-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 630 matched at least 6 of the 16 largest peaks in the unknown

Rank In.	Name
1	3241 2-PENTANONE, 4-HYDROXY-4-METHYL-
2	5791 1,3-DIOXOLANE-2-METHANOL, 2,4-DIMETHYL-
3	3366 2-PENTANOL, 2,4-DIMETHYL-
4	3288 ACETIC ACID, 1,1-DIMETHYLETHYL ESTER
5	5594 2-PROPANOL, 2-NITROSO-, ACETATE (ESTER)
6	919 TERT-BUTYL HYDROPEROXIDE
7	3349 2-HEXANOL, 2-METHYL-
8	8432 2-HEXANONE, 3-HYDROXY-3,5-DIMETHYL-
9	8425 2-HEPTANONE, 3-HYDROXY-3-METHYL-

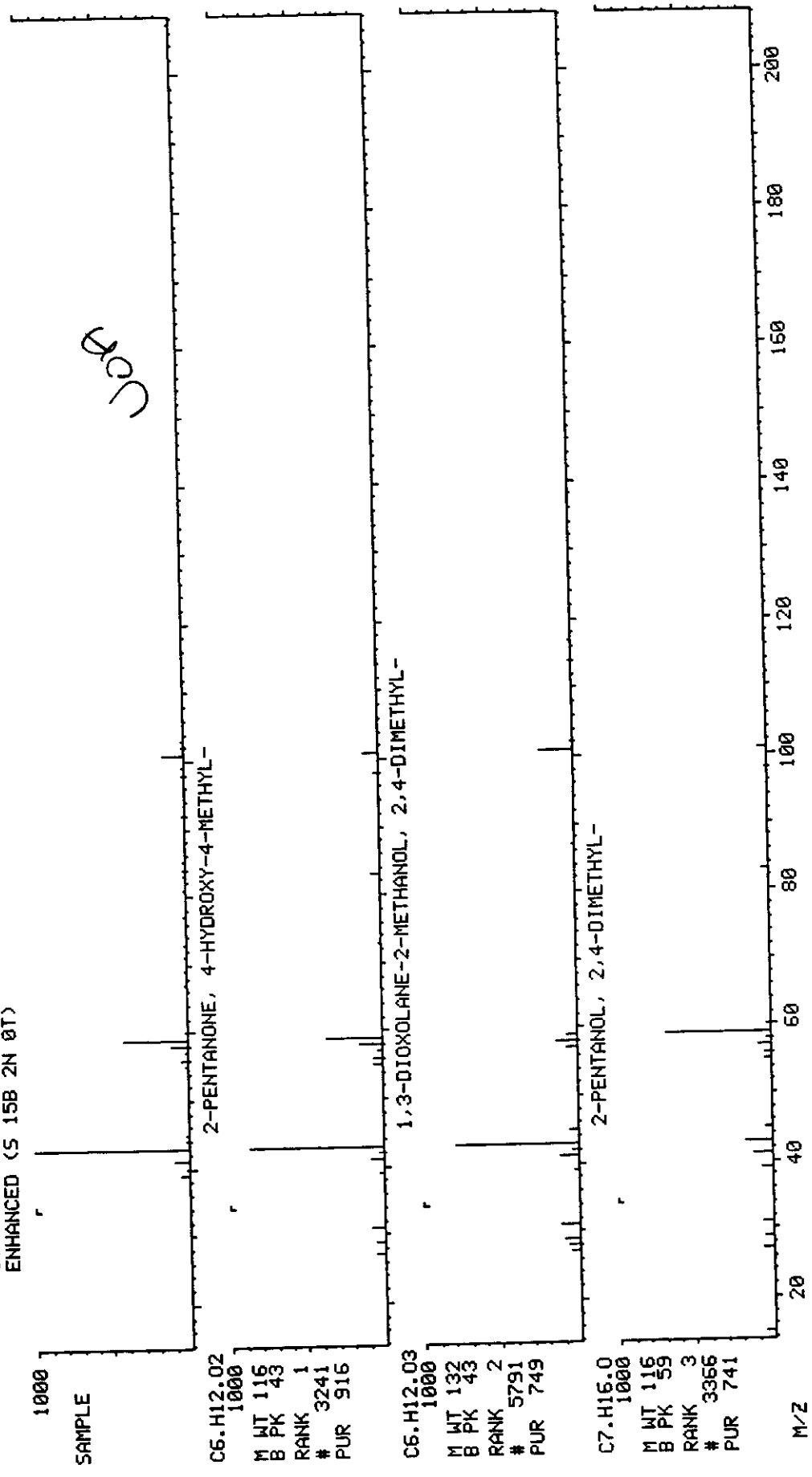
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C6.H12.O2	116	43	916	955	929
2	C6.H12.O3	132	43	749	823	793
3	C7.H16.O	116	59	741	764	741
4	C6.H12.O2	116	43	725	774	739
5	C5.H9.O3.N	131	43	718	821	810
6	C4.H10.O2	90	59	705	856	727
7	C7.H16.O	116	59	704	738	704
8	C8.H16.O2	144	59	662	725	688
9	C8.H16.O2	144	59	659	715	675

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	123-42-2
2	—	—	—	—	53951-43-2
3	—	—	—	—	625-06-9
4	—	—	—	—	540-88-5
5	—	—	—	—	6931-04-0
6	—	—	—	—	75-91-2
7	—	—	—	—	625-23-0
8	—	—	—	—	6321-14-8
9	—	—	—	—	13757-91-0

DATA: 30058108 # 170
CALI: 30058108 # 3

BASE M/Z: 43
RIC: 278528.

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 22:29:00 + 1:53
SAMPLE: 5-MM5-FB 1/35A/1ML INST. ID: F16
COND5: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)



Library Search Data: 30068108 # 222 Base m/z: 43
 08/31/98 22:29:00 + 2:28 Cali: 30068108 # 3 RIC: 253952.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 921 matched at least 7 of the 16 largest peaks in the unknown

Rank	In.	Name
1	5159	NONANE
2	5151	HEPTANE, 3,4-DIMETHYL-
3	3086	HEXANE, 2,4-DIMETHYL-
4	5141	HEPTANE, 2,4-DIMETHYL-
5	19015	DECANE, 2,5,6-TRIMETHYL-
6	8085	HEPTANE, 2,4,6-TRIMETHYL-
7	5144	HEPTANE, 4-ETHYL-
8	11607	UNDECANE
9	3081	OCTANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C9.H20	128	43	935	986	935
2	C9.H20	128	43	897	926	897
3	C8.H18	114	43	887	963	898
4	C9.H20	128	43	885	938	890
5	C13.H28	184	57	868	910	887
6	C10.H22	142	43	864	909	879
7	C9.H20	128	43	863	914	871
8	C11.H24	156	43	859	921	908
9	C8.H18	114	43	857	935	878

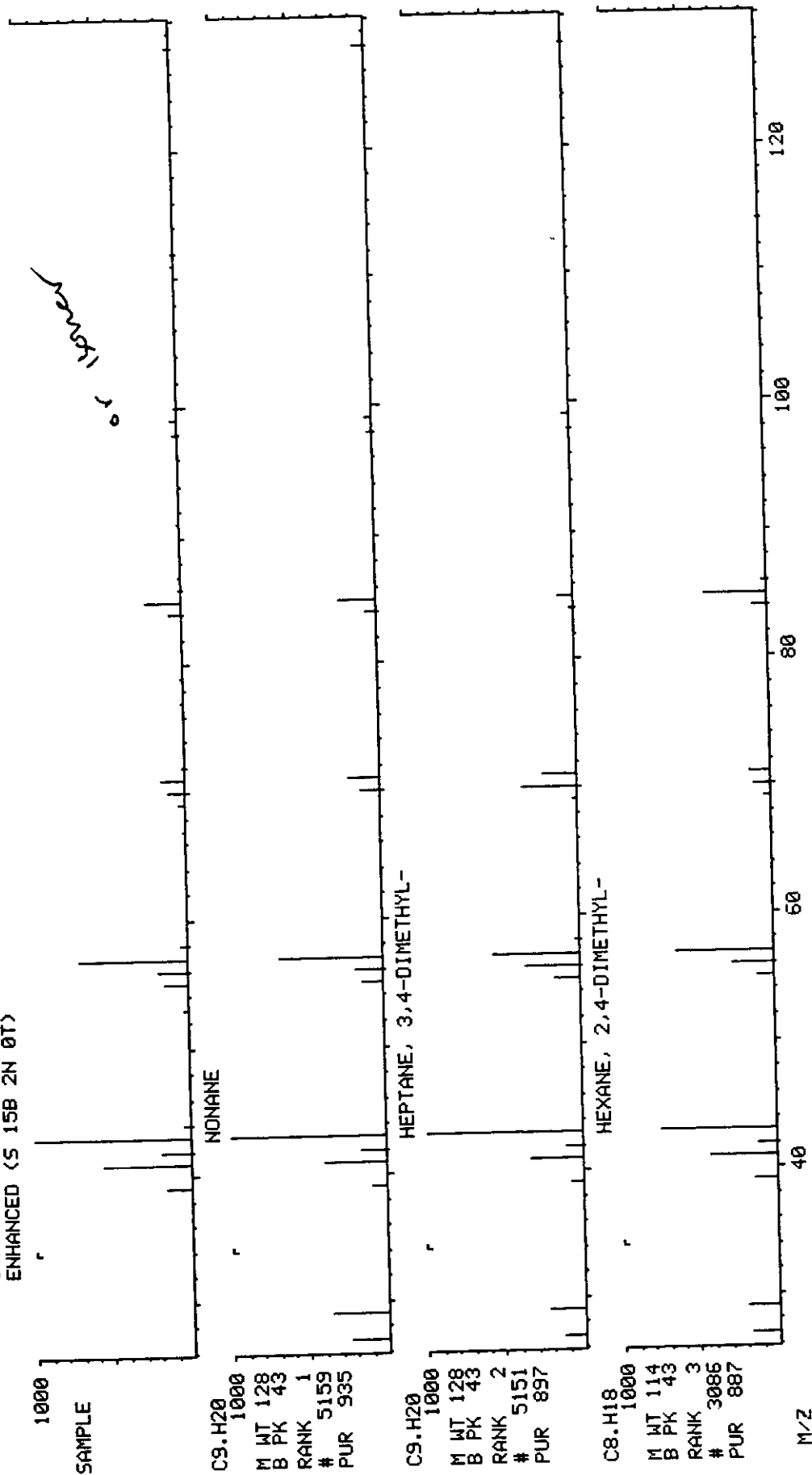
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	111-84-2
2	—	—	—	—	922-28-1
3	—	—	—	—	589-43-5
4	—	—	—	—	2213-23-2
5	—	—	—	—	62108-23-0
6	—	—	—	—	2613-61-8
7	—	—	—	—	2216-32-2
8	—	—	—	—	1120-21-4
9	—	—	—	—	111-65-9

BASE M/Z: 43
RIC: 253952.

DATA: 30058108 # 222
CALI: 30058108 # 3

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 22:29:00 + 2:28
SAMPLE: 5-MM5-FB 1/35A/1ML INST. ID: F16
CONDS.: UG/ML *1ML *100%/100% *(NA/NA) >1/35A NA M
ENHANCED (S 15B 2N 0T)



Library Search Data: 30068108 # 285 Base m/z: 77
 08/31/98 22:29:00 + 3:10 Cali: 30068108 # 3 RIC: 13744.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 366 matched at least 5 of the 16 largest peaks in the unknown

Rank In.	Name
1	2017 BENZALDEHYDE
2	7284 BENZOYL CHLORIDE
3	6466 ETHANONE, 2-HYDROXY-1-PHENYL-
4	2014 2,4-HEPTADIEN-6-YNAL, (E,E)-
5	9645 BENZOYLFORMIC ACID
6	13112 BENZOYL ISOTHIOCYANATE
7	6857 BENZENECARBOTHIOIC ACID
8	13359 BENZENEACETIC ACID, .ALPHA.-OXO-, METHYL ESTER
9	6111 1,2-BENZENEDICARBOXALDEHYDE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C7.H6.O	106	77	923	967	944
2	C7.H5.O.CL	140	105	817	874	830
3	C8.H8.O2	136	105	806	842	822
4	C7.H6.O	106	105	805	846	810
5	C8.H6.O3	150	105	779	846	804
6	C8.H5.O.N.S	163	105	773	820	788
7	C7.H6.O.S	138	77	763	845	810
8	C9.H8.O3	164	105	759	832	787
9	C8.H6.O2	134	105	758	807	915

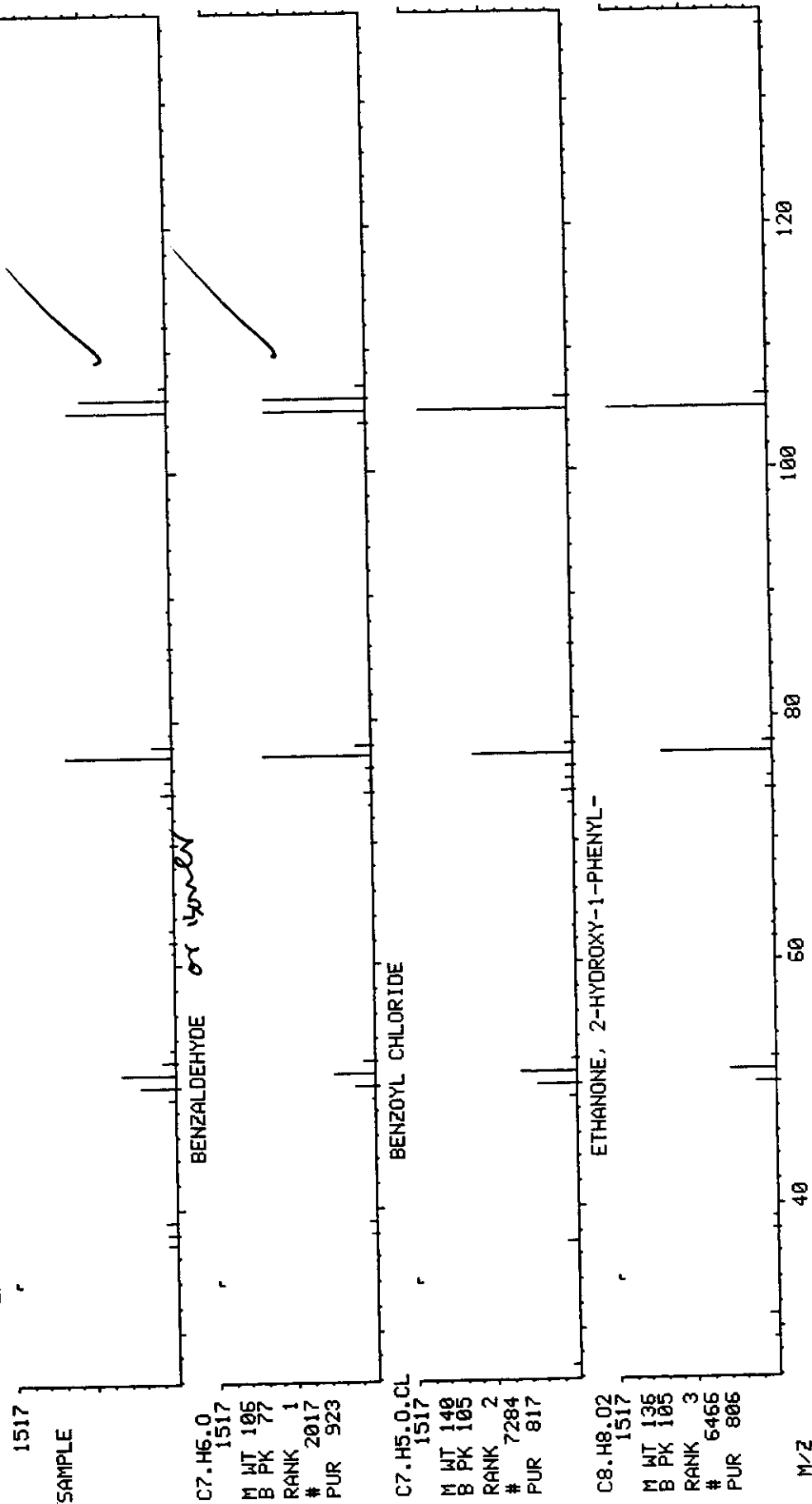
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	100-52-7
2	—	—	—	—	98-88-4
3	—	—	—	—	582-24-1
4	—	—	—	—	7200-04-6
5	—	—	—	—	611-73-4
6	—	—	—	—	532-55-8
7	—	—	—	—	98-91-9
8	—	—	—	—	15206-55-0
9	—	—	—	—	643-79-8

BASE M/Z: 77
RIC: 13744.

DATA: 30068108 # 285
CALI: 30068108 # 3

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 22:29:00 + 3:10
SAMPLE: 5-MM5-FB 1/35A/1ML INST. ID: F16
CONDS.: UG/ML *1ML *100%/100% *(NA/NA) 1/35A NA M
ENHANCED (S 15B 2N 0T)



Library Search Data: 30068108 # 932 Base m/z: 71
 08/31/98 22:29:00 + 10:22 Cali: 30068108 # 3 RIC: 18880.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 540 matched at least 6 of the 16 largest peaks in the unknown

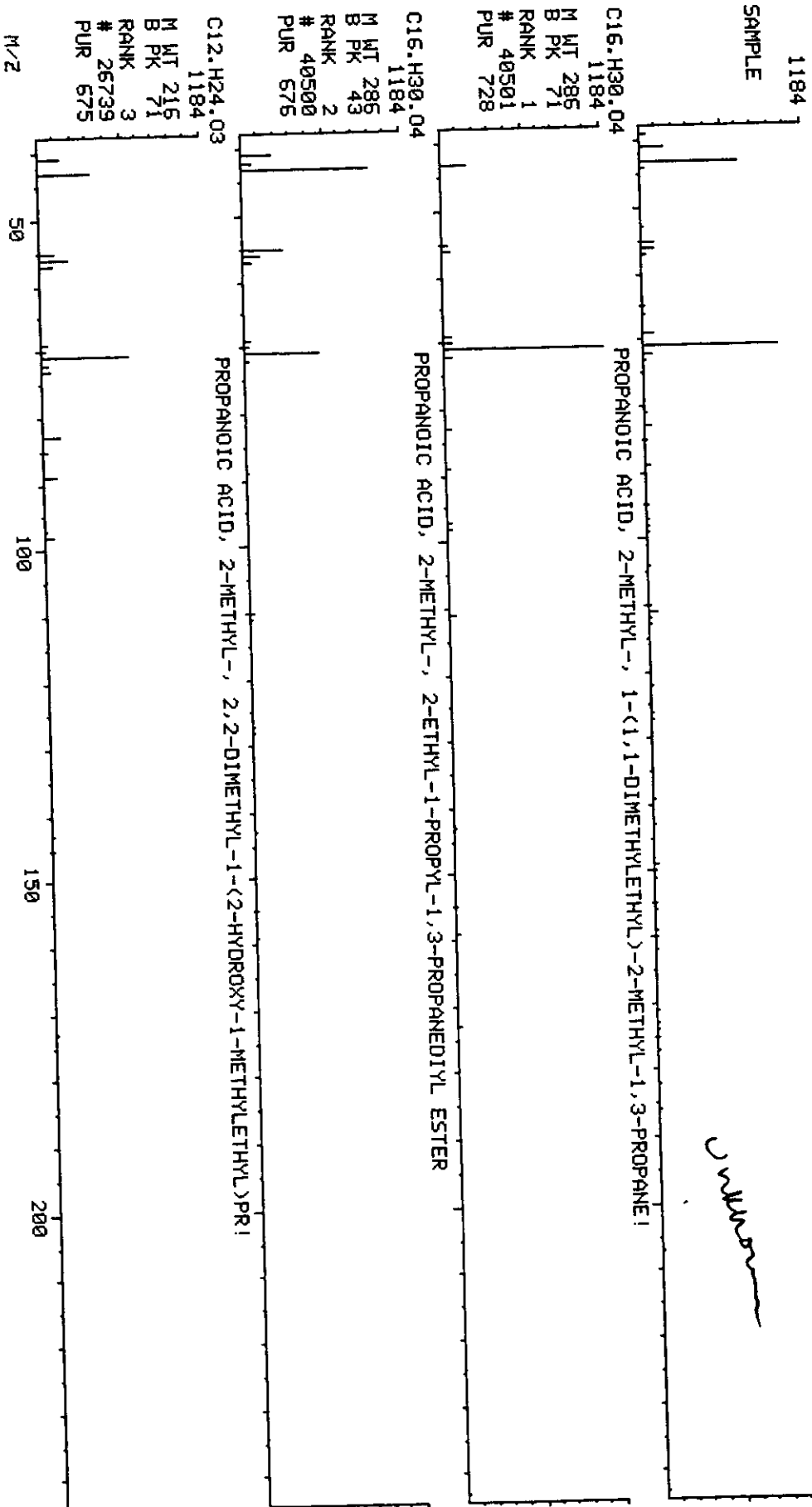
Rank In. Name
 1 40501 PROPANOIC ACID, 2-METHYL-, 1-(1,1-DIMETHYLETHYL)-2-METHYL-1,3-PROPA*
 2 40500 PROPANOIC ACID, 2-METHYL-, 2-ETHYL-1-PROPYL-1,3-PROPANEDIYL ESTER
 3 26739 PROPANOIC ACID, 2-METHYL-, 2,2-DIMETHYL-1-(2-HYDROXY-1-METHYLETHYL)*
 4 26740 PROPANOIC ACID, 2-METHYL-, 3-HYDROXY-2,4,4-TRIMETHYLPENTYL ESTER
 5 4625 1-HEXENE, 3,4,5-TRIMETHYL-
 6 2720 4,4-DIMETHYL-1-HEXENE
 7 4556 ETHANONE, 1-(3-ETHYLCYCLOBUTYL)-
 8 4677 2,4,4-TRIMETHYL-1-HEXENE
 9 26737 PROPANOIC ACID, 2-METHYL-, 2-ETHYL-3-HYDROXYHEXYL ESTER

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C16.H30.O4	286	71	728	915	748
2	C16.H30.O4	286	43	676	875	712
3	C12.H24.O3	216	71	675	888	693
4	C12.H24.O3	216	71	626	800	631
5	C9.H18	126	43	557	798	652
6	C8.H16	112	71	552	788	617
7	C8.H14.O	126	43	546	735	625
8	C9.H18	126	71	527	721	637
9	C12.H24.O3	216	71	514	694	538

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	74381-40-1
2	—	—	—	—	74367-30-9
3	—	—	—	—	74367-33-2
4	—	—	—	—	74367-34-3
5	—	—	—	—	56728-10-0
6	—	—	—	—	1647-08-1
7	—	—	—	—	56335-71-8
8	—	—	—	—	51174-12-0
9	—	—	—	—	74367-31-0

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 10:22 INST. ID: F16
 SAMPLE: S-MM5-FB 1/35A/1ML
 COND5.: UG/ML *1ML *1002/1002 *(NA/NA)/1/35A NA M
 ENHANCED (S 1SB 2N 0T)

DATA: 30068108 # 932 BASE M/Z: 71
 CALI: 30068108 # 3 RIC: 18880.



Library Search Data: 30068108 # 985 Base m/z: 57
 08/31/98 22:29:00 + 10:57 Cali: 30068108 # 3 RIC: 12368.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 548 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37465 NONADECANE
 2 22527 DODECANE, 2,5-DIMETHYL-
 3 22530 TETRADECANE
 4 29264 TRIDECANE, 5-PROPYL-
 5 25997 PENTADECANE
 6 29263 HEXADECANE
 7 32059 HEPTADECANE
 8 37462 HEPTADECANE, 2,6-DIMETHYL-
 9 19054 UNDECANE, 2,6-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	763	926	787
2	C14.H30	198	57	754	934	759
3	C14.H30	198	43	752	933	759
4	C16.H34	226	57	750	977	757
5	C15.H32	212	57	750	934	775
6	C16.H34	226	57	749	916	793
7	C17.H36	240	57	749	915	787
8	C19.H40	268	57	747	936	763
9	C13.H28	184	57	740	943	746

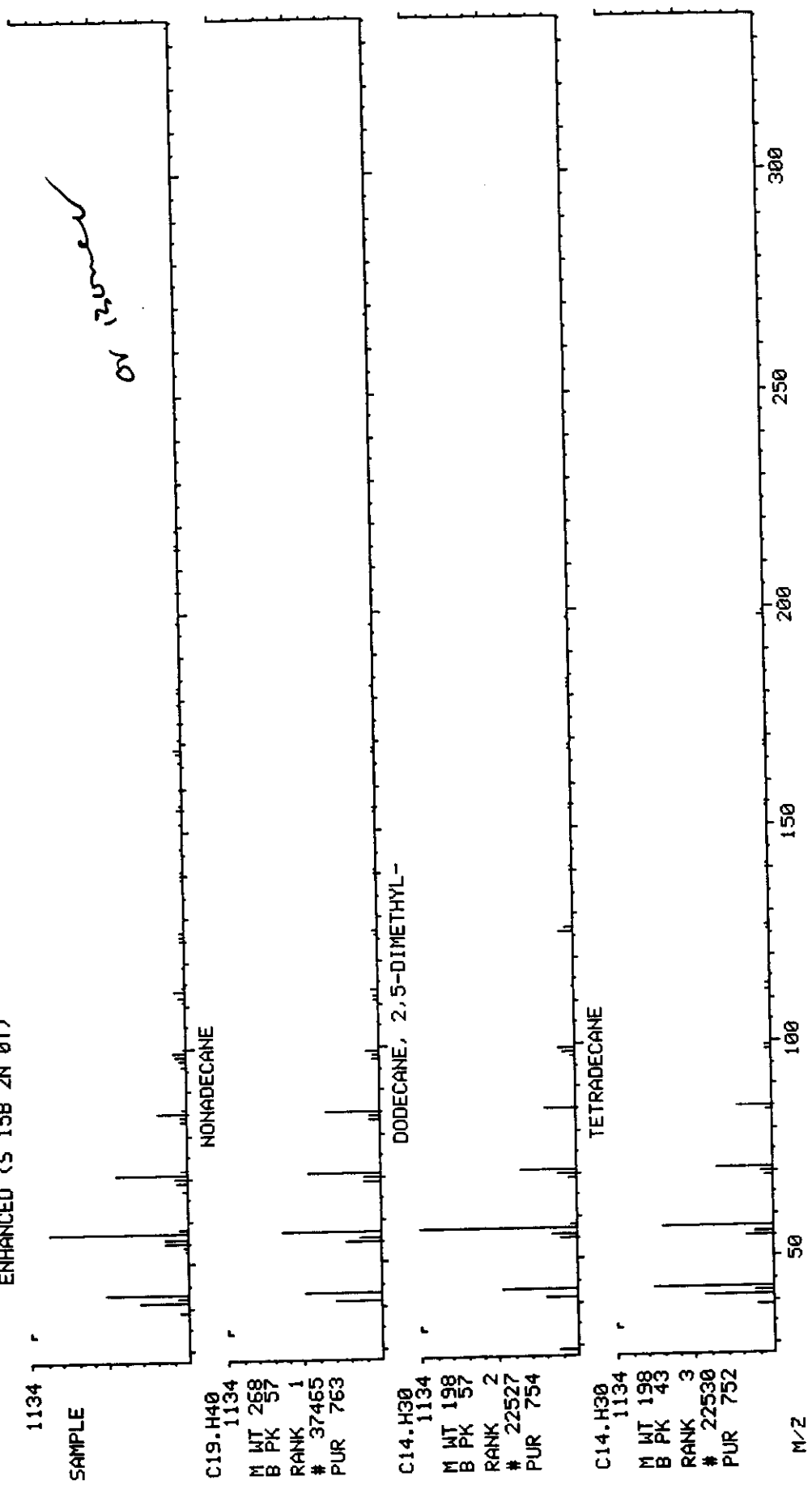
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	56292-65-0
3	—	—	—	—	629-59-4
4	—	—	—	—	55045-11-9
5	—	—	—	—	629-62-9
6	—	—	—	—	544-76-3
7	—	—	—	—	629-78-7
8	—	—	—	—	54105-67-8
9	—	—	—	—	17301-23-4

BASE M/Z: 57
RIC: 12368.

DATA: 30068108 # 985
CALI: 30068108 # 3

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 22:29:00 + 10:57
SAMPLE: S-MMS-FB 1/35A/1ML
CONDS.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



Library Search Data: 30068108 #1033 Base m/z: 57
 08/31/98 22:29:00 + 11:29 Cali: 30068108 # 3 RIC: 23872.
 Sample: S-MMS-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 450 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 3 29253 PENTADECANE, 2-METHYL-
 4 32058 HEXADECANE, 3-METHYL-
 5 37466 PENTADECANE, 2,6,10,14-TETRAMETHYL-
 6 37465 NONADECANE
 7 25991 DODECANE, 2,6,10-TRIMETHYL-
 8 22535 DODECANE, 4,6-DIMETHYL-
 9 42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	746	939	758
2	C21.H44	296	57	727	932	744
3	C16.H34	226	43	702	925	723
4	C17.H36	240	57	701	932	714
5	C19.H40	268	71	697	933	721
6	C19.H40	268	57	697	930	730
7	C15.H32	212	57	692	950	696
8	C14.H30	198	57	690	952	690
9	C21.H44	296	57	687	915	720

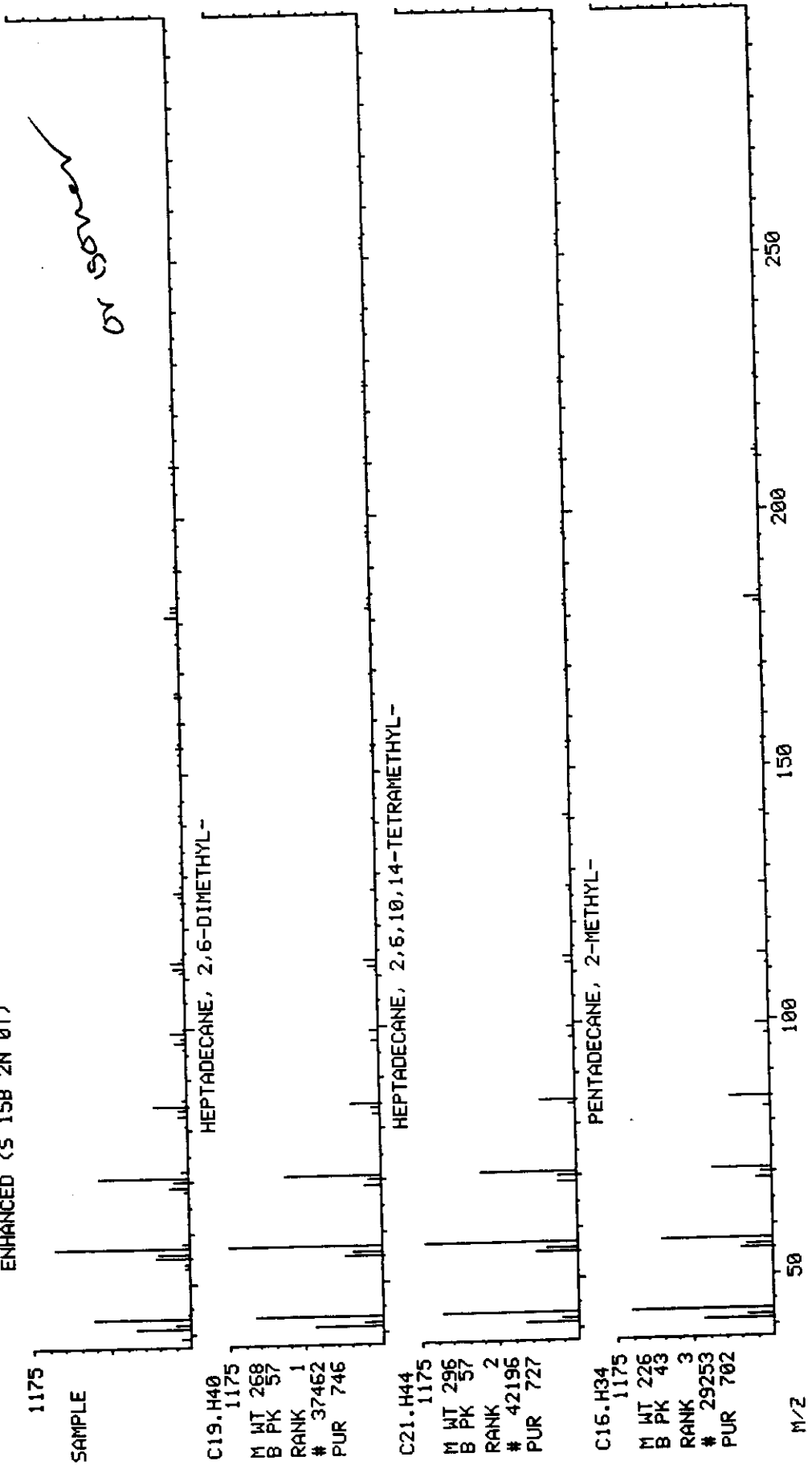
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	18344-37-1
3	—	—	—	—	1560-93-6
4	—	—	—	—	6418-43-5
5	—	—	—	—	1921-70-6
6	—	—	—	—	629-92-5
7	—	—	—	—	3891-98-3
8	—	—	—	—	61141-72-8
9	—	—	—	—	54833-48-6

BASE M/Z: 57
RIC: 23872.

DATA: 30068108 #1033
CALI: 30068108 # 3

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 22:29:00 + 11:29
SAMPLE: S-MMS-FB 1/35A/1ML
CONDS.: UG/ML *100%/100% *(NA/NA)/1/35A NA M
ENHANCED (S 15B 2N 0T)

INST. ID: F16



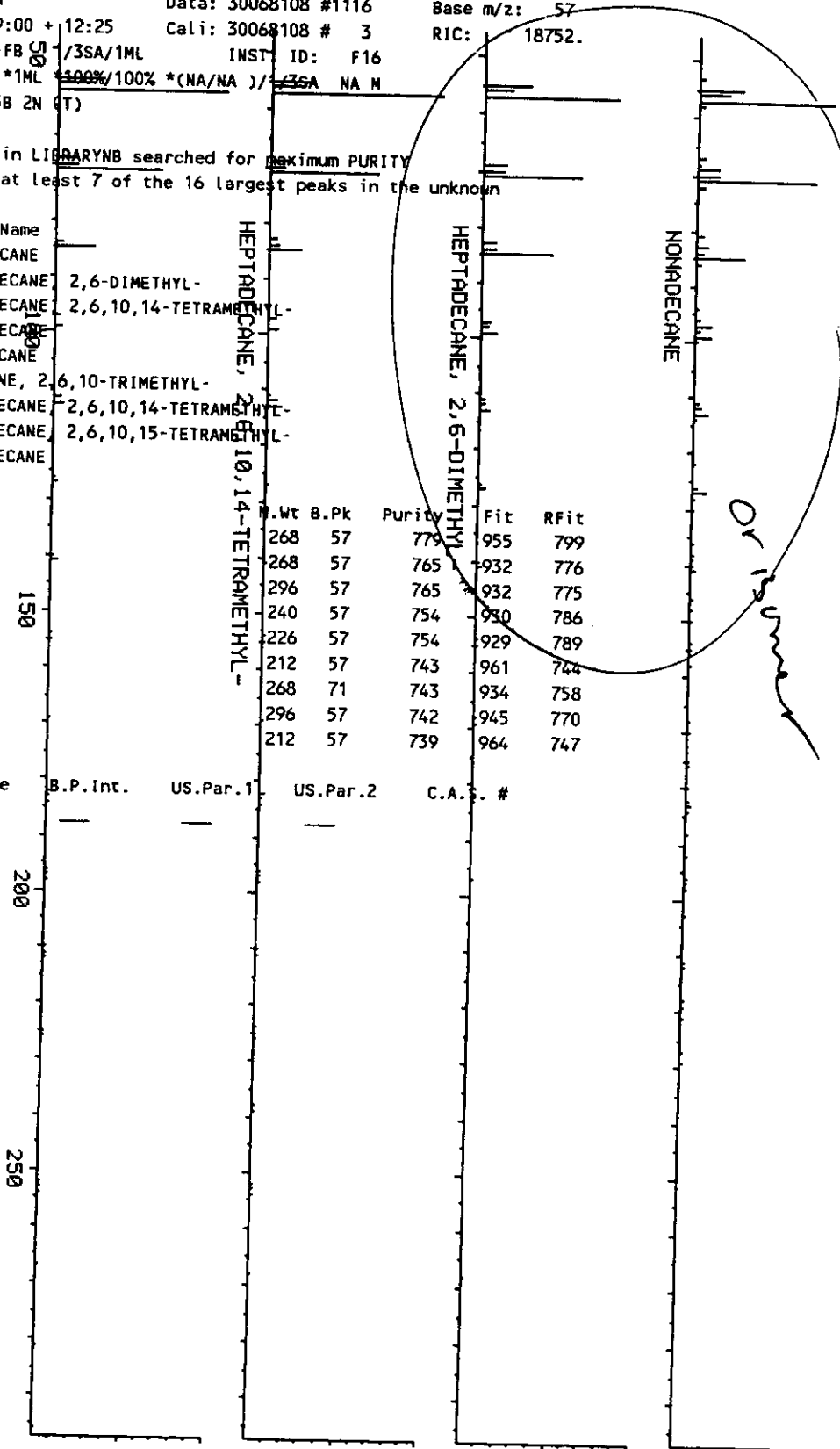
Library Search Data: 30068108 #1116 Base m/z: 57
 08/31/98 22:29:00 + 12:25 Cali: 30068108 # 3 RIC: 18752.
 Sample: S-MM5-FB 1/35A/1ML INST ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 848 matched at least 7 of the 16 largest peaks in the unknown

Rank In.	Name
1	37465 NONADECANE
2	37462 HEPTADECANE, 2,6-DIMETHYL-
3	42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
4	32059 HEPTADECANE
5	29263 HEXADECANE
6	25991 DODECANE, 2,6,10-TRIMETHYL-
7	37466 PENTADECANE, 2,6,10,14-TETRAMETHYL-
8	42192 HEPTADECANE, 2,6,10,15-TETRAMETHYL-
9	25997 PENTADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	779	955	799
2	C19.H40	268	57	765	932	776
3	C21.H44	296	57	765	932	775
4	C17.H36	240	57	754	930	786
5	C16.H34	226	57	754	929	789
6	C15.H32	212	57	743	961	744
7	C19.H40	268	71	743	934	758
8	C21.H44	296	57	742	945	770
9	C15.H32	212	57	739	964	747

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1					



MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 12:25
 SAMPLE: S-MM5-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)

INST. ID: F16

DATA: 30068108 #1116
 CALI: 30068108 # 3

BASE M/Z: 57
 RIC: 18752.

Library Search Data: 30068108 #1235 Base m/z: 43
 08/31/98 22:29:00 + 13:44 Cali: 30068108 # 3 RIC: 19456.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 912 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 35182 HEXADECANOIC ACID
 2 29642 TETRADECANOIC ACID
 3 37913 GLYCINE, N-METHYL-N-(1-OXODODECYL)-
 4 40669 DODECANAMIDE, N,N-BIS(2-HYDROXYETHYL)-
 5 22949 DODECANOIC ACID
 6 40184 OCTADECANOIC ACID
 7 26351 TRIDECANOIC ACID
 8 19469 UNDECANOIC ACID
 9 32392 PENTADECANOIC ACID

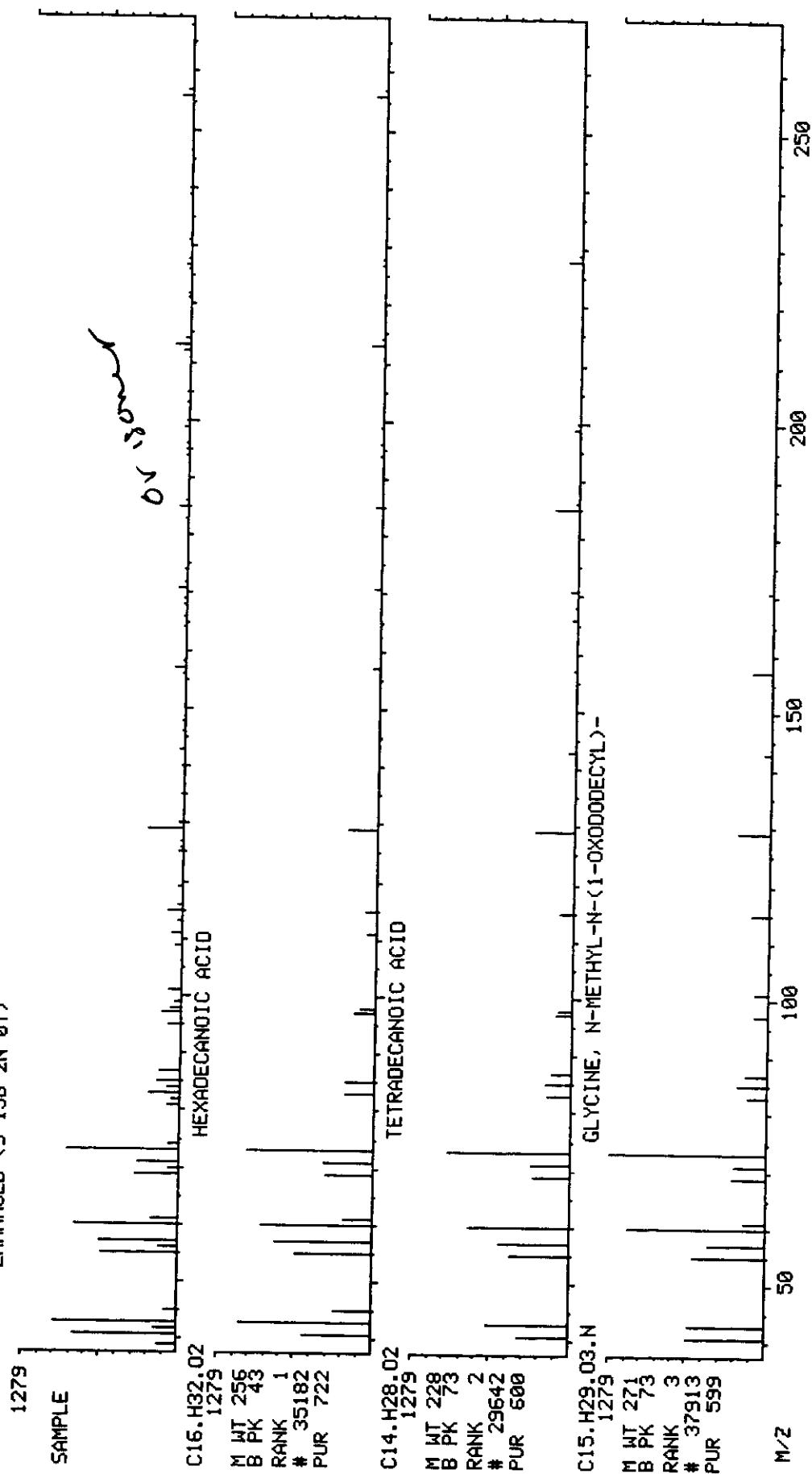
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C16.H32.O2	256	43	722	977	722
2	C14.H28.O2	228	73	600	936	607
3	C15.H29.O3.N	271	73	599	975	599
4	C16.H33.O3.N	287	73	585	967	589
5	C12.H24.O2	200	60	579	941	590
6	C18.H36.O2	284	43	561	776	707
7	C13.H26.O2	214	73	550	885	568
8	C11.H22.O2	186	60	539	940	548
9	C15.H30.O2	242	43	525	810	626

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	57-10-3
2	—	—	—	—	544-63-8
3	—	—	—	—	97-78-9
4	—	—	—	—	120-40-1
5	—	—	—	—	143-07-7
6	—	—	—	—	57-11-4
7	—	—	—	—	638-53-9
8	—	—	—	—	112-37-8
9	—	—	—	—	1002-84-2

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 13:44
 SAMPLE: S-MMS-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA) /1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068108 #1235
 CALI: 30068108 # 3
 BASE M/Z: 43
 RIC: 19456.

INST. ID: F16
 NA M



Library Search Data: 30068108 #1334 Base m/z: 57
 08/31/98 22:29:00 + 14:50 Cali: 30068108 # 3 RIC: 13856.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (\$ 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 416 matched at least 7 of the 16 largest peaks in the unknown

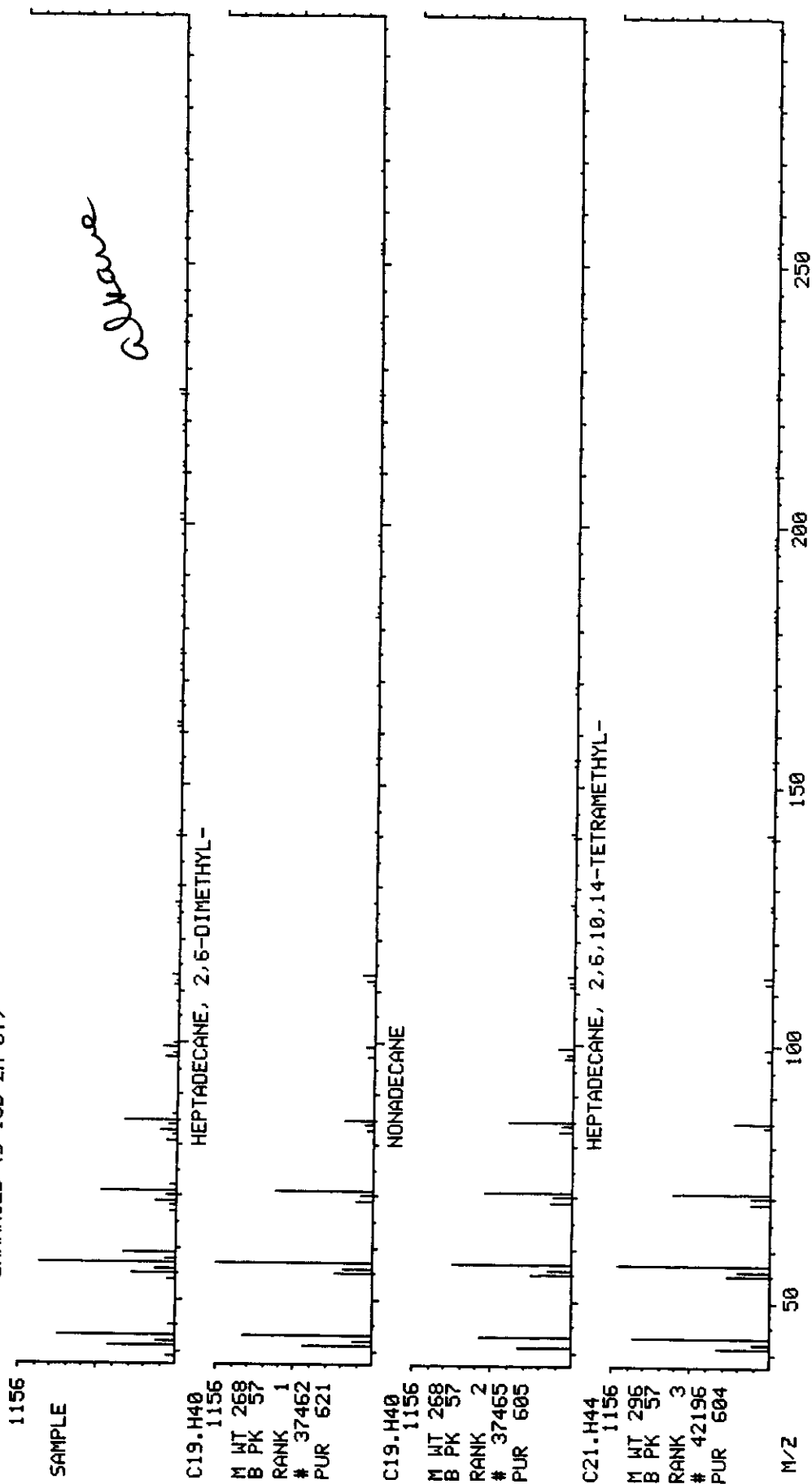
Rank In. Name
 1 37462 HEPTADECANE, 2,6-DIMETHYL-
 2 37465 NONADECANE
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 15969 HYDROXYLAMINE, O-DECYL-
 5 46161 TRICOSANE
 6 32058 HEXADECANE, 3-METHYL-
 7 39858 EICOSANE
 8 37456 2-METHYLOCTADECANE
 9 49555 PENTACOSANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	621	892	637
2	C19.H40	268	57	605	954	627
3	C21.H44	296	57	604	898	622
4	C10.H23.O.N	173	43	598	968	598
5	C23.H48	324	43	598	885	664
6	C17.H36	240	57	594	922	610
7	C20.H42	282	57	594	867	653
8	C19.H40	268	43	591	903	633
9	C25.H52	352	43	589	864	672

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	54105-67-8
2	—	—	—	—	629-92-5
3	—	—	—	—	18344-37-1
4	—	—	—	—	29812-79-1
5	—	—	—	—	638-67-5
6	—	—	—	—	6418-43-5
7	—	—	—	—	112-95-8
8	—	—	—	—	- -
9	—	—	—	—	629-99-2

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 14:50
 SAMPLE: S-MM5-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA) /1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068108 #1334
 CALI: 30068108 # 3
 BASE M/Z: 57
 RIC: 13856.



1214

Library Search Data: 30068108 #1361 Base m/z: 59
 08/31/98 22:29:00 + 15:08 Cali: 30068108 # 3 RIC: 4168.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 483 matched at least 3 of the 16 largest peaks in the unknown

Rank In. Name
 1 20265 3-HEXANOL, 1,5-DIMETHOXY-2,4-DIMETHYL-
 2 36537 2,5,8,11-TETRAOXATETRADECAN-13-OL, 4,7,10-TRIMETHYL-
 3 16192 2-PROPANOL, 1-[(1-METHYL-2-(2-PROPENYLOXY)ETHOXY)-
 4 20626 2-PROPANOL, 1,1'-[(1-METHYL-1,2-ETHANEDIYL)BIS(OXY)]BIS-
 5 9235 1-PROPANOL, 2-(2-METHOXY-1-METHYLETHOXY)-
 6 5826 2-PROPANOL, 1-ISOPROPOXY-2-METHYL-
 7 5786 PROPANOIC ACID, 2-HYDROXY-2-METHYL-, ETHYL ESTER
 8 12385 BUTYRIC ACID, 4-ISOPROPOXY-, METHYL ESTER
 9 5790 BUTANOIC ACID, 2-HYDROXY-, ETHYL ESTER

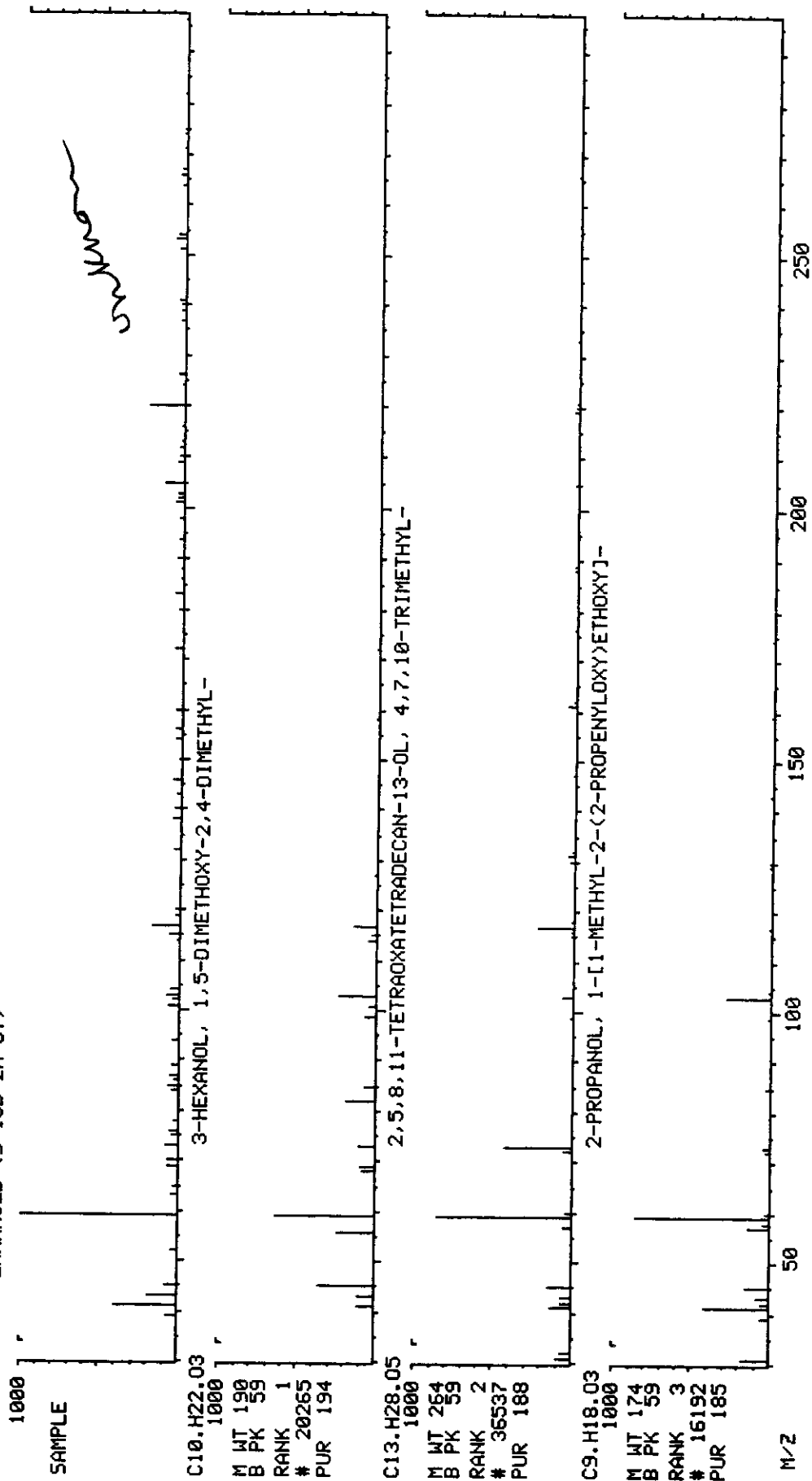
Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C10.H22.O3	190	59	194	664	256
2	C13.H28.O5	264	59	188	462	250
3	C9.H18.O3	174	59	185	715	214
4	C9.H20.O4	192	59	184	763	211
5	C7.H16.O3	148	59	182	756	189
6	C7.H16.O2	132	59	178	717	197
7	C6.H12.O3	132	59	174	778	192
8	C8.H16.O3	160	85	173	663	199
9	C6.H12.O3	132	59	172	797	193

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	13897-22-8
2	—	—	—	—	20324-34-9
3	—	—	—	—	55956-25-7
4	—	—	—	—	1638-16-0
5	—	—	—	—	55956-21-3
6	—	—	—	—	3587-75-5
7	—	—	—	—	80-55-7
8	—	—	—	—	29006-05-1
9	—	—	—	—	52089-54-0

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 15:08
 SAMPLE: S-MM5-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30058108 #1361
 CALI: 30058108 # 3

BASE M/Z: 59
 RIC: 4168.



Library Search Data: 30068108 #1401 Base m/z: 57
 08/31/98 22:29:00 + 15:35 Cali: 30068108 # 3 RIC: 11632.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 743 matched at least 7 of the 16 largest peaks in the unknown

Rank In. Name
 1 46161 TRICOSANE
 2 49555 PENTACOSANE
 3 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 4 37462 HEPTADECANE, 2,6-DIMETHYL-
 5 37465 NONADECANE
 6 58739 PENTATRIACONTANE
 7 42197 HENEICOSANE
 8 42193 EICOSANE, 10-METHYL-
 9 44314 DOCOSANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C23.H48	324	43	624	903	676
2	C25.H52	352	43	615	883	684
3	C21.H44	296	57	604	915	623
4	C19.H40	268	57	603	896	627
5	C19.H40	268	57	601	944	630
6	C35.H72	492	57	596	891	654
7	C21.H44	296	57	591	877	646
8	C21.H44	296	57	589	878	636
9	C22.H46	310	57	588	869	659

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	638-67-5
2	—	—	—	—	629-99-2
3	—	—	—	—	18344-37-1
4	—	—	—	—	54105-67-8
5	—	—	—	—	629-92-5
6	—	—	—	—	630-07-9
7	—	—	—	—	629-94-7
8	—	—	—	—	54833-23-7
9	—	—	—	—	629-97-0

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 22:29:00 + 15:35

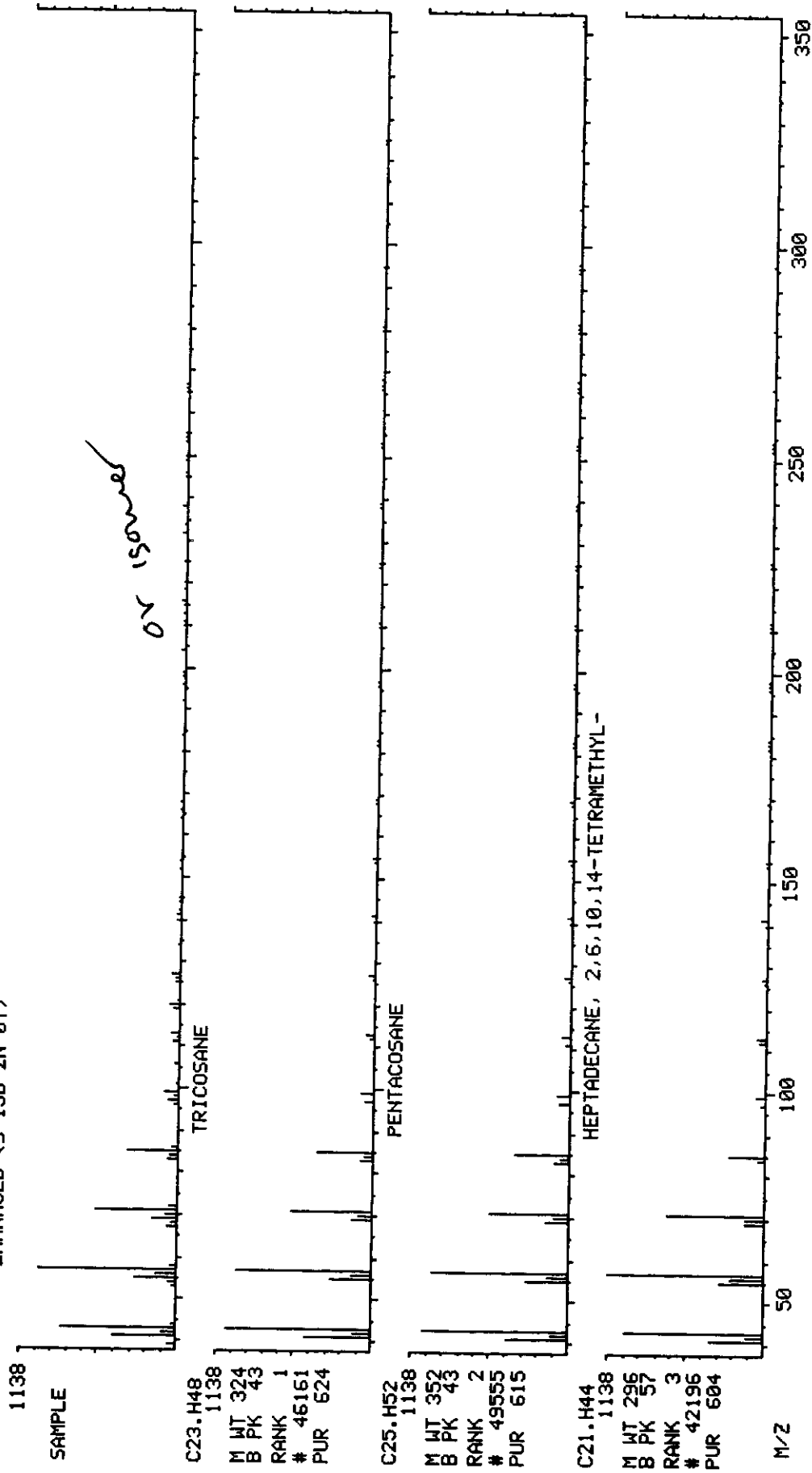
SAMPLE: S-MMS-FB 1/35A/1ML

COND.: UG/ML *1ML *100%/100% *(NA/NA) 1/35A NA M

ENHANCED (S 15B 2N 0T)

DATA: 30068108 #1401
CALI: 30068108 # 3

BASE M/Z: 57
RIC: 11632.



Library Search Data: 30068108 #1467 Base m/z: 57
 08/31/98 22:29:00 + 16:19 Cali: 30068108 # 3 RIC: 10384.
 Sample: S-MMS-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N OT)

62231 spectra in LIBRARYNB searched for maximum PURITY
 928 matched at least 7 of the 16 largest peaks in the unknown

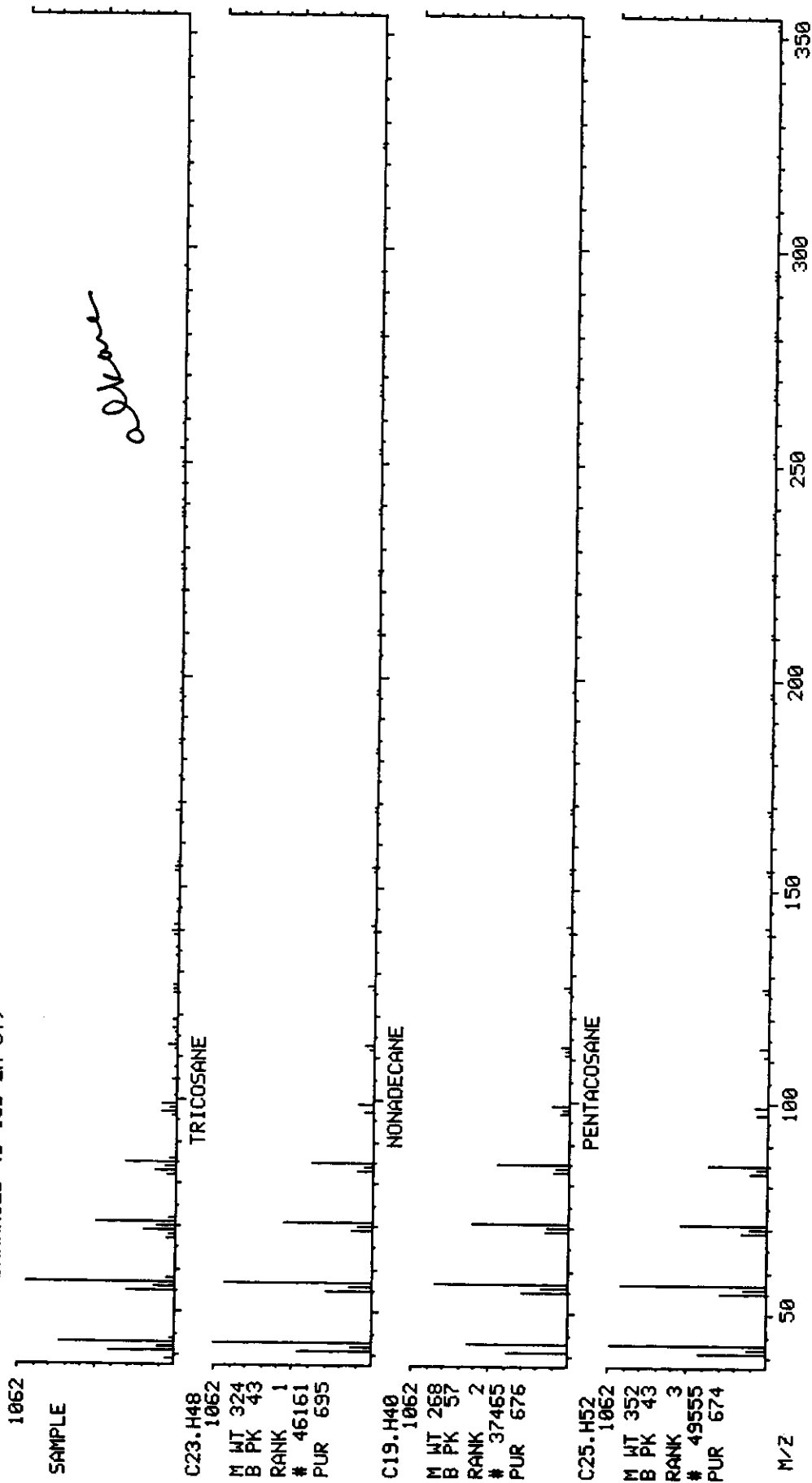
Rank In. Name
 1 46161 TRICOSANE
 2 37465 NONADECANE
 3 49555 PENTACOSANE
 4 39858 EICOSANE
 5 42193 EICOSANE, 10-METHYL-
 6 37462 HEPTADECANE, 2,6-DIMETHYL-
 7 42197 HENEICOSANE
 8 42196 HEPTADECANE, 2,6,10,14-TETRAMETHYL-
 9 58739 PENTATRIACONTANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C23.H48	324	43	695	900	753
2	C19.H40	268	57	676	941	705
3	C25.H52	352	43	674	875	757
4	C20.H42	282	57	665	870	745
5	C21.H44	296	57	661	877	706
6	C19.H40	268	57	659	883	687
7	C21.H44	296	57	657	882	704
8	C21.H44	296	57	649	895	673
9	C35.H72	492	57	647	879	722

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	638-67-5
2	—	—	—	—	629-92-5
3	—	—	—	—	629-99-2
4	—	—	—	—	112-95-8
5	—	—	—	—	54833-23-7
6	—	—	—	—	54105-67-8
7	—	—	—	—	629-94-7
8	—	—	—	—	18344-37-1
9	—	—	—	—	630-07-9

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 16:19
 SAMPLE: S-MM5-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA) 1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30058108 #1467
 CALI: 30058108 # 3
 BASE M/Z: 57
 RIC: 10384.



11220

Library Search Data: 30068108 #1501 Base m/z: 55
 08/31/98 22:29:00 + 16:42 Cali: 30068108 # 3 RIC: 4408.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 742 matched at least 5 of the 16 largest peaks in the unknown

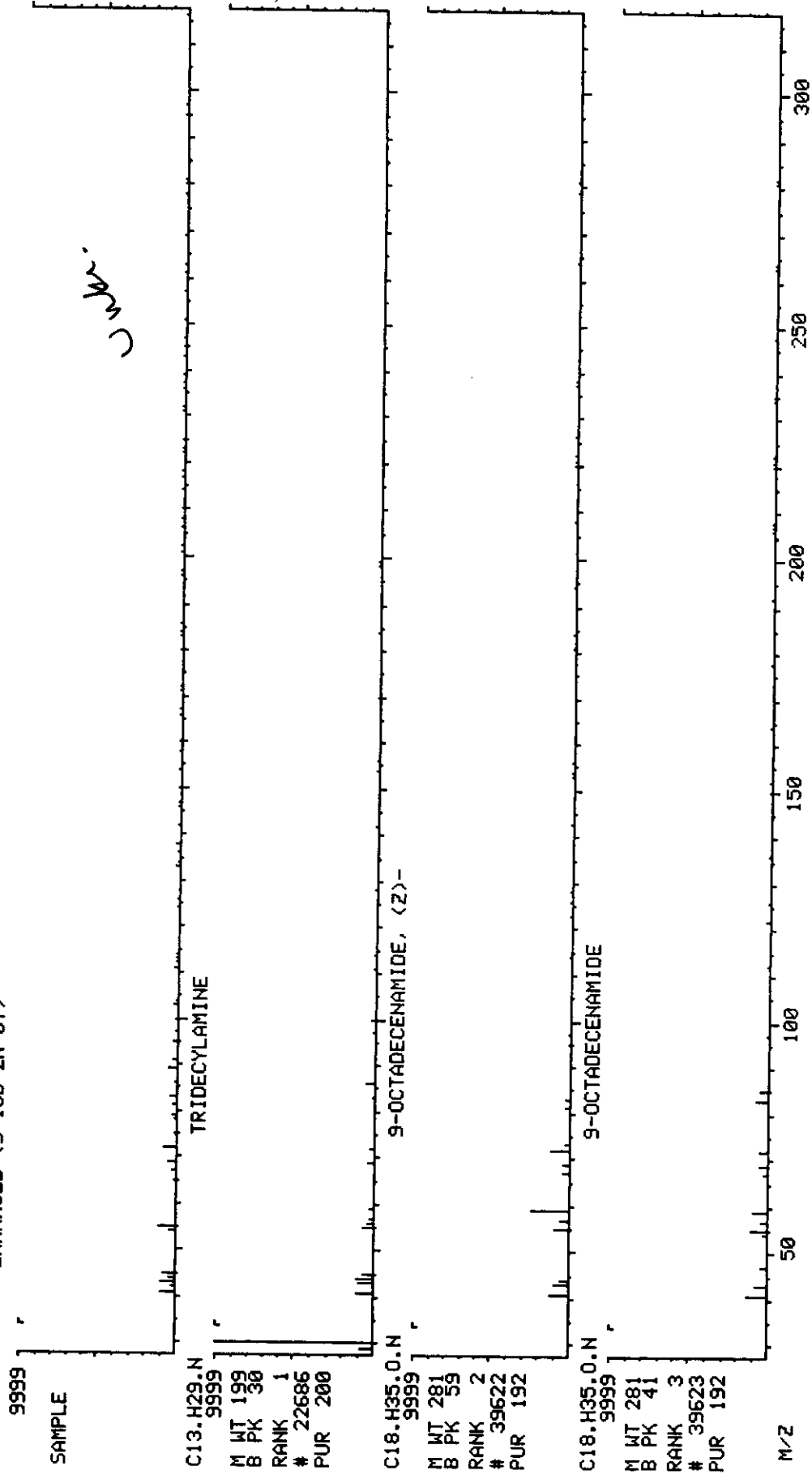
Rank In. Name
 1 22686 TRIDECYLAMINE
 2 39622 9-OCTADECENAMIDE, (Z)-
 3 39623 9-OCTADECENAMIDE
 4 29380 PENTADECYLAMINE
 5 58185 9-HEXADECENOIC ACID, 9-HEXADECENYL ESTER, (Z,Z)-
 6 36536 PHOSPHONOCHLORIDOUS ACID, (1-METHYLETHYL)-, 5-METHYL-2-(1-METHYLETH*
 7 38358 HEXANAL, 4-BROMO-6,6-DICHLORO-2,2-DIMETHYL-
 8 8851 CYCLOPROPANE, 1-CHLORO-1-ETHYL-2,2,3-TRIMETHYL-
 9 48422 1,22-DOCOSANEDIOL

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C13.H29.N	199	30	200	696	256
2	C18.H35.O.N	281	59	192	586	308
3	C18.H35.O.N	281	41	192	594	296
4	C15.H33.N	227	30	189	683	243
5	C32.H60.O2	476	55	184	656	260
6	C13.H26.O.CL.P	264	83	181	687	228
7	C8.H13.O.CL2.8R	274	72	176	621	249
8	C8.H15.CL	146	41	171	744	210
9	C22.H46.O2	342	55	169	587	272

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	2869-34-3
2	—	—	—	—	301-02-0
3	—	—	—	—	3322-62-1
4	—	—	—	—	2570-26-5
5	—	—	—	—	22393-97-1
6	—	—	—	—	74630-89-0
7	—	—	—	—	- -
8	—	—	—	—	61142-56-1
9	—	—	—	—	22513-81-1

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 16:42
 SAMPLE: S-MM5-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA) 1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30058108 #1501
 CALI: 30058108 # 3
 BASE M/Z: 55
 RIC: 4408.



1222

Library Search Data: 30068108 #1509 Base m/z: 59
 08/31/98 22:29:00 + 16:47 Cali: 30068108 # 3 RIC: 6432.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 388 matched at least 6 of the 16 largest peaks in the unknown

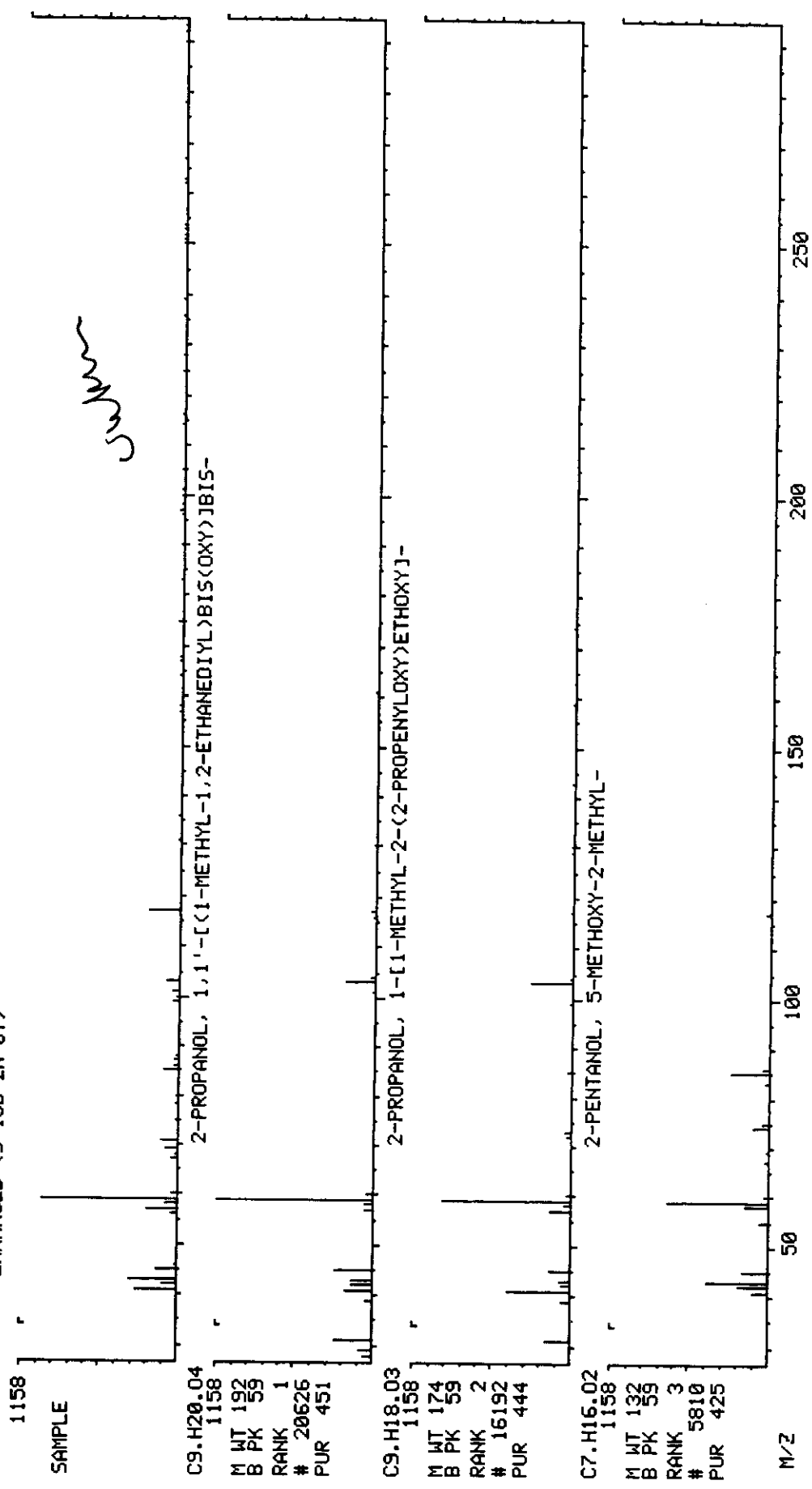
Rank	In.	Name
1	20626	2-PROPANOL, 1,1'-[(1-METHYL-1,2-ETHANEDIYL)BIS(OXY)]BIS-
2	16192	2-PROPANOL, 1-[1-METHYL-2-(2-PROPENYLOXY)ETHOXY]-
3	5810	2-PENTANOL, 5-METHOXY-2-METHYL-
4	3502	BUTANOIC ACID, 3-HYDROXY-3-METHYL-
5	3522	2,3-BUTANEDIOL, 2,3-DIMETHYL-
6	8432	2-HEXANONE, 3-HYDROXY-3,5-DIMETHYL-
7	5832	2-PROPANOL, 1-(1-METHYLPROPOXY)-
8	5809	1-ETHOXYPENTAN-3-OL
9	8805	1,3-DIOXAN-5-OL, 4,4,5-TRIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C9.H20.O4	192	59	451	816	477
2	C9.H18.O3	174	59	444	753	478
3	C7.H16.O2	132	59	425	784	466
4	C5.H10.O3	118	59	423	778	489
5	C6.H14.O2	118	59	419	893	434
6	C8.H16.O2	144	59	415	899	432
7	C7.H16.O2	132	43	414	722	440
8	C7.H16.O2	132	59	408	802	449
9	C7.H14.O3	146	59	407	866	442

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	1638-16-0
2	—	—	—	—	55956-25-7
3	—	—	—	—	55724-04-4
4	—	—	—	—	625-08-1
5	—	—	—	—	76-09-5
6	—	—	—	—	6321-14-8
7	—	—	—	—	53907-95-2
8	—	—	—	—	-
9	—	—	—	—	54063-14-8

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 16:47
 SAMPLE: S-MM5-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *1002/1002 *(NA/NA >1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068108 #1509
 CALI: 30068108 # 3
 BASE M/Z: 59
 RIC: 6432.



1224

Library Search Data: 30068108 #1529 Base m/z: 57
 08/31/98 22:29:00 + 17:00 Cal: 30068108 # 3 RIC: 10960.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 867 matched at least 7 of the 16 largest peaks in the unknown

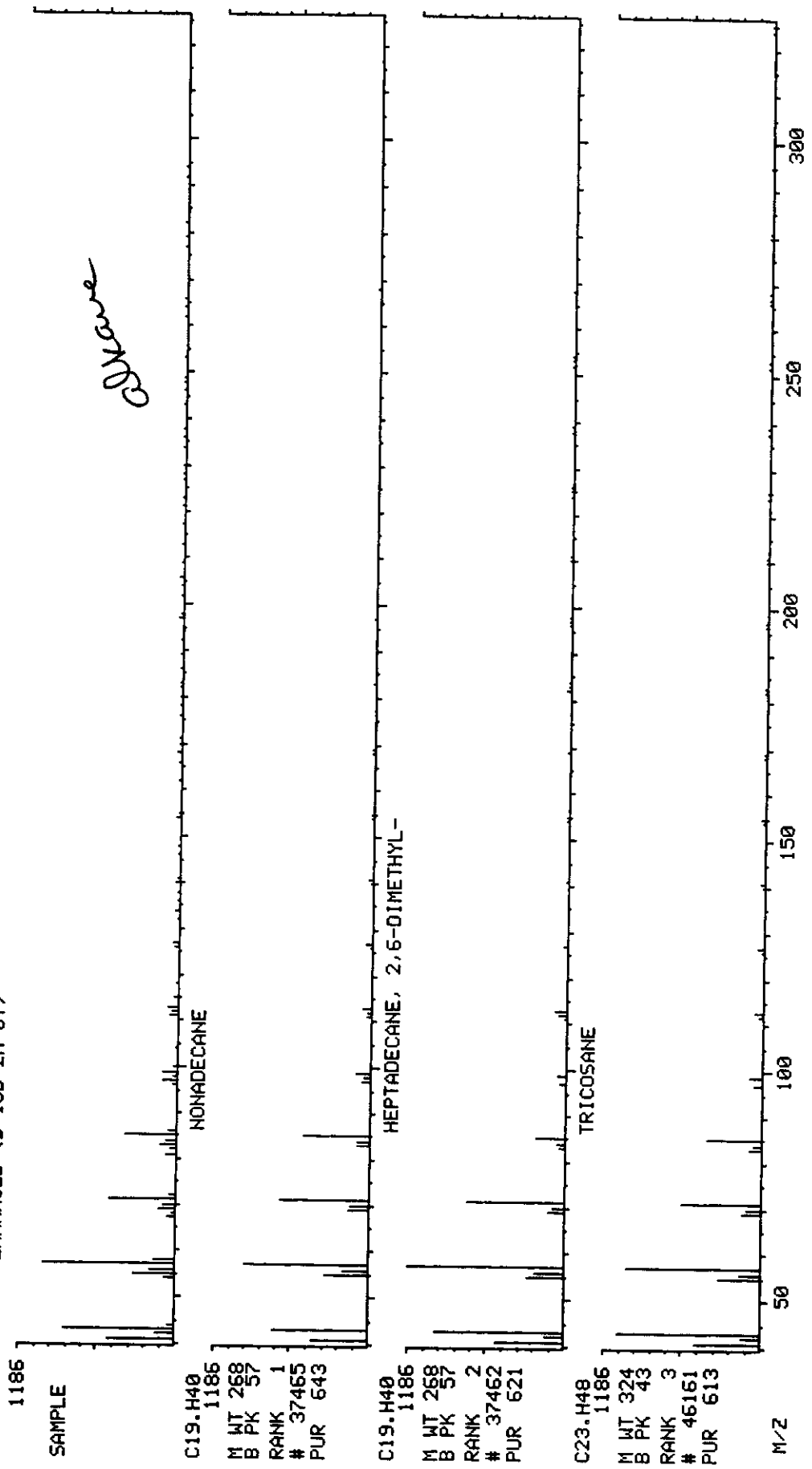
Rank	In.	Name
1	37465	NONADECANE
2	37462	HEPTADECANE, 2,6-DIMETHYL-
3	46161	TRICOSANE
4	42193	EICOSANE, 10-METHYL-
5	49555	PENTACOSANE
6	32059	HEPTADECANE
7	29263	HEXADECANE
8	42196	HEPTADECANE, 2,6,10,14-TETRAMETHYL-
9	39858	EICOSANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	643	955	659
2	C19.H40	268	57	621	934	630
3	C23.H48	324	43	613	904	664
4	C21.H44	296	57	611	909	643
5	C25.H52	352	43	606	880	675
6	C17.H36	240	57	597	928	626
7	C16.H34	226	57	597	924	627
8	C21.H44	296	57	596	929	610
9	C20.H42	282	57	593	893	643

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	54105-67-8
3	—	—	—	—	638-67-5
4	—	—	—	—	54833-23-7
5	—	—	—	—	629-99-2
6	—	—	—	—	629-78-7
7	—	—	—	—	544-76-3
8	—	—	—	—	18344-37-1
9	—	—	—	—	112-95-8

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 17:00
 SAMPLE: S-MNS-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA) 1/1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068108 #1529
 CALI: 30068108 # 3
 BASE M/Z: 57
 RIC: 10960.



1230

Library Search Data: 30068108 #1588 Base m/z: 43
 08/31/98 22:29:00 + 17:40 Cali: 30068108 # 3 RIC: 39616.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 950 matched at least 7 of the 16 largest peaks in the unknown

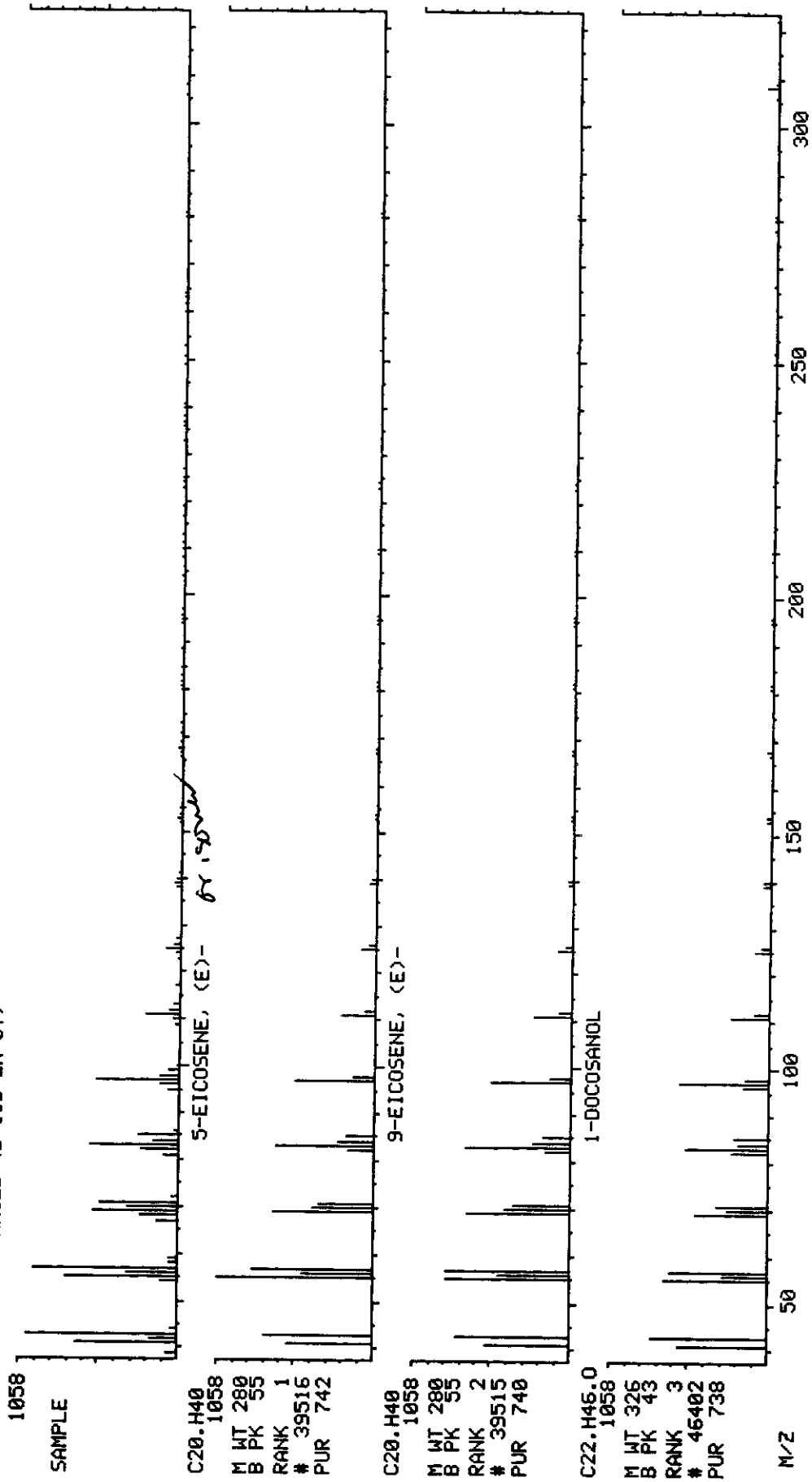
Rank In. Name
 1 39516 5-EICOSENE, (E)-
 2 39515 9-EICOSENE, (E)-
 3 46402 1-DOCOSANOL
 4 58701 17-PENTATRIACONTENE
 5 34410 3-OCTADECENE, (E)-
 6 31653 1-HEPTADECENE
 7 35209 1-HEXADECANOL, 2-METHYL-
 8 39517 3-EICOSENE, (E)-
 9 34411 5-OCTADECENE, (E)-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C20.H40	280	55	742	975	743
2	C20.H40	280	55	740	973	741
3	C22.H46.0	326	43	738	942	755
4	C35.H70	490	43	738	959	765
5	C18.H36	252	69	736	962	747
6	C17.H34	238	55	736	974	738
7	C17.H36.0	256	57	736	977	741
8	C20.H40	280	57	731	979	733
9	C18.H36	252	55	731	956	741

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	74685-30-6
2	—	—	—	—	74685-29-3
3	—	—	—	—	661-19-8
4	—	—	—	—	6971-40-0
5	—	—	—	—	7206-19-1
6	—	—	—	—	6765-39-5
7	—	—	—	—	2490-48-4
8	—	—	—	—	74685-33-9
9	—	—	—	—	7206-21-5

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 17:40
 SAMPLE: S-MM5-FB 1/35A/1ML
 CONDS.: UG/ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068108 #1588
 CALL: 30068108 # 3
 BASE M/Z: 43
 RIC: 39616.



Library Search Data: 30068108 #1661 Base m/z: 59
 08/31/98 22:29:00 + 18:29 Cali: 30068108 # 3 RIC: 5184.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 894 matched at least 5 of the 16 largest peaks in the unknown

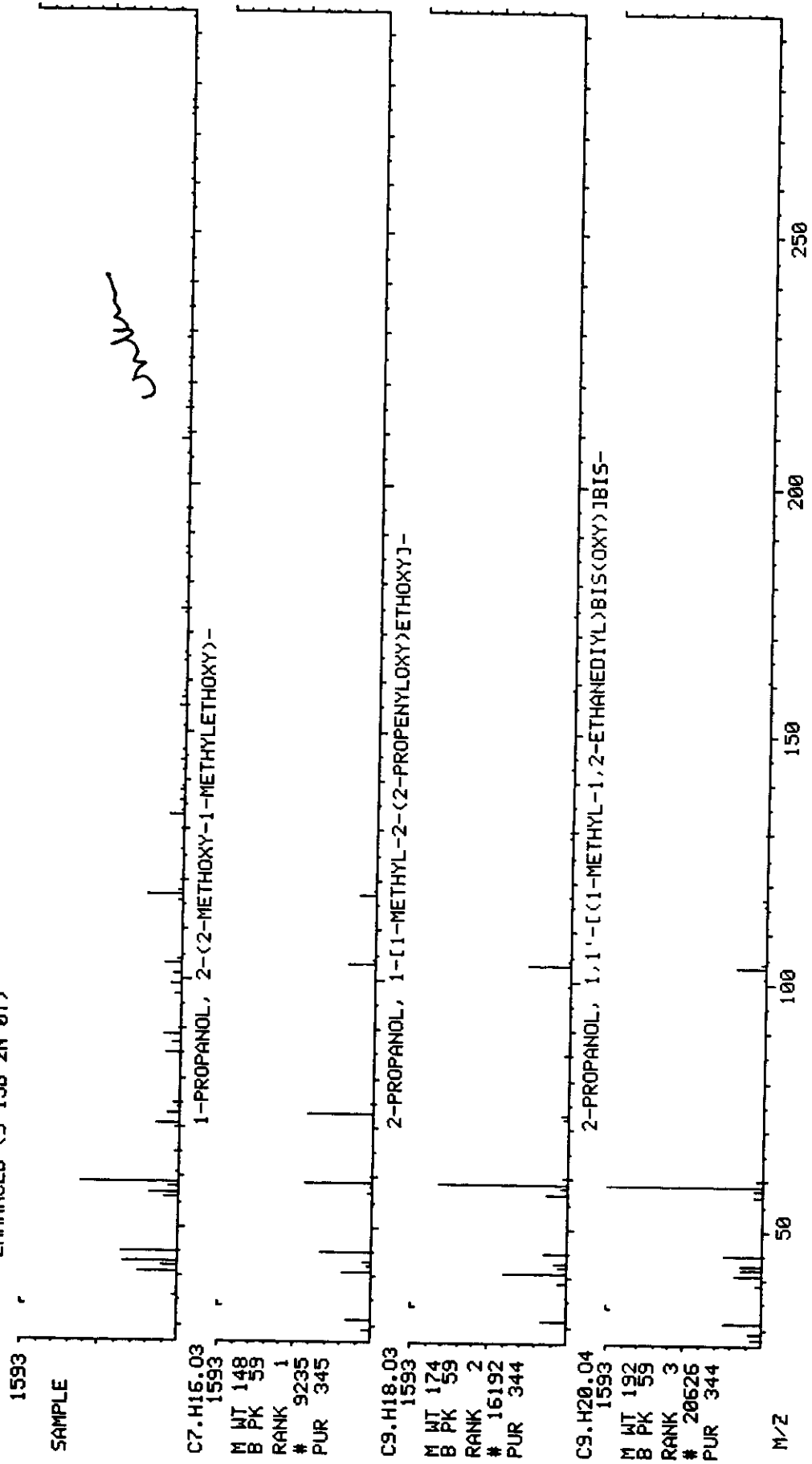
Rank In. Name
 1 9235 1-PROPANOL, 2-(2-METHOXY-1-METHYLETHOXY)-
 2 16192 2-PROPANOL, 1-[1-METHYL-2-(2-PROPENYLOXY)ETHOXY]-
 3 20626 2-PROPANOL, 1,1'-[1-(1-METHYL-1,2-ETHANEDIYL)BIS(OXY)]BIS-
 4 5832 2-PROPANOL, 1-(1-METHYLPROPOXY)-
 5 31381 6-O-METHYL-2,4-METHYLENE-.BETA.-SEDOHEPTITOL
 6 24247 2-PROPANOL, 1-[2-(2-METHOXY-1-METHYLETHOXY)-1-METHYLETHOXY]-
 7 9227 1,3,3-TRIMETHOXYBUTANE
 8 6083 2-PROPANOL, 1,1'-OXYBIS-
 9 9228 1-PROPANOL, 2-(2-METHOXYPROPOXY)-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C7.H16.O3	148	59	345	800	348
2	C9.H18.O3	174	59	344	745	361
3	C9.H20.O4	192	59	344	806	355
4	C7.H16.O2	132	43	344	770	367
5	C9.H18.O7	238	58	340	661	388
6	C10.H22.O4	206	59	339	724	383
7	C7.H16.O3	148	45	325	689	340
8	C6.H14.O3	134	59	319	905	324
9	C7.H16.O3	148	59	318	800	328

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	55956-21-3
2	—	—	—	—	55956-25-7
3	—	—	—	—	1638-16-0
4	—	—	—	—	53907-95-2
5	—	—	—	—	- -
6	—	—	—	—	20324-33-8
7	—	—	—	—	6607-66-5
8	—	—	—	—	110-98-5
9	—	—	—	—	13588-28-8

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 18:29
 SAMPLE: S-MM5-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100%/100% *(NA/NA) /1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068108 #1661
 CALL: 30068108 # 3
 BASE M/Z: 59
 RIC: 5184.



Library Search Data: 30068108 #1688 Base m/z: 41
 08/31/98 22:29:00 + 18:47 Cali: 30068108 # 3 RIC: 2660.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 259 matched at least 5 of the 16 largest peaks in the unknown

Rank In. Name
 1 37446 3-OCTADECANONE
 2 54746 1,3-DIOXOLANE, 4-[[[2-METHOXY-4-OCTADECENYL)OXY]METHYL]-2,2-DIMETHY*
 3 33330 1,3:2,5:4,6-TRIMETHYLENE-D-GLYCERO-D-MANNOHEPTITOL
 4 19529 4-DODECANOL
 5 19875 2-BUTANOL, 3-(1,3,3-TRIMETHYLBUTOXY)-
 6 33472 PHOSPHINOUS CHLORIDE, (1-METHYLETHYL)[METHYL(1-METHYLETHYL)CYCLOHEX*
 7 5555 2,2,4-TRIMETHYL-3-PENTANOL
 8 29673 3-DODECANOL, 3,7,11-TRIMETHYL-
 9 5557 3-HEXANOL, 2,2-DIMETHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C18.H36.O	268	57	232	687	283
2	C25.H48.O4	412	71	205	692	268
3	C10.H16.O7	248	45	203	585	319
4	C12.H26.O	186	55	193	694	226
5	C11.H24.O2	188	57	192	715	219
6	C13.H26.CL.P	248	83	188	535	260
7	C8.H18.O	130	73	185	792	212
8	C15.H32.O	228	73	181	595	214
9	C8.H18.O	130	55	179	717	234

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	18261-92-2
2	—	—	—	—	16725-41-0
3	—	—	—	—	- -
4	—	—	—	—	10203-32-4
5	—	—	—	—	74810-44-9
6	—	—	—	—	74710-02-4
7	—	—	—	—	5162-48-1
8	—	—	—	—	7278-65-1
9	—	—	—	—	4209-90-9

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 22:29:00 + 18:47

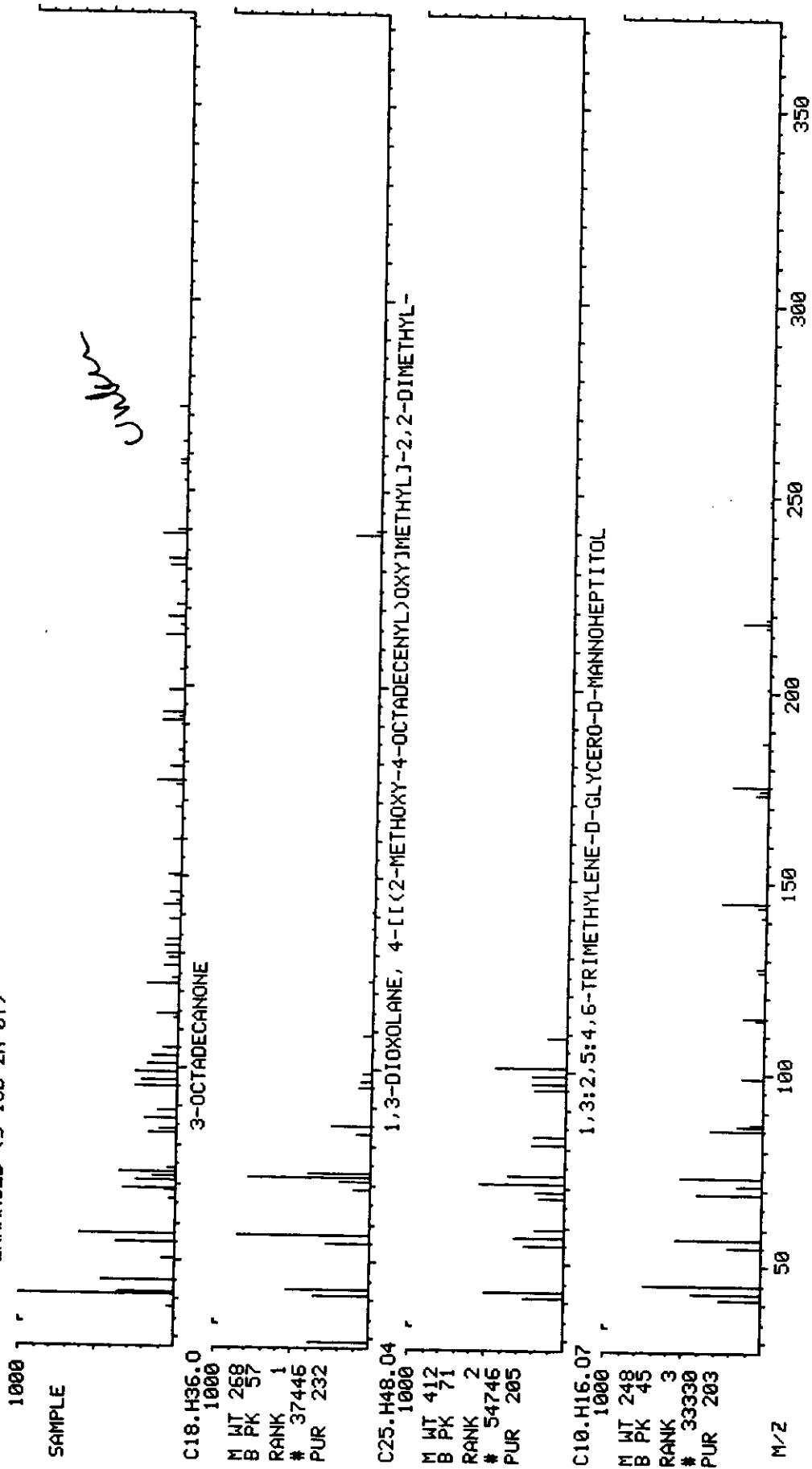
SAMPLE: 5-MMS-FB 1/35A/1ML

CONDS.: UG/ML *1ML *100%/100% *(NA/NA) >1/35A NA M

ENHANCED (5 15B 2N 0T)

DATA: 30068108 #1688
CALI: 30068108 # 3

BASE M/Z: 41
RIC: 2650.



Library Search Data: 30068108 #1703 Base m/z: 57
 08/31/98 22:29:00 + 18:57 Cali: 30068108 # 3 RIC: 7064.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 538 matched at least 7 of the 16 largest peaks in the unknown

Rank	In.	Name
1	37465	NONADECANE
2	46161	TRICOSANE
3	25997	PENTADECANE
4	32059	HEPTADECANE
5	29263	HEXADECANE
6	42196	HEPTADECANE, 2,6,10,14-TETRAMETHYL-
7	22530	TETRADECANE
8	29249	TETRADECANE, 6,9-DIMETHYL-
9	34810	OCTADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	723	941	753
2	C23.H48	324	43	712	901	776
3	C15.H32	212	57	707	950	731
4	C17.H36	240	57	701	931	744
5	C16.H34	226	57	700	929	745
6	C21.H44	296	57	695	897	721
7	C14.H30	198	43	687	917	718
8	C16.H34	226	57	678	906	702
9	C18.H38	254	57	677	867	773

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	638-67-5
3	—	—	—	—	629-62-9
4	—	—	—	—	629-78-7
5	—	—	—	—	544-76-3
6	—	—	—	—	18344-37-1
7	—	—	—	—	629-59-4
8	—	—	—	—	55045-13-1
9	—	—	—	—	593-45-3

MID LIBRARY SEARCH (LIBRARYNB)

08/31/98 22:29:00 + 18:57

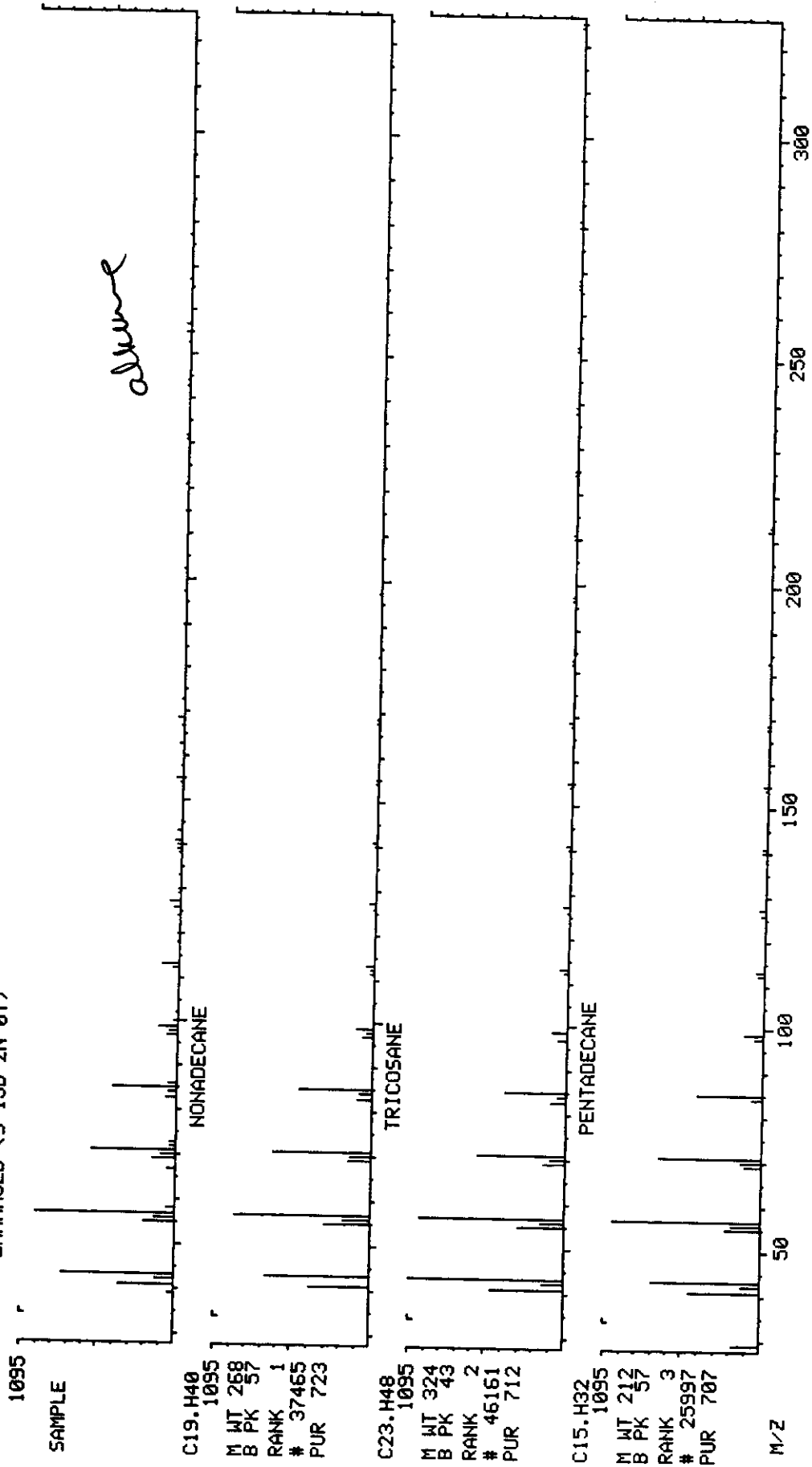
SAMPLE: S-MM5-FB 1/35A/1ML

CONDS.: UG/ML *1ML *100%/100% *(NA/NA) >1/35A NA M

ENHANCED (S 15B 2N 0T)

DATA: 30058108 #1703
CALI: 30058108 # 3

BASE M/Z: 57
RIC: 7064.



Library Search Data: 30068108 #1741 Base m/z: 59
 08/31/98 22:29:00 + 19:22 Cali: 30068108 # 3 RIC: 19520.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 483 matched at least 6 of the 16 largest peaks in the unknown

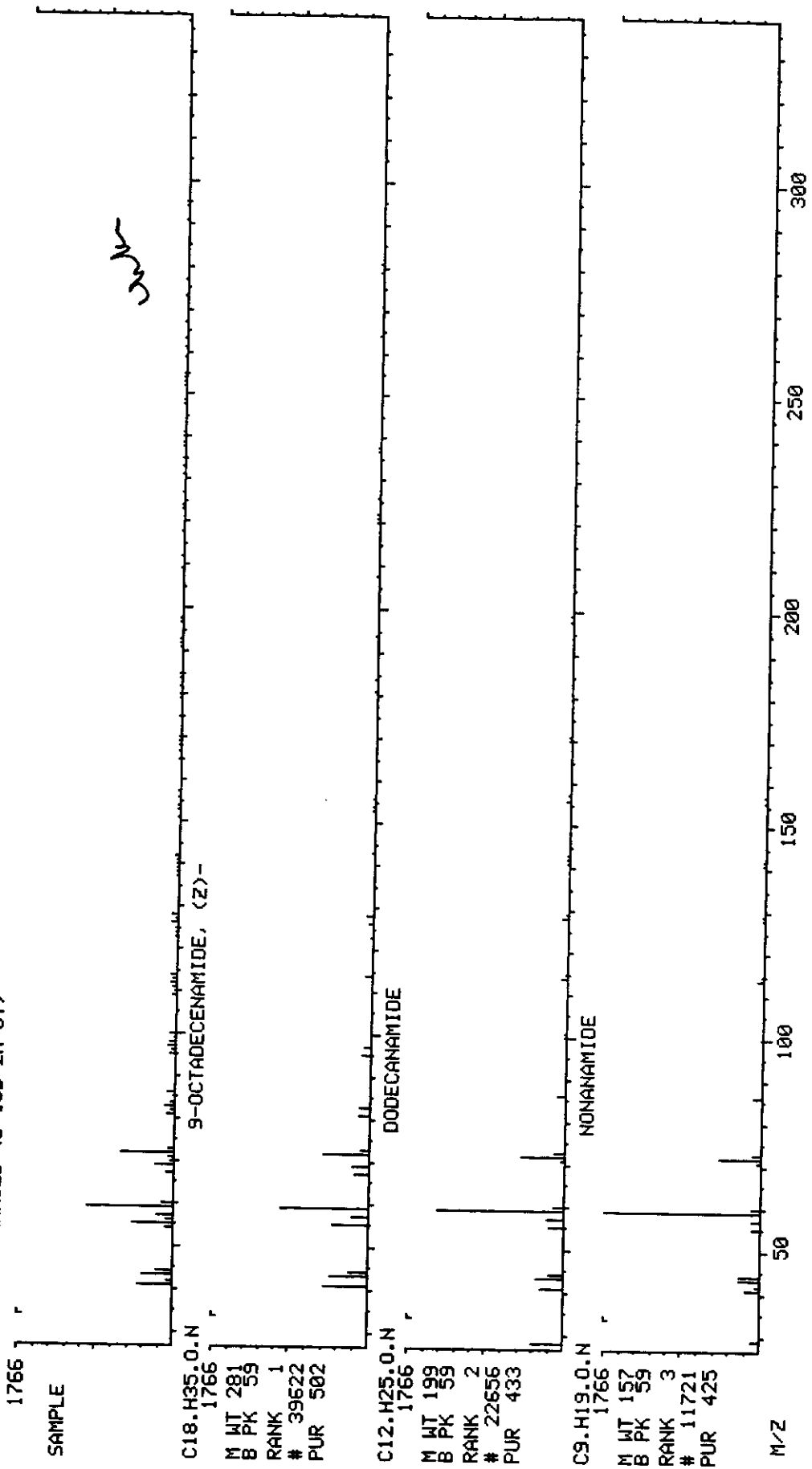
Rank In. Name
 1 39622 9-OCTADECENAMIDE, (Z)-
 2 22656 DODECANAMIDE
 3 11721 NONANAMIDE
 4 19516 1-DECANOL, 5,9-DIMETHYL-
 5 34956 HEXADECANAMIDE
 6 15482 HEPTANAMIDE, 4-ETHYL-5-METHYL-
 7 7998 TRANS-3,4-EPOXYNONANE
 8 3133 PENTANAMIDE, 4-METHYL-
 9 1617 PENTANAMIDE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C18.H35.O.N	281	59	502	895	556
2	C12.H25.O.N	199	59	433	896	461
3	C9.H19.O.N	157	59	425	882	428
4	C12.H26.O	186	41	415	819	419
5	C16.H33.O.N	255	59	406	846	439
6	C10.H21.O.N	171	59	406	865	419
7	C9.H18.O	142	41	397	807	414
8	C6.H13.O.N	115	59	392	910	397
9	C5.H11.O.N	101	59	363	872	363

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	301-02-0
2	—	—	—	—	1120-16-7
3	—	—	—	—	1120-07-6
4	—	—	—	—	91482-38-1
5	—	—	—	—	629-54-9
6	—	—	—	—	54789-40-1
7	—	—	—	—	- -
8	—	—	—	—	1119-29-5
9	—	—	—	—	626-97-1

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 19:22
 SAMPLE: S-MM5-FB 1/35A/1ML
 COND5.: UG/ML *1ML *100%/100% *(NA/NA) 1/35A NA M
 ENHANCED (S 158 2N 0T)

DATA: 30068108 #1741
 CALI: 30068108 # 3
 BASE M/Z: 59
 RIC: 19520.



Library Search Data: 30068108 #1809 Base m/z: 57
 08/31/98 22:29:00 + 20:07 Cali: 30068108 # 3 RIC: 9184.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 249 matched at least 7 of the 16 largest peaks in the unknown

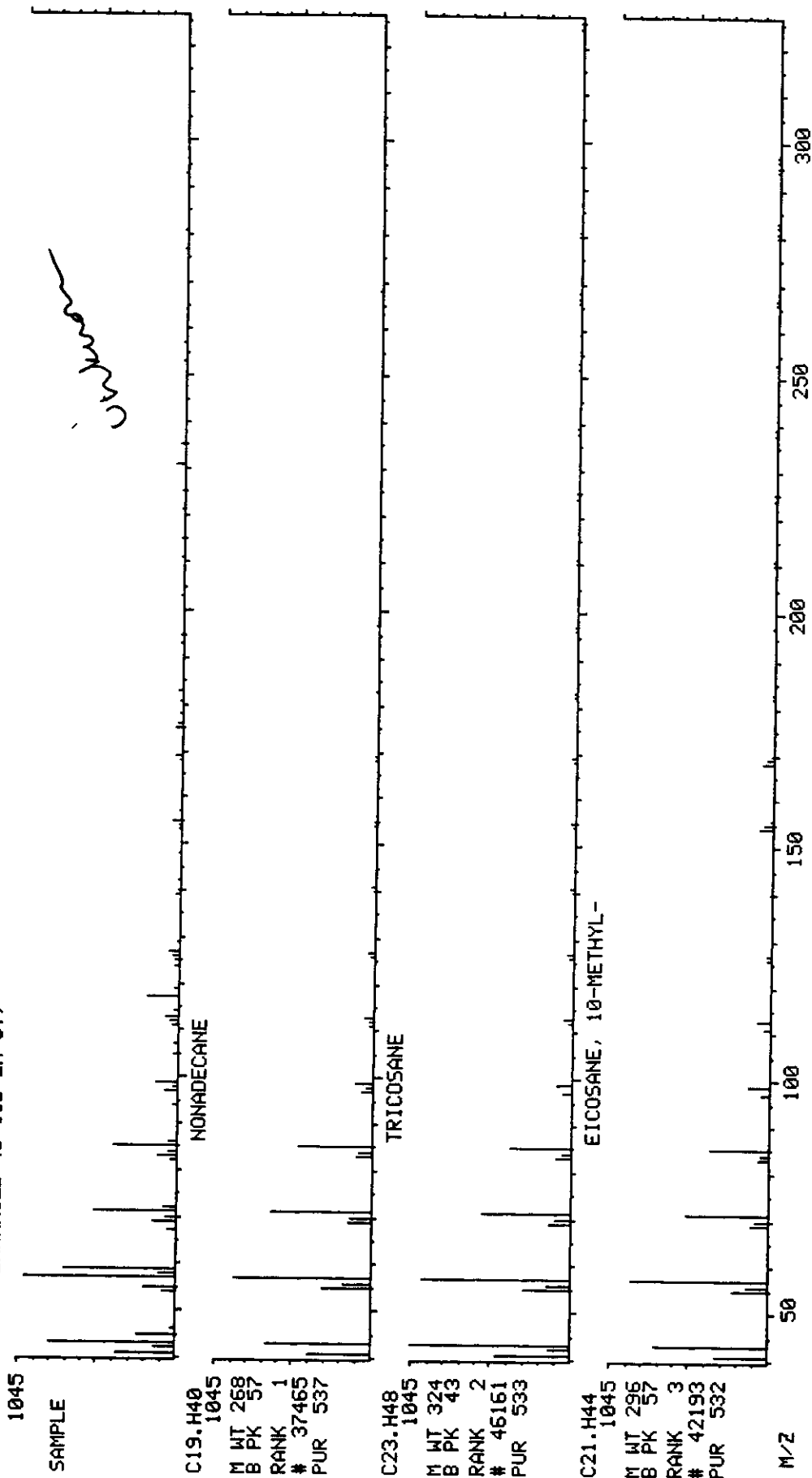
Rank In. Name
 1 37465 NONADECANE
 2 46161 TRICOSANE
 3 42193 EICOSANE, 10-METHYL-
 4 49555 PENTACOSANE
 5 37462 HEPTADECANE, 2,6-DIMETHYL-
 6 32059 HEPTADECANE
 7 44314 DOCOSANE
 8 37456 2-METHYLOCTADECANE
 9 25997 PENTADECANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C19.H40	268	57	537	904	576
2	C23.H48	324	43	533	875	595
3	C21.H44	296	57	532	898	561
4	C25.H52	352	43	517	856	592
5	C19.H40	268	57	517	872	557
6	C17.H36	240	57	510	897	555
7	C22.H46	310	57	510	866	578
8	C19.H40	268	43	507	891	551
9	C15.H32	212	57	506	899	546

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	629-92-5
2	—	—	—	—	638-67-5
3	—	—	—	—	54833-23-7
4	—	—	—	—	629-99-2
5	—	—	—	—	54105-67-8
6	—	—	—	—	629-78-7
7	—	—	—	—	629-97-0
8	—	—	—	—	- -
9	—	—	—	—	629-62-9

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 20:07
 SAMPLE: S-MM5-FB 1/35A/1ML
 CONDS.: UC/ML *1ML *1002/1002 *(NA/NA >1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30058108 #1809
 CALI: 30058108 # 3
 BASE M/Z: 57
 RIC: 9184.



1230

Library Search Data: 30068108 #1917 Base m/z: 43
 08/31/98 22:29:00 + 21:19 Cali: 30068108 # 3 RIC: 15024.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 818 matched at least 7 of the 16 largest peaks in the unknown

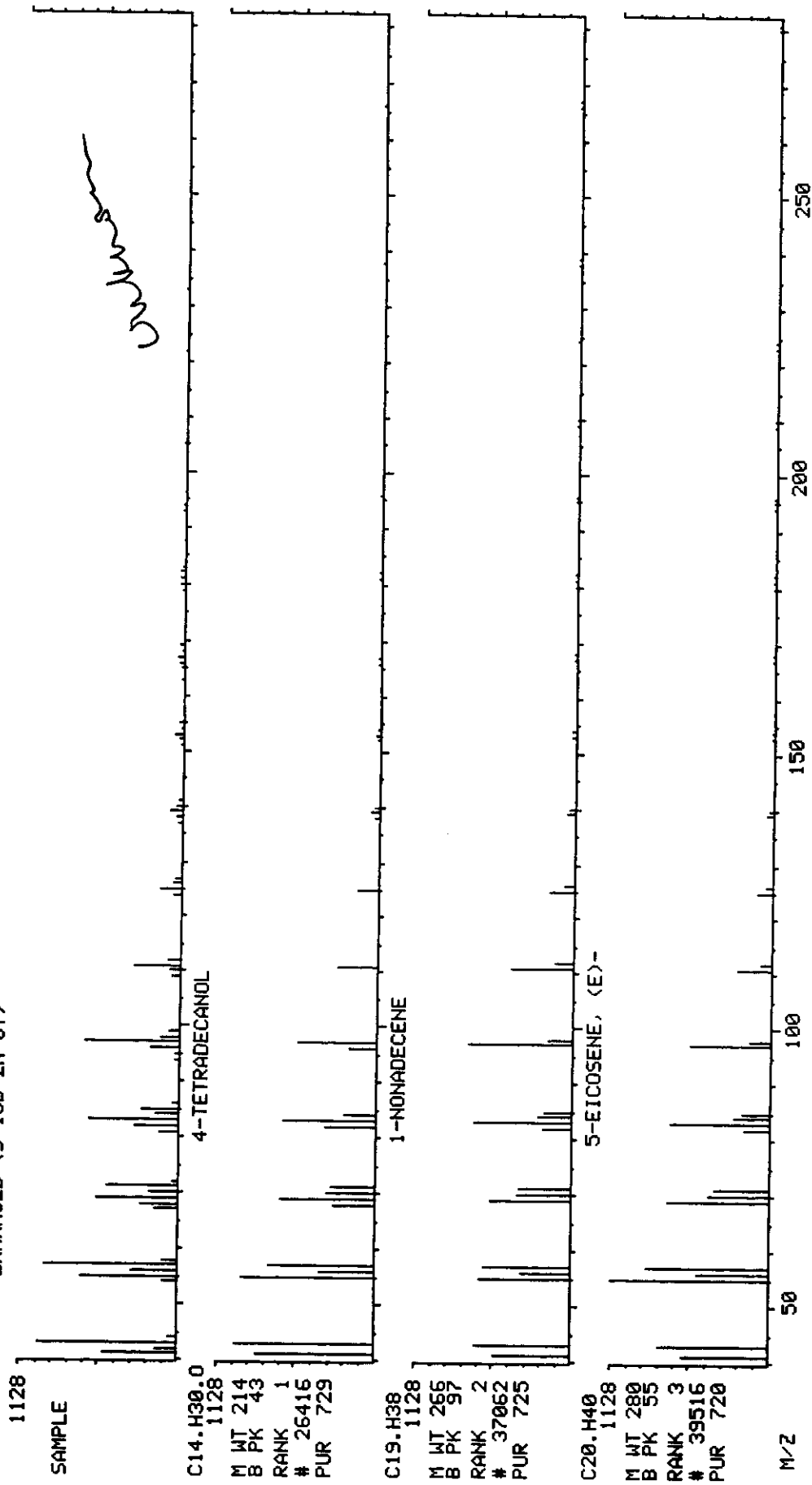
Rank In. Name
 1 26416 4-TETRADECANOL
 2 37062 1-NONADECENE
 3 39516 5-EICOSENE, (E)-
 4 34410 3-OCTADECENE, (E)-
 5 39515 9-EICOSENE, (E)-
 6 31653 1-HEPTADECENE
 7 34418 9-OCTADECENE, (E)-
 8 34411 5-OCTADECENE, (E)-
 9 48203 1-HENEICOSYL FORMATE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C14.H30.O	214	43	729	966	736
2	C19.H38	266	97	725	951	741
3	C20.H40	280	55	720	942	733
4	C18.H36	252	69	720	927	734
5	C20.H40	280	55	718	947	730
6	C17.H34	238	55	714	937	726
7	C18.H36	252	55	714	925	727
8	C18.H36	252	55	712	923	726
9	C22.H44.O2	340	57	711	937	749

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	1653-33-4
2	—	—	—	—	18435-45-5
3	—	—	—	—	74685-30-6
4	—	—	—	—	7206-19-1
5	—	—	—	—	74685-29-3
6	—	—	—	—	6765-39-5
7	—	—	—	—	7206-25-9
8	—	—	—	—	7206-21-5
9	—	—	—	—	77899-03-7

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 21:19
 SAMPLE: 5-MMS-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068108 #1917
 CALI: 30068108 # 3
 BASE M/Z: 43
 RIC: 15024.



1240

Library Search Data: 30068108 #1934 Base m/z: 45
 08/31/98 22:29:00 + 21:31 Cali: 30068108 # 3 RIC: 4880.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 226 matched at least 6 of the 16 largest peaks in the unknown

Rank In. Name
 1 9227 1,3,3-TRIMETHOXYBUTANE
 2 24247 2-PROPANOL, 1-[2-(2-METHOXY-1-METHYLETHOXY)-1-METHYLETHOXY]-
 3 16192 2-PROPANOL, 1-[1-METHYL-2-(2-PROPENYLOXY)ETHOXY]-
 4 9235 1-PROPANOL, 2-(2-METHOXY-1-METHYLETHOXY)-
 5 3528 2-METHYL-2,3-PENTANEDIOL
 6 36510 1,4,7,10,13,16-HEXAOXACYCLOOCTADECANE
 7 9228 1-PROPANOL, 2-(2-METHOXYPROPOXY)-
 8 5809 1-ETHOXYPENTAN-3-OL
 9 5810 2-PENTANOL, 5-METHOXY-2-METHYL-

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C7.H16.O3	148	45	399	722	437
2	C10.H22.O4	206	59	394	725	434
3	C9.H18.O3	174	59	380	693	394
4	C7.H16.O3	148	59	375	794	382
5	C6.H14.O2	118	59	360	824	374
6	C12.H24.O6	264	45	357	653	422
7	C7.H16.O3	148	59	353	767	365
8	C7.H16.O2	132	59	341	813	381
9	C7.H16.O2	132	59	334	728	383

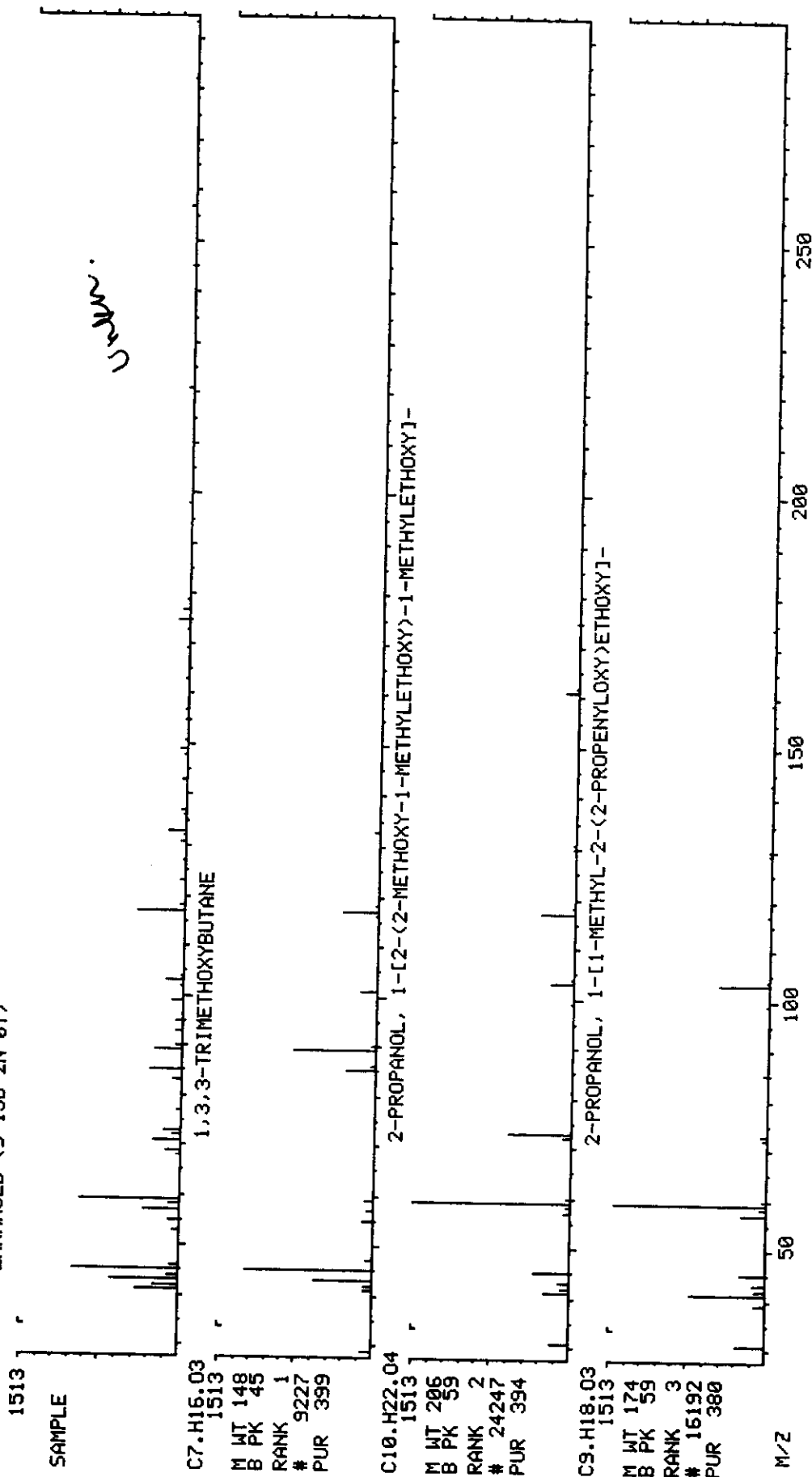
Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	6607-66-5
2	—	—	—	—	20324-33-8
3	—	—	—	—	55956-25-7
4	—	—	—	—	55956-21-3
5	—	—	—	—	7795-80-4
6	—	—	—	—	17455-13-9
7	—	—	—	—	13588-28-8
8	—	—	—	—	- -
9	—	—	—	—	55724-04-4

MID LIBRARY SEARCH (LIBRARYNB)
 08/31/98 22:29:00 + 21:31
 SAMPLE: S-MMS-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100% *(NA/NA) /1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30068108 #1934
 CALI: 30068108 # 3

BASE M/Z: 45
 RIC: 4890.

INST. ID: F16



Library Search Data: 30068108 #2016 Base m/z: 57
 08/31/98 22:29:00 + 22:25 Cali: 30068108 # 3 RIC: 5768.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 158 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 818 matched at least 7 of the 16 largest peaks in the unknown

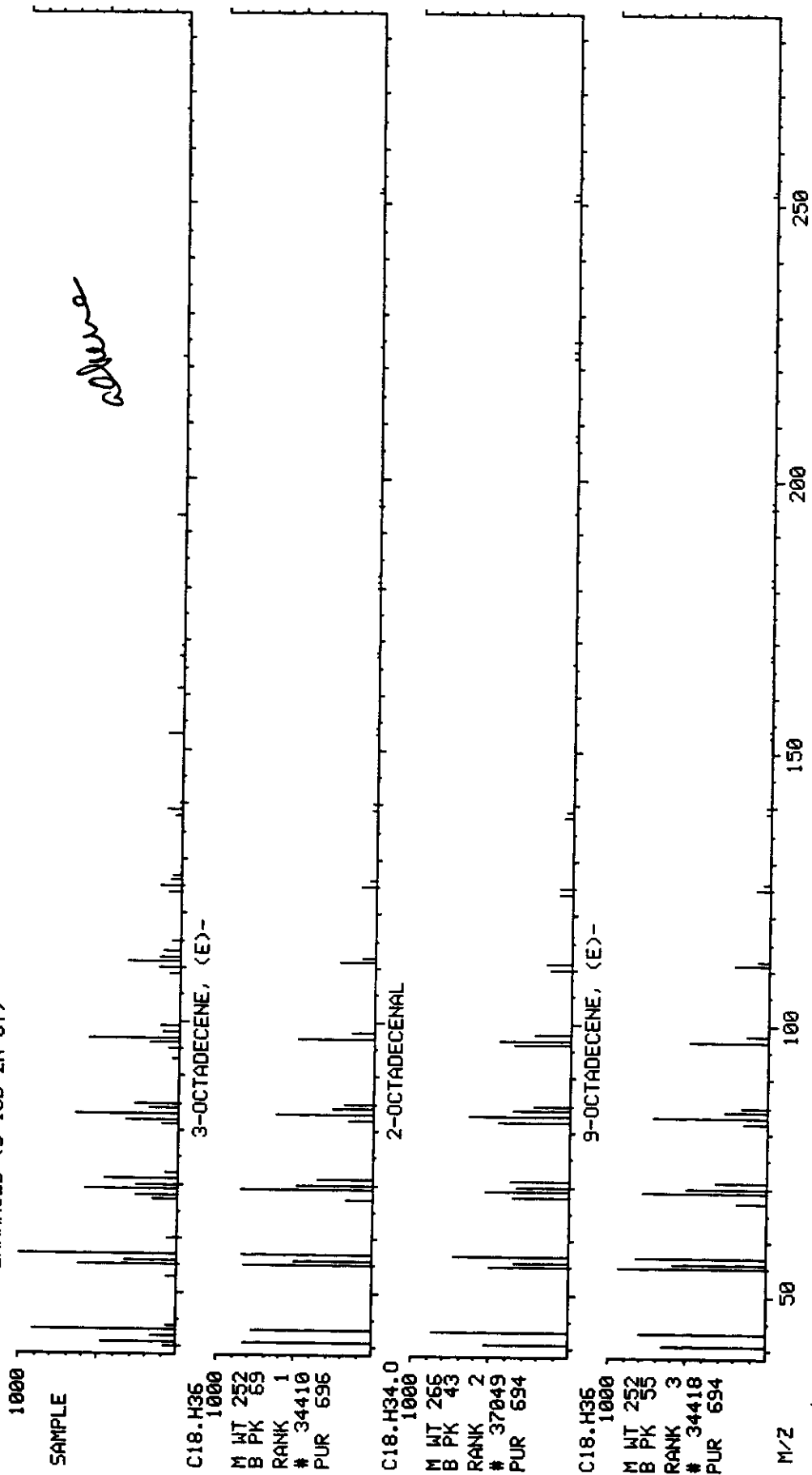
Rank In. Name
 1 34410 3-OCTADECENE, (E)-
 2 37049 2-OCTADECENAL
 3 34418 9-OCTADECENE, (E)-
 4 39515 9-EICOSENE, (E)-
 5 34411 5-OCTADECENE, (E)-
 6 37062 1-NONADECENE
 7 39516 5-EICOSENE, (E)-
 8 40232 1-NONADECANOL
 9 58701 17-PENTATRIACONTENE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C18.H36	252	69	696	927	716
2	C18.H34.O	266	43	694	900	747
3	C18.H36	252	55	694	925	712
4	C20.H40	280	55	691	940	712
5	C18.H36	252	55	691	921	709
6	C19.H38	266	97	690	932	720
7	C20.H40	280	55	688	934	710
8	C19.H40.O	284	55	688	966	700
9	C35.H70	490	43	683	905	738

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	7206-19-1
2	—	—	—	—	56554-96-2
3	—	—	—	—	7206-25-9
4	—	—	—	—	74685-29-3
5	—	—	—	—	7206-21-5
6	—	—	—	—	18435-45-5
7	—	—	—	—	74685-30-6
8	—	—	—	—	1454-84-8
9	—	—	—	—	6971-40-0

MID LIBRARY SEARCH <LIBRARYNB>
 08/31/98 22:29:00 + 22:25
 SAMPLE: S-MM5-FB 1/35A/1ML
 CONDS.: UG/ML *1ML *100Z/100Z *(NA/NA)/1/35A NA M
 ENHANCED (S 15B 2N 0T)

DATA: 30058108 #2016
 CALI: 30058108 # 3
 BASE M/Z: 57
 RIC: 5768.



1244

Library Search Data: 30068108 #2052 Base m/z: 59
 08/31/98 22:29:00 + 22:49 Cali: 30068108 # 3 RIC: 3200.
 Sample: S-MM5-FB 1/3SA/1ML INST. ID: F16
 Conds.: UG/ML *1ML *100%/100% *(NA/NA)/1/3SA NA M
 Enhanced (S 15B 2N 0T)

62231 spectra in LIBRARYNB searched for maximum PURITY
 214 matched at least 5 of the 16 largest peaks in the unknown

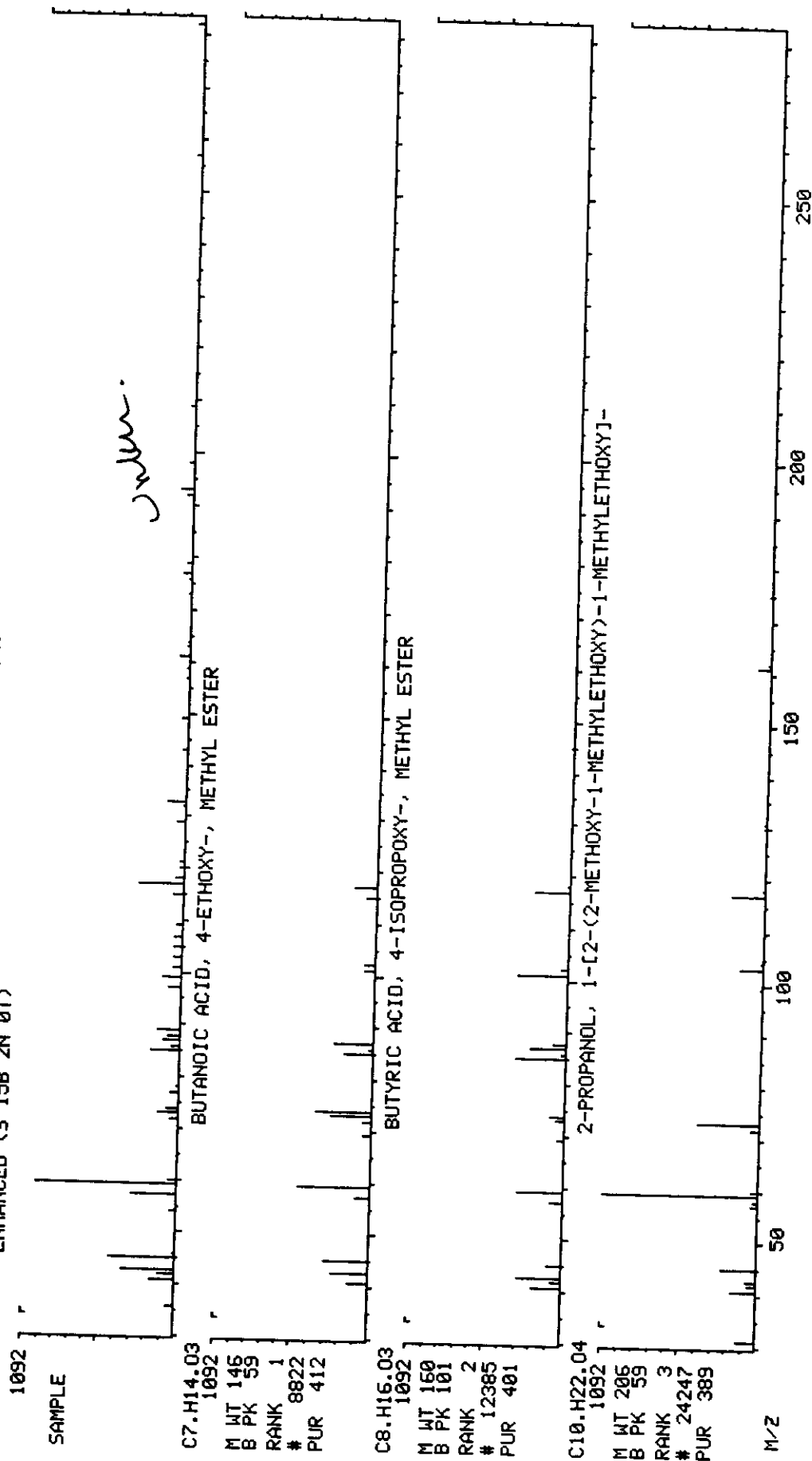
Rank In. Name
 1 8822 BUTANOIC ACID, 4-ETHOXY-, METHYL ESTER
 2 12385 BUTYRIC ACID, 4-ISOPROPOXY-, METHYL ESTER
 3 24247 2-PROPANOL, 1-[2-(2-METHOXY-1-METHYLETHOXY)-1-METHYLETHOXY]-
 4 36537 2,5,8,11-TETRAOXATETRADECAN-13-OL, 4,7,10-TRIMETHYL-
 5 16199 BUTYRIC ACID, 4-BUTOXY-, METHYL ESTER
 6 9228 1-PROPANOL, 2-(2-METHOXYPROPOXY)-
 7 16192 2-PROPANOL, 1-[1-METHYL-2-(2-PROPENYLOXY)ETHOXY]-
 8 27502 15-CROWN-5
 9 9227 1,3,3-TRIMETHOXYBUTANE

Rank	Formula	M.Wt	B.Pk	Purity	Fit	RFit
1	C7.H14.O3	146	59	412	839	435
2	C8.H16.O3	160	101	401	795	417
3	C10.H22.O4	206	59	389	823	405
4	C13.H28.O5	264	59	383	752	436
5	C9.H18.O3	174	57	361	748	398
6	C7.H16.O3	148	59	350	829	360
7	C9.H18.O3	174	59	335	664	352
8	C10.H20.O5	220	45	334	650	389
9	C7.H16.O3	148	45	333	673	351

Rank	Ret.Time	B.P.Int.	US.Par.1	US.Par.2	C.A.S. #
1	—	—	—	—	29006-04-0
2	—	—	—	—	29006-05-1
3	—	—	—	—	20324-33-8
4	—	—	—	—	20324-34-9
5	—	—	—	—	29006-06-2
6	—	—	—	—	13588-28-8
7	—	—	—	—	55956-25-7
8	—	—	—	—	33100-27-5
9	—	—	—	—	6607-66-5

MID LIBRARY SEARCH (LIBRARYNB)
08/31/98 22:29:00 + 22:49
SAMPLE: S-MMS-FB 1/35A/1ML
CONDS.: UG/ML *1ML *1002/1002 *(NA/NA) 1/35A NA M
ENHANCED (5 15B 2N 0T)

DATA: 30068108 #2052
CALI: 30068108 # 3
BASE M/Z: 59
RIC: 3200.



1246

[Handwritten signature]

TIC SELECTION REPORT

DATA FILE: 30068108

THE FOLLOWING PEAKS WERE REJECTED BECAUSE
AT LEAST 40 % OF THEIR SIZE WAS ACCOUNTED FOR BY
TARGET COMPOUNDS ELUTING WITHIN 4 SCANS OF THE
PEAK TOP.

SCAN	SIZE	AMOUNT
198	230232.	32.201
308	286752.	40.106
340	226092.	31.622
421	132760.	18.568
529	285996.	40.000
722	314784.	36.205
827	347784.	40.000
968	289424.	32.602
1086	355104.	40.000
1383	311936.	57.604
1383	311936.	57.604
1551	216608.	40.000
1603	32336.	5.971
1782	175064.	40.000

TOTAL NUMBER OF UNIDENTIFIED PEAKS WITH SIZE
GREATER THAN 10 % OF THE CLOSEST INTERNAL STANDARD
THAT DOES NOT HAVE INTERFERENCES = 32

INTERNAL STANDARDS THAT HAVE RIC SIZE LESS THAN
50 % OR GREATER THAN 200 % OF THE ESTIMATED RIC SIZE
ARE CONSIDERED TO HAVE INTERFERENCES AND WILL NOT BE USED
FOR QUANTITATION.

#	INTERNAL STANDARD	RIC SIZE	PERCENT OF ESTIMATED RIC SIZE	SCAN
1	C140 NAPHTHALENE-D8	285996.	129	529
2	C150 ACENAPHTHENE-D10	347784.	129	827
3	C160 PHENANTHRENE-D10	355104.	128	1086
4	C170 CHRYSENE-D12	216608.	134	1551
5	C175 PERYLENE-D12	175064.	123	1782

* INDICATES INTERFERENCE

SIZE = AREA

TECHNICAL REPORT DATA Please read instructions on the reverse before completing		
1. REPORT NO. EPA-454/R-00-025E	2.	3. RECIPIENT'S ACCESSION NO.
4. TITLE AND SUBTITLE Final Report Hot Mix Asphalt Plants, Truck Loading and Silo Filling, Manual Methods Testing, Asphalt Plant C, Los Angeles, California Volume 5 of 8	5. REPORT DATE May 2000	
	6. PERFORMING ORGANIZATION CODE	
7. AUTHOR(S) Frank J. Phoenix	8. PERFORMING ORGANIZATION REPORT NO.	
9. PERFORMING ORGANIZATION NAME AND ADDRESS Pacific Environmental Services, Inc. Post Office Box 12077 Research Triangle Park, North Carolina 27709-2077	10. PROGRAM ELEMENT NO.	
	11. CONTRACT/GRANT NO. 68-D-98004	
12. SPONSORING AGENCY NAME AND ADDRESS U.S. Environmental Protection Agency Office of Air Quality Planning and Standards Emissions, Monitoring and Analysis Division Research Triangle Park, North Carolina 27711	13. TYPE OF REPORT AND PERIOD COVERED Final	
	14. SPONSORING AGENCY CODE EPA/200/04	
15. SUPPLEMENTARY NOTES		
16. ABSTRACT The United States Environmental Protection Agency (EPA) Office of Air Quality Planning and Standards (OAQPS) is investigating hot mix asphalt plants to identify and quantify particulate matter (PM), methylene chloride extractable matter (MCEM), and organic hazardous air pollutant (HAP) emissions during asphalt concrete loading operations. In support of this investigation, the OAQPS issued Pacific Environmental Services, Inc. (PES) a series of work assignments to conduct emissions testing at a hot mix asphalt plant during load-out operations. The primary objective of the emissions testing was to characterize the uncontrolled emissions of PM, MCEM, polynuclear aromatic hydrocarbons (PAHs), semi-volatile organic hazardous air pollutants (SVOHAPS), and volatile organic hazardous air pollutants (VOHAPS) from a hot mix production plant during loading operations. An asphalt plant south of Los Angeles, California was selected by EPA as the host facility. Testing was performed over five consecutive days beginning on July 24, 1998. Testing was performed under two conditions. Under normal operations, testing was performed to characterize load-out emissions from the tunnel exhaust and load-in emissions from the asphalt concrete storage silo. Under background conditions, testing was performed to characterize emissions from the combustion of diesel fuel in transport trucks. The entire report consists of eight volumes totaling 4,234 pages, Vol. 1 (388 pages), Vol. 2 (308 pages), Vol. 3 (573 pages), Vol. 4 (694 pages), Vol. 5 (606 pages), Vol. 6 (564 pages), Vol. 7 (570 pages), and Vol. 8 (531 pages).		
17. KEY WORDS AND DOCUMENT ANALYSIS		
a. DESCRIPTIONS	b. IDENTIFIERS/OPEN ENDED TERMS	c. COASTI Field/Group
Hazardous Air Pollutants Methylene Chloride Extractable Matter Particulate Matter Polynuclear Aromatic Hydrocarbons Semi-volatile Organic Hazardous Air Pollutants Volatile Organic Hazardous Air Pollutants		
18. DISTRIBUTION STATEMENT Unlimited	19. SECURITY CLASS (<i>This Report</i>) Unclassified	21. NO. OF PAGES Vol. 5 - 606
	20. SECURITY CLASS (<i>This page</i>) Unclassified	22. PRICE