User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 03/28/2002 Time: 17:01:40

Sample: AE03508
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 3/28/02 17:01 Ended: 3/28/02 17:01 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		87		5.0

Comments for Sample AE03508 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 03-28-2002 Time: 17:02:35

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. The recoveries for 6 of the 43 compounds in the LCS Standard were high.

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: SPLIT.P
Quant Time: Mar 15 10:05:08 2002 Quant Results File: HP022102.RES

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Last Update : Mon Feb 25 10:47:59 2002
Response via : Initial Calibration
DataAcq Meth : HP022102

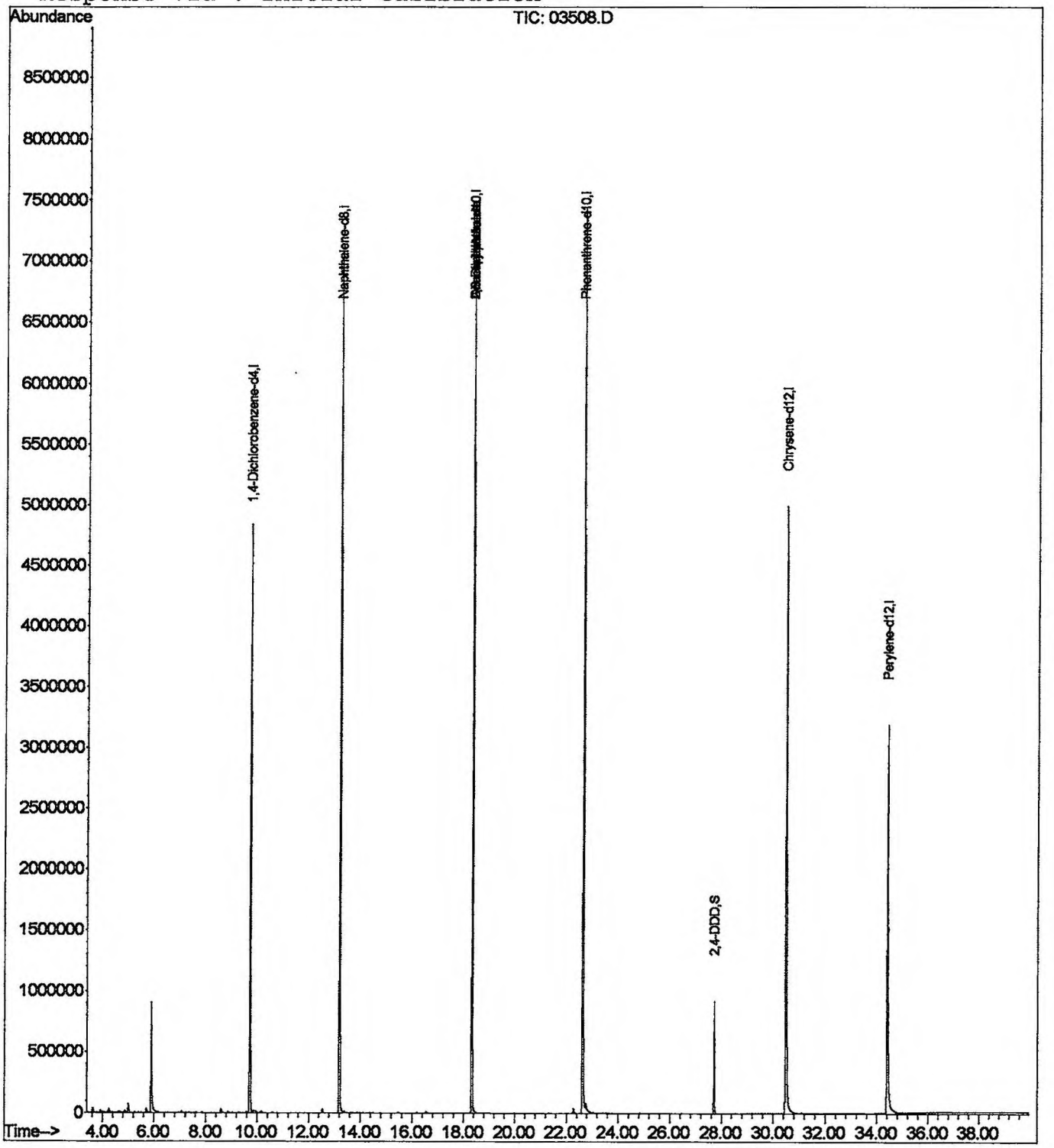
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.72	152	1694626	20.00	ug/L	0.00
10) Naphthalene-d8	13.20	136	7304283	20.00	ug/L	0.00
17) Acenaphthene-d10	18.33	164	3620508	20.00	ug/L	0.00
29) Phenanthrene-d10	22.63	188	6007536	20.00	ug/L	-0.01
38) Chrysene-d12	30.49	240	4278787	20.00	ug/L	-0.03
44) Perylene-d12	34.41	264	2450370	20.00	ug/L	-0.04
System Monitoring Compounds						
37) 2,4-DDD	27.72	235	347695	4.34	ug/L	0.00
Spiked Amount 5.000			Recovery = 86.80%			
Target Compounds						Qvalue
20) Dimethylphthalate	18.33	163	587338	2.67	ug/L	# 58
21) 2,6-Dinitrotoluene	18.33	165	465976	9.19	ug/L	# 27

(#) = qualifier out of range (m) = manual integration
03508.D HP022102.M Fri Mar 15 10:05:09 2002 Page 1

Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: SPLIT.P
Quant Time: Mar 15 10:05 2002 Quant Results File: HP022102.R

Method : C:\MSDCHEM\1\METHODS\METHODS\HP022102.M (RTE Integrator)
Title : 508 scan method
Last Update : Mon Feb 25 10:47:59 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\MSDCHEM\1\METHODS\METHODS\HP022102.M (RTE Integrator)
Title : 508 scan method
Smoothing : OFF Filtering: 5
Sampling : 2 Min Area: 100 Area counts
Start Thrs: 0.2 Max Peaks: 20
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

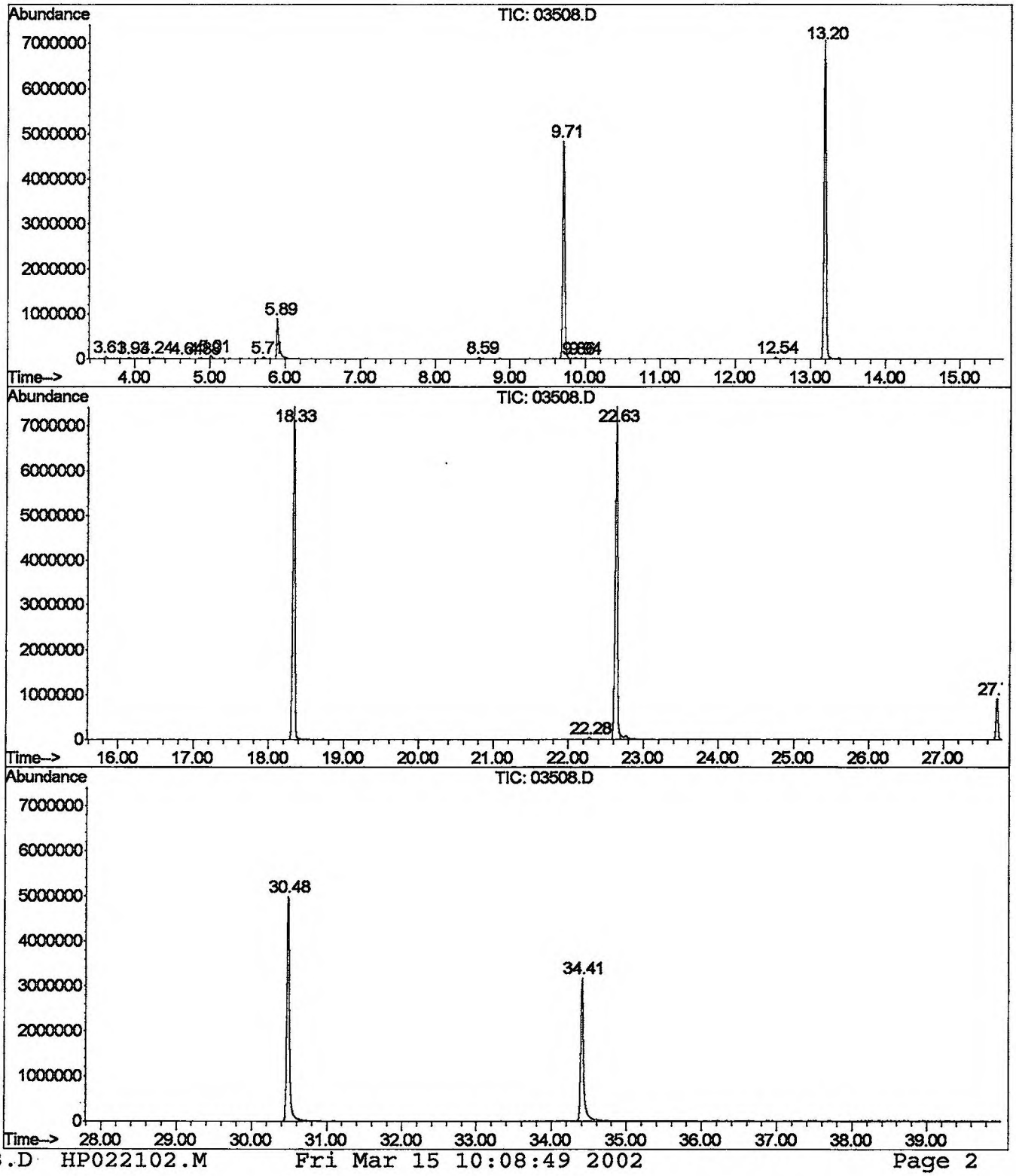
Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.611	19	21	29	rVB	41760	72967	0.47%	0.092%
2	3.927	45	49	53	rVB	21382	37878	0.24%	0.048%
3	4.243	75	77	85	rVV3	30861	70114	0.45%	0.088%
4	4.638	107	112	125	rBV4	10966	53206	0.34%	0.067%
5	4.875	129	133	143	rBV3	17690	62723	0.40%	0.079%
6	5.011	143	145	153	rVB	70715	127911	0.82%	0.161%
7	5.710	203	207	213	rVB	34701	61888	0.39%	0.078%
8	5.891	219	223	243	rBV	897500	1874101	11.96%	2.353%
9	8.589	457	462	465	rBV	34718	58135	0.37%	0.073%
10	9.707	557	561	567	rBV	4832674	9633318	61.47%	12.095%
11	9.865	573	575	579	rBV	16049	39077	0.25%	0.049%
12	9.944	579	582	591	rVB2	16491	50278	0.32%	0.063%
13	12.540	807	812	817	rVB	33686	59189	0.38%	0.074%
14	13.195	865	870	875	rBV	7073817	13944385	88.98%	17.508%
15	18.331	1319	1325	1331	rBV	7427508	14958021	95.45%	18.781%
16	22.283	1671	1675	1681	rBV	41317	84814	0.54%	0.106%
17	22.633	1701	1706	1711	rBV2	7427132	15671413	100.00%	19.677%
18	27.724	2153	2157	2161	rVV	916590	1710909	10.92%	2.148%
19	30.478	2395	2401	2435	rBV	4986565	12592213	80.35%	15.810%
20	34.407	2743	2749	2785	rBV	3180400	8482083	54.12%	10.650%

Sum of corrected areas: 79644623

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\5973N\02MAR08\03508.D
Operator :
Acquired : 8 Mar 2002 1:07 pm using AcqMethod HP022102
Instrument : Instrumen
Sample Name: 508 scan
Misc Info : March 8, 2002
Vial Number: 4
Quant File :HP022102.RES (RTE Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

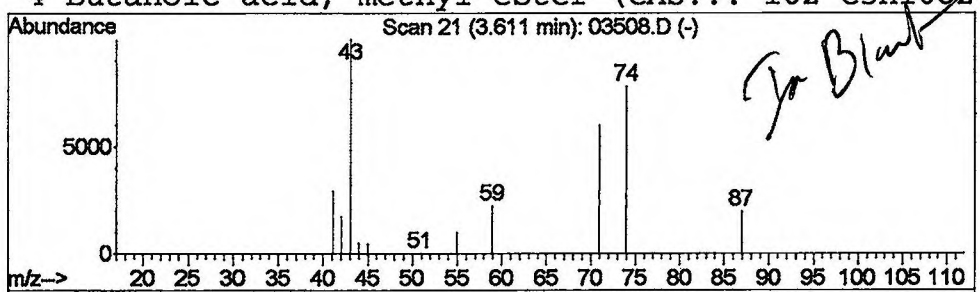
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

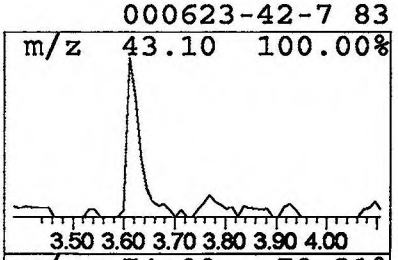
****

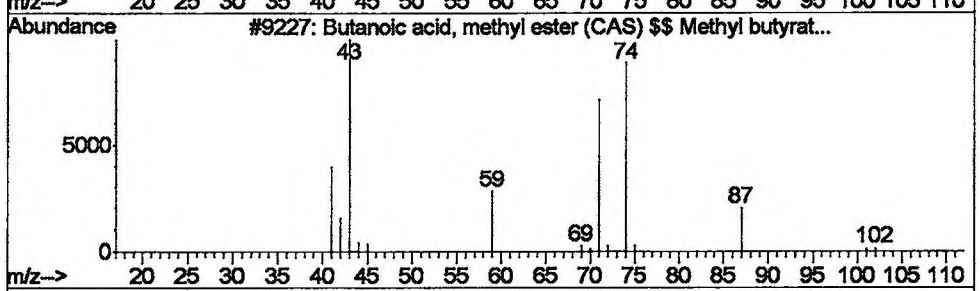
Peak Number 1 Butanoic acid, methyl ester... Concentration Rank 3

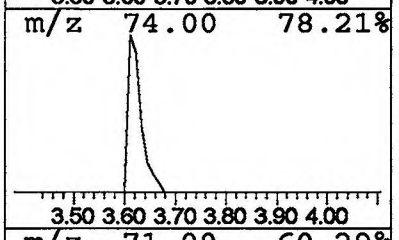
R.T.	EstConc	Area	Relative to ISTD	R.T.
3.61	0.15 ug/L	72967	1,4-Dichlorobenzene-d4	9.72

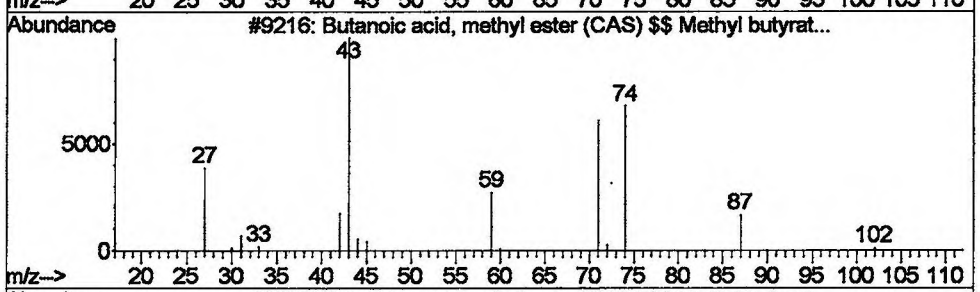
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	90
2	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83
3	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83
4	Butanoic acid, methyl ester (CAS...	102	C5H10O2	000623-42-7	83

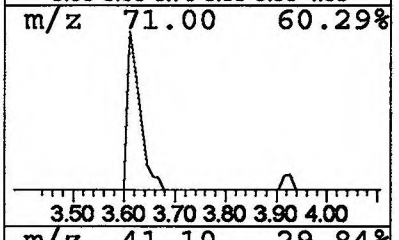


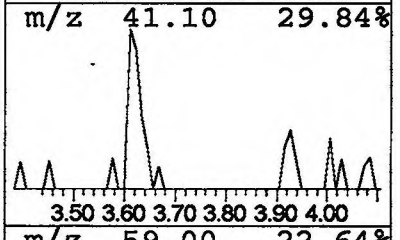


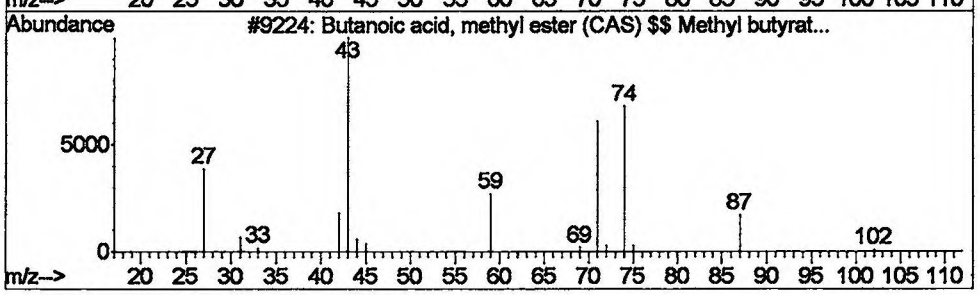


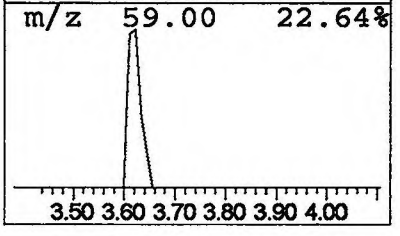












Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

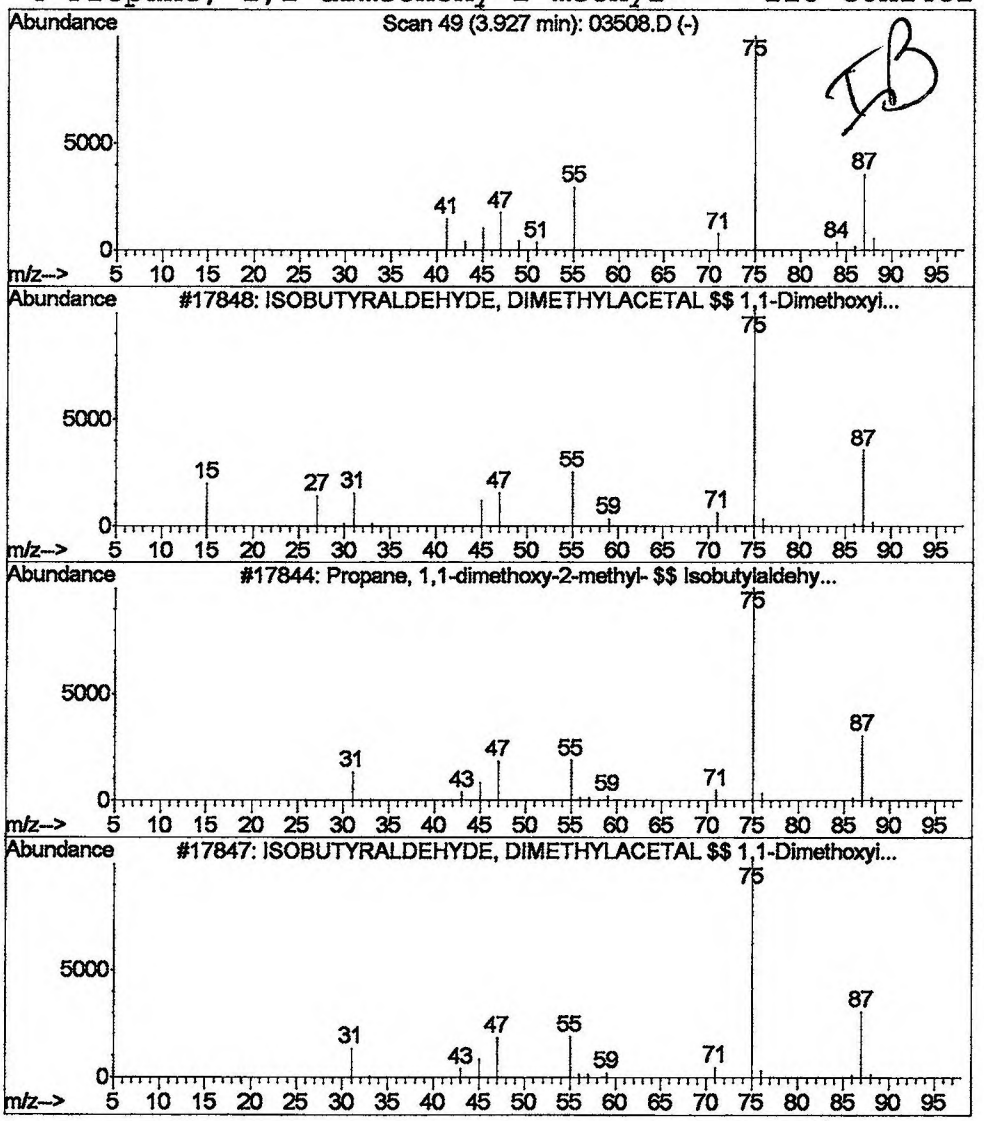
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

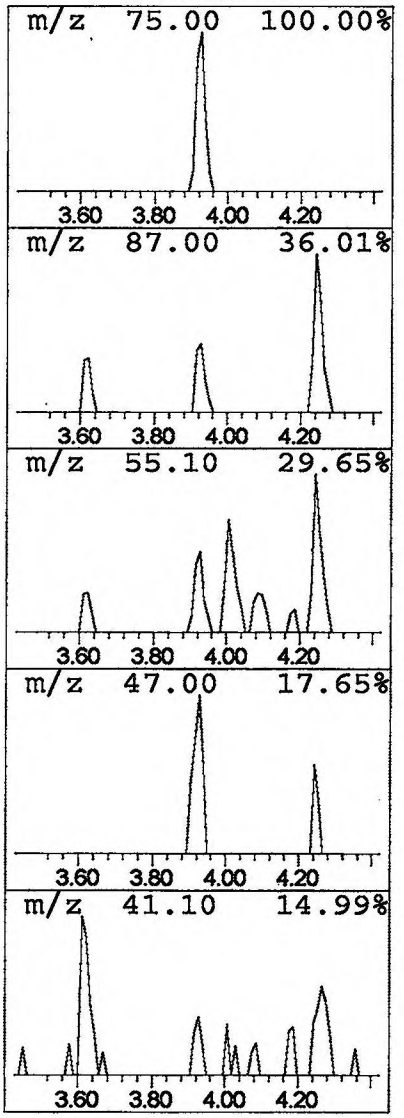
****

Peak Number 2 ISOBUTYRALDEHYDE, DIMETHYLA... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.93	0.08 ug/L	37878	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	78
2	Propane, 1,1-dimethoxy-2-methyl-...	118	C6H14O2	041632-89-7	78
3	ISOBUTYRALDEHYDE, DIMETHYLACETAL...	118	C6H14O2	000000-00-0	78
4	Propane, 1,1-dimethoxy-2-methyl-	118	C6H14O2	041632-89-7	64





Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

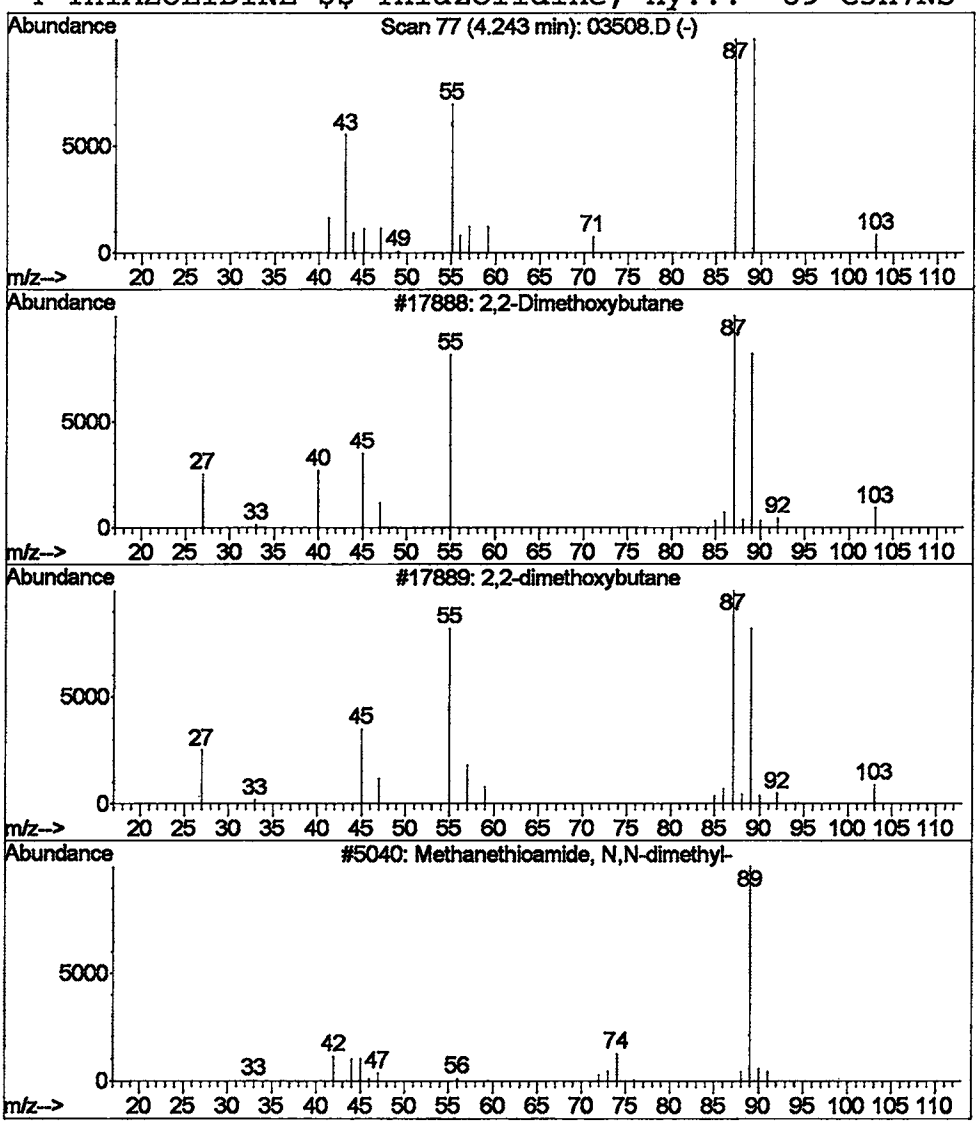
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

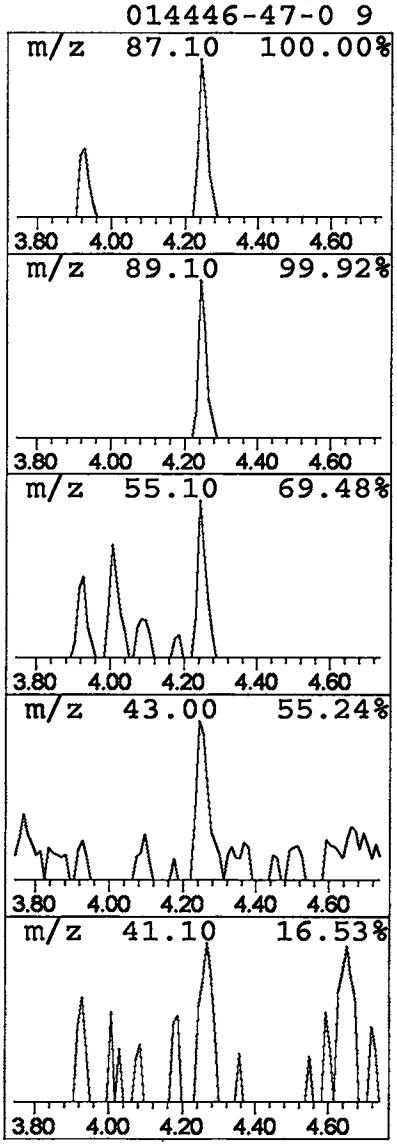
****

Peak Number 3 2,2-Dimethoxybutane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.24	0.15 ug/L	70114	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Dimethoxybutane	118	C6H14O2	003453-99-4	64
2	2,2-dimethoxybutane	118	C6H14O2	003453-99-4	64
3	Methanethioamide, N,N-dimethyl-	89	C3H7NS	000758-16-7	9
4	THIAZOLIDINE $$ Thiazolidine, hy...	89	C3H7NS	014446-47-0	9





Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

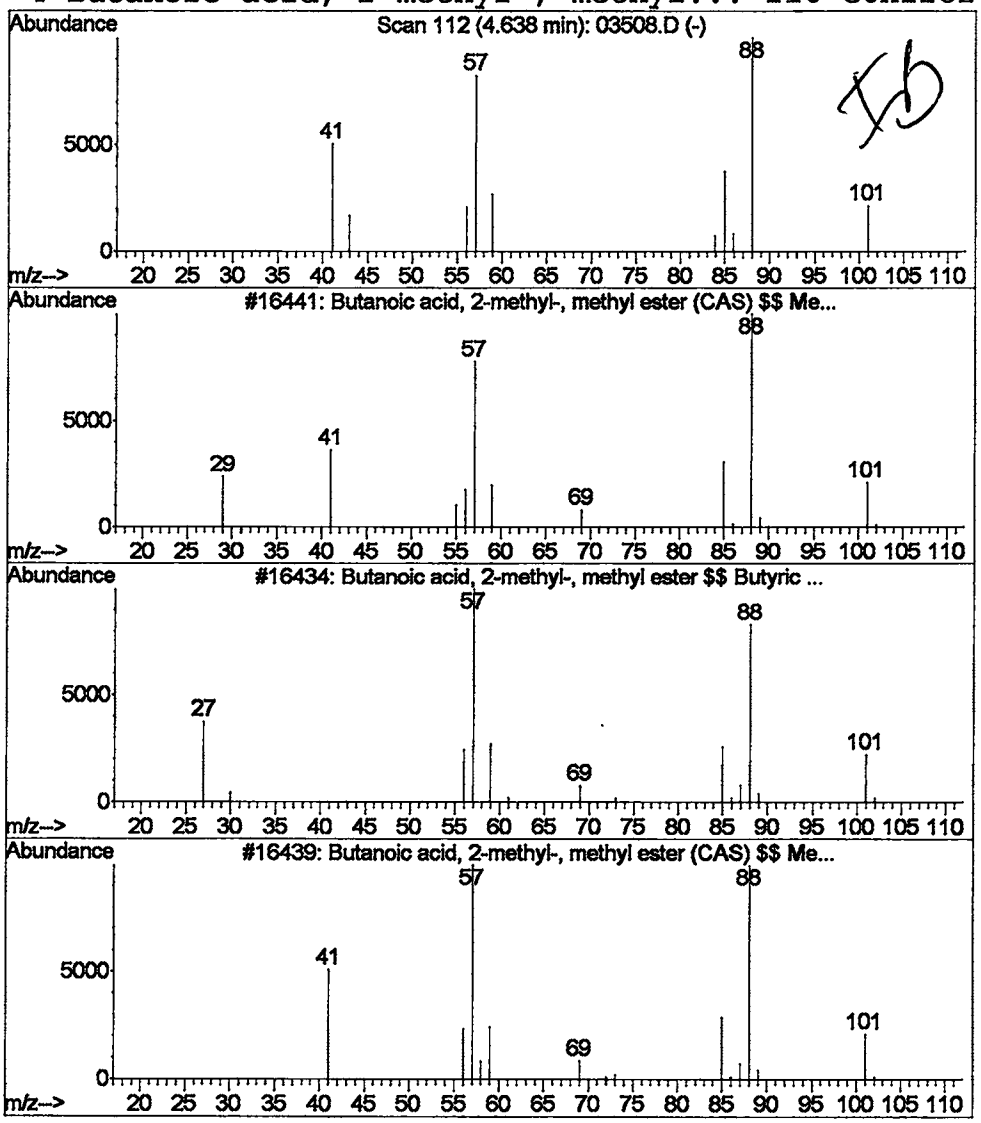
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

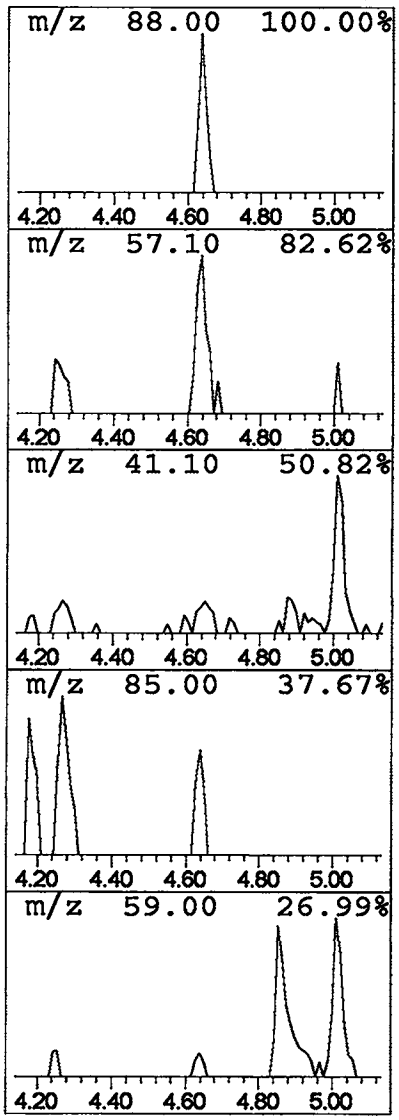
****

Peak Number 4 Butanoic acid, 2-methyl-, m... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.64	0.11 ug/L	53206	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 2-methyl-, methyl...	116	C6H12O2	000868-57-5	83
2	Butanoic acid, 2-methyl-, methyl...	116	C6H12O2	000868-57-5	74
3	Butanoic acid, 2-methyl-, methyl...	116	C6H12O2	000868-57-5	74
4	Butanoic acid, 2-methyl-, methyl...	116	C6H12O2	053955-81-0	64





Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

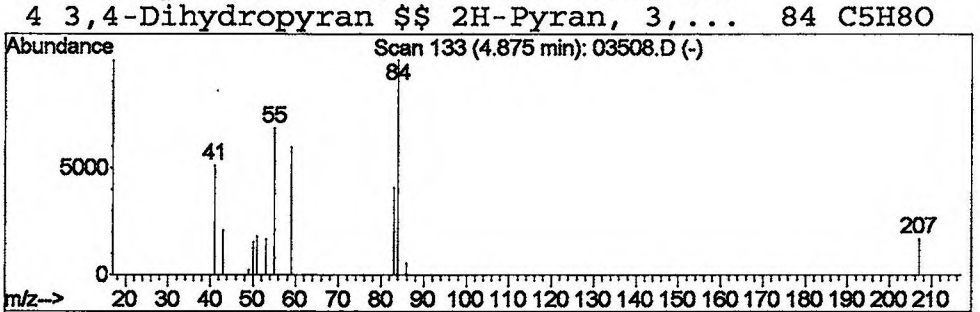
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

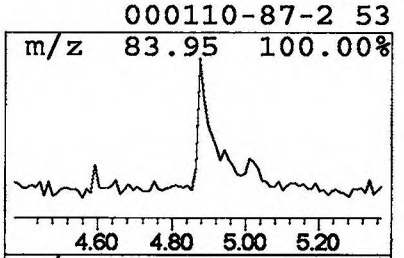
****

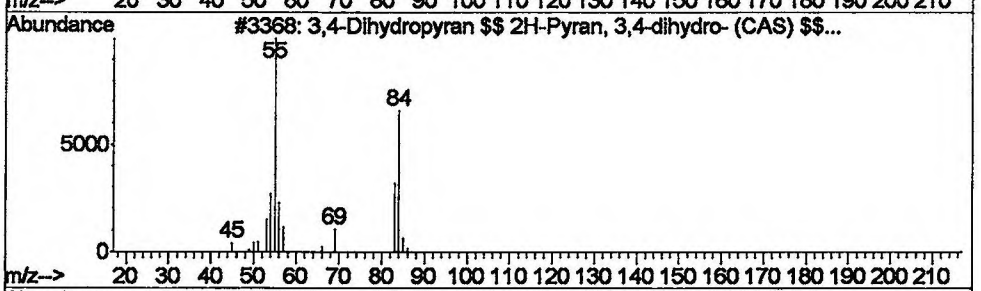
Peak Number 5 3,4-Dihydropyran $$ 2H-Pyra... Concentration Rank 5

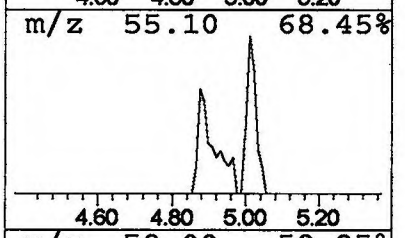
R.T.	EstConc	Area	Relative to ISTD	R.T.
4.88	0.13 ug/L	62723	1,4-Dichlorobenzene-d4	9.72

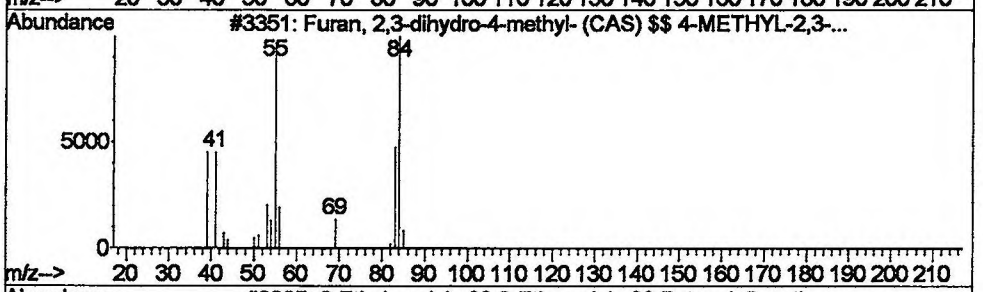
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,4-Dihydropyran $$ 2H-Pyran, 3,...	84	C5H8O	000110-87-2	58
2	Furan, 2,3-dihydro-4-methyl- (CA...	84	C5H8O	034314-83-5	53
3	2-Ethylacrolein $$ 2-Ethacrolein...	84	C5H8O	000922-63-4	53
4	3,4-Dihydropyran $$ 2H-Pyran, 3,...	84	C5H8O	000110-87-2	53

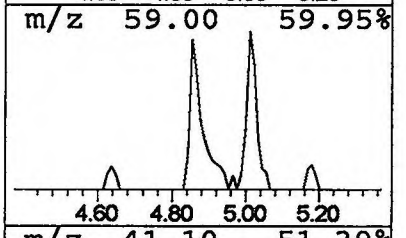


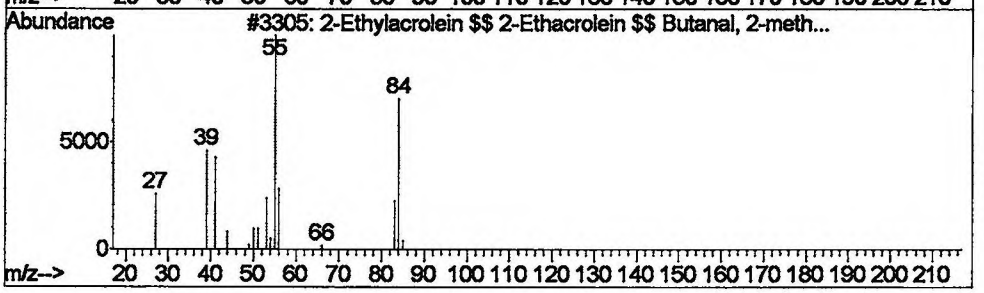


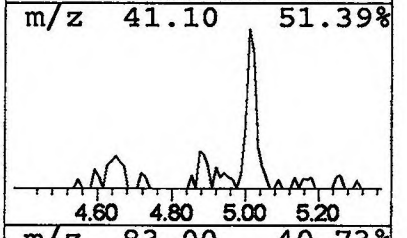


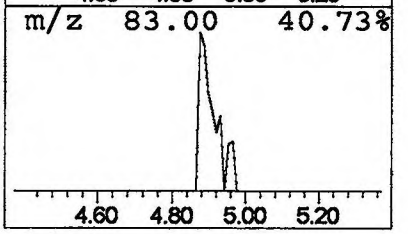












Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

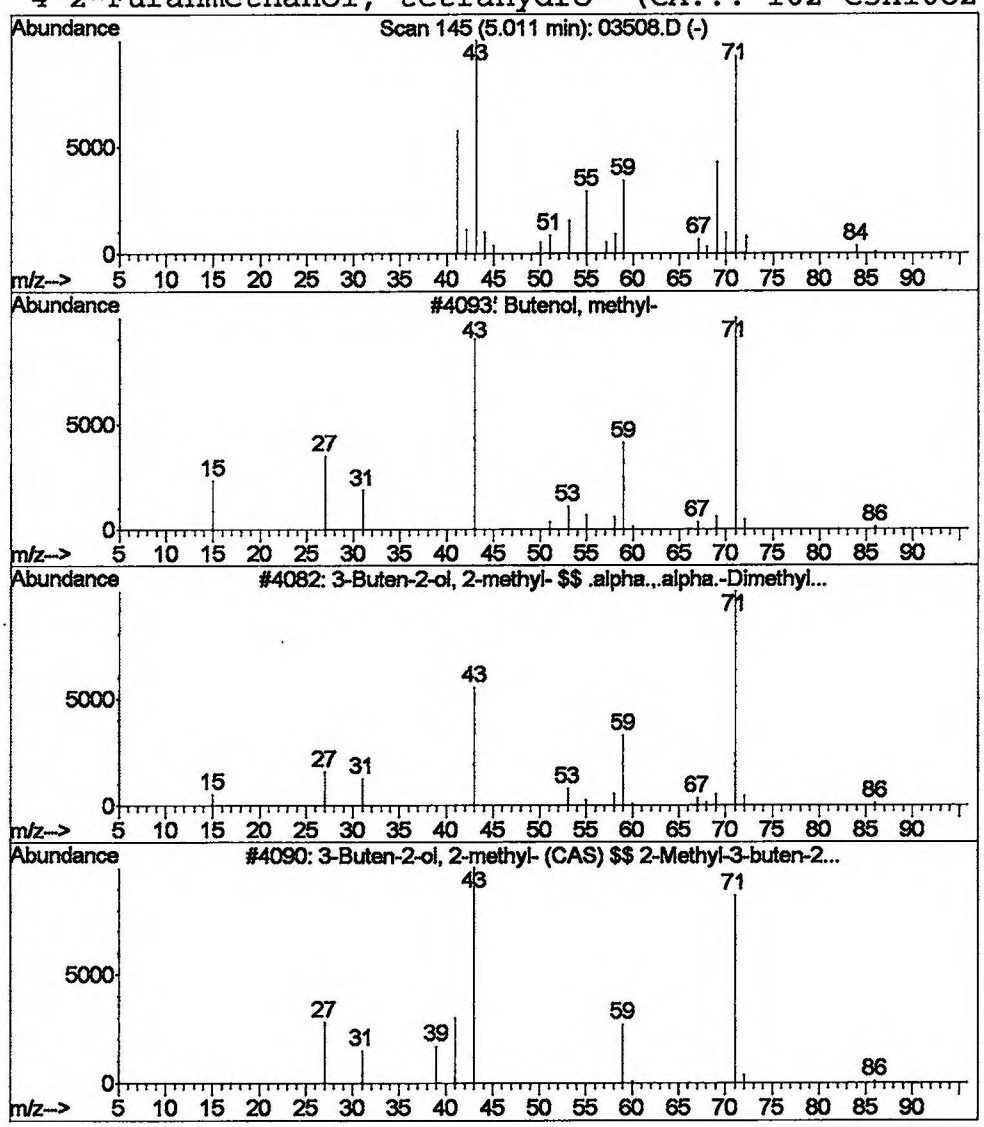
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

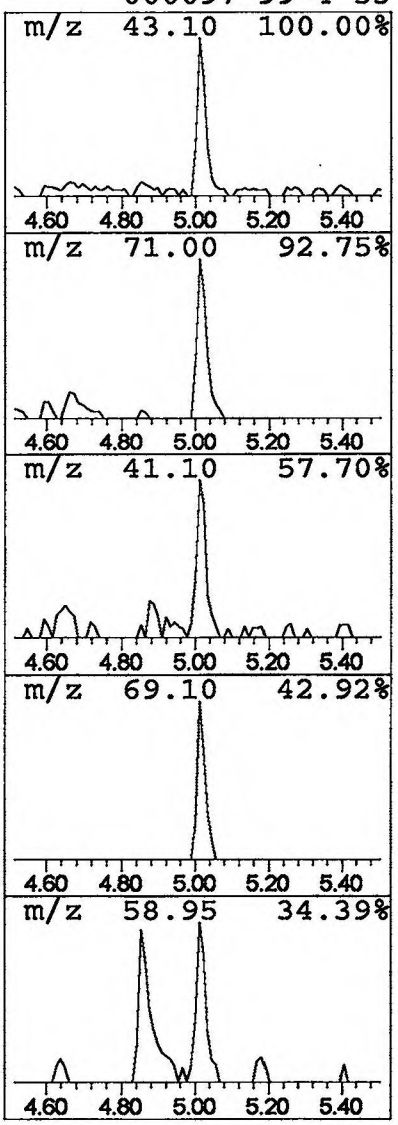
****

Peak Number 6 Butenol, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	0.27 ug/L	127911	1,4-Dichlorobenzene-d4	9.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butenol, methyl-	86	C5H10O	060766-00-9	64
2	3-Buten-2-ol, 2-methyl- $$ .alph...	86	C5H10O	000115-18-4	59
3	3-Buten-2-ol, 2-methyl- (CAS) $$...	86	C5H10O	000115-18-4	53
4	2-Furanmethanol, tetrahydro- (CA...	102	C5H10O2	000097-99-4	53





Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

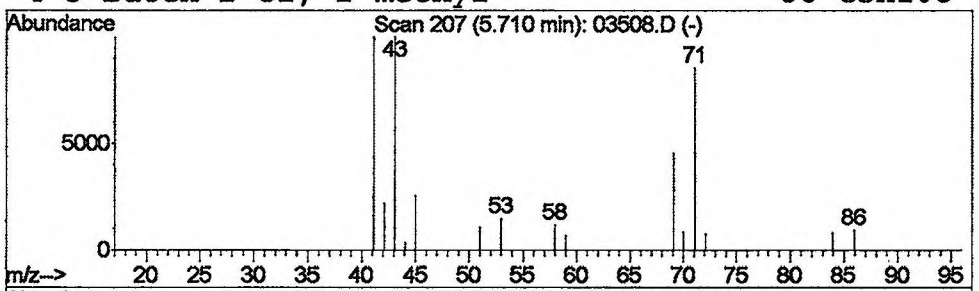
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

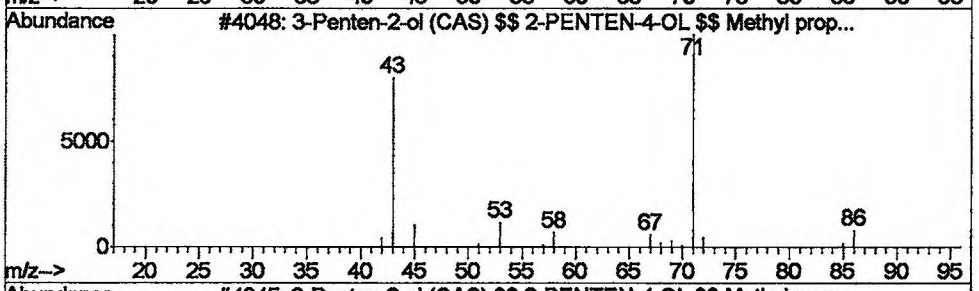
****

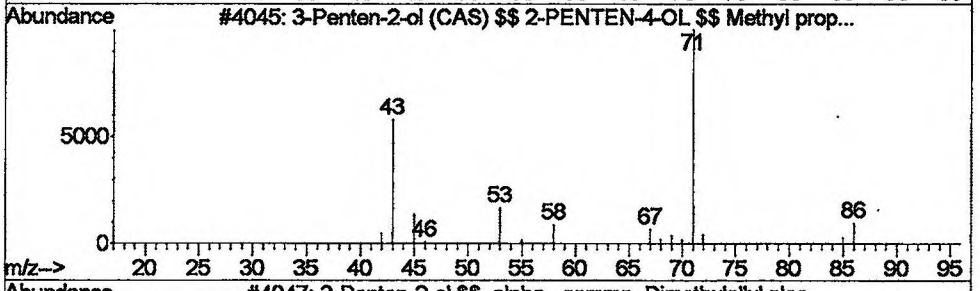
Peak Number 7 3-Penten-2-ol (CAS) $$ 2-PE... Concentration Rank 6

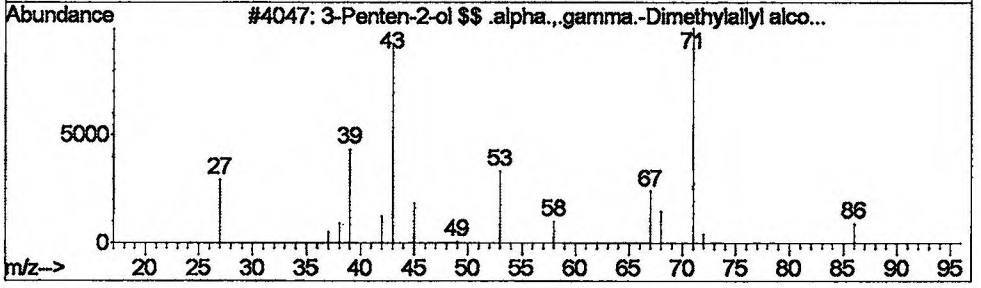
R.T.	EstConc	Area	Relative to ISTD	R.T.
5.71	0.13 ug/L	61888	1,4-Dichlorobenzene-d4	9.72

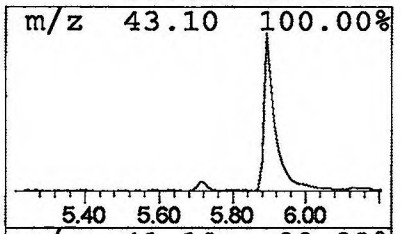
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Penten-2-ol (CAS) $$ 2-PENTEN-...	86	C5H10O	001569-50-2	59
2	3-Penten-2-ol (CAS) $$ 2-PENTEN-...	86	C5H10O	001569-50-2	59
3	3-Penten-2-ol $$ .alpha.,.gamma....	86	C5H10O	001569-50-2	53
4	3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	50

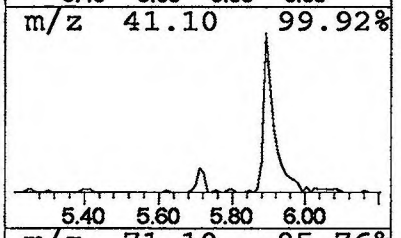


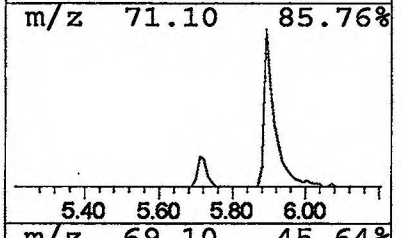


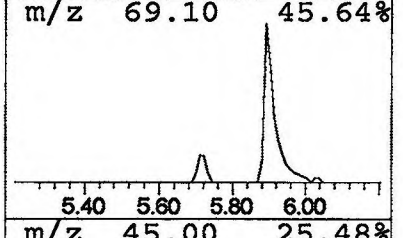


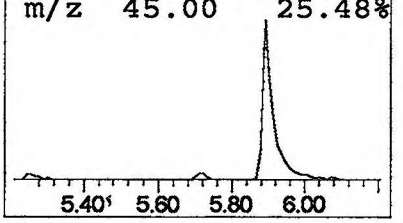












Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

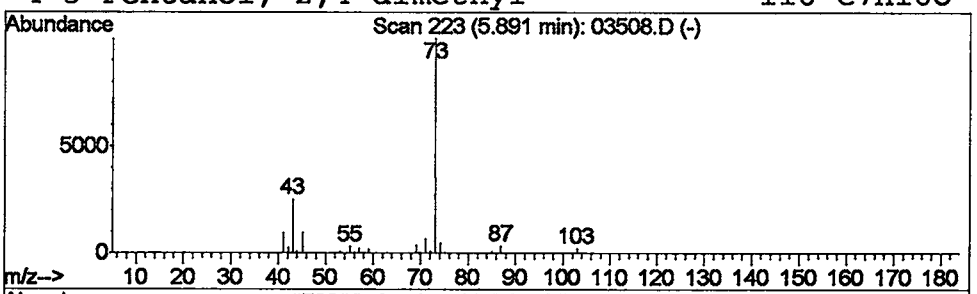
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

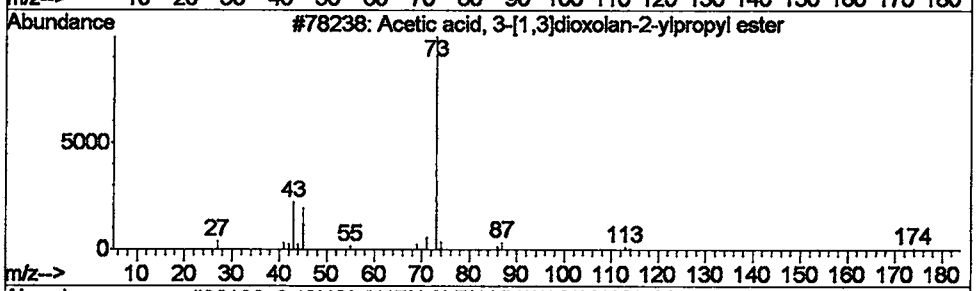
****

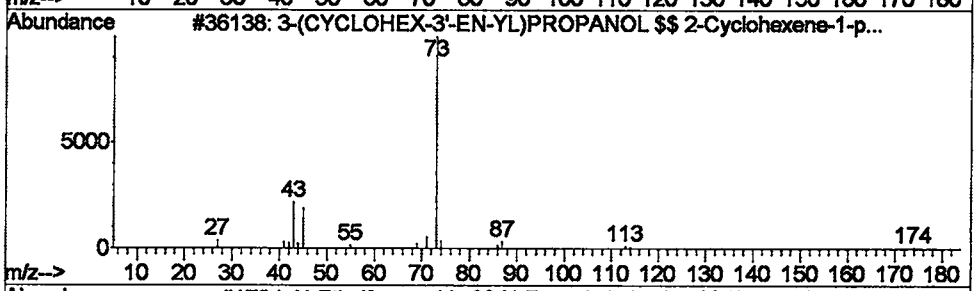
Peak Number 8 Acetic acid, 3-[1,3]dioxola... Concentration Rank 1

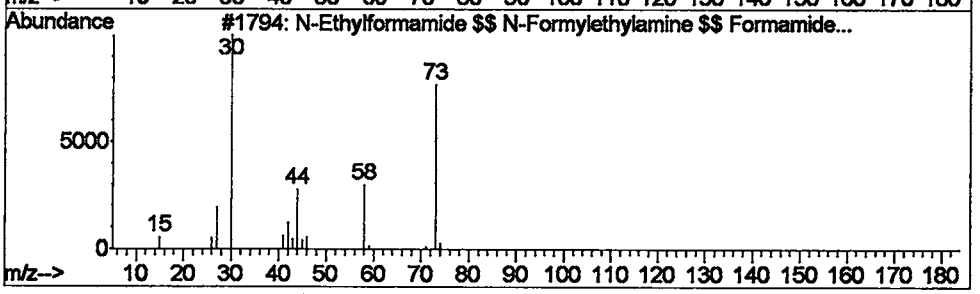
R.T.	EstConc	Area	Relative to ISTD	R.T.
5.89	3.89 ug/L	1874100	1,4-Dichlorobenzene-d4	9.72

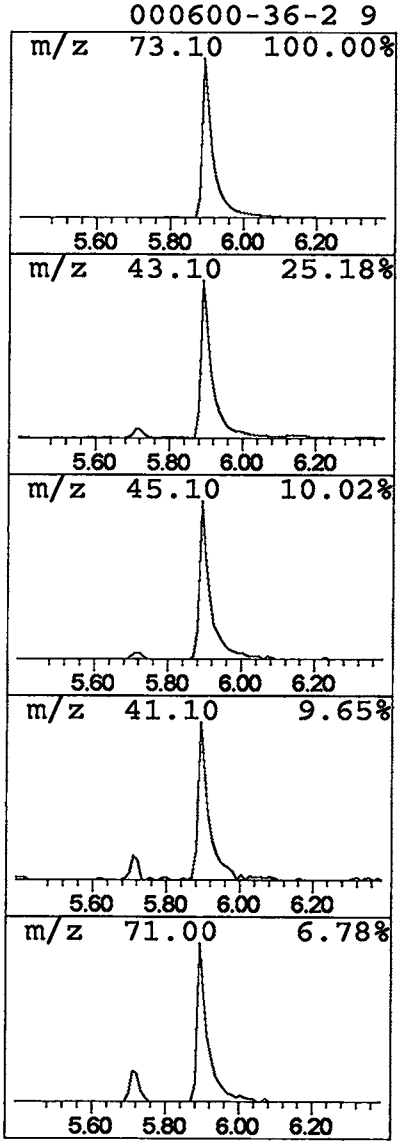
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, 3-[1,3]dioxolan-2-y...	174	C8H14O4	000000-00-0	33
2	3-(CYCLOHEX-3'-EN-YL) PROPANOL $$...	174	C8H14O4	015745-87-6	33
3	N-Ethylformamide $$ N-Formylethy...	73	C3H7NO	000627-45-2	9
4	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	9











Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

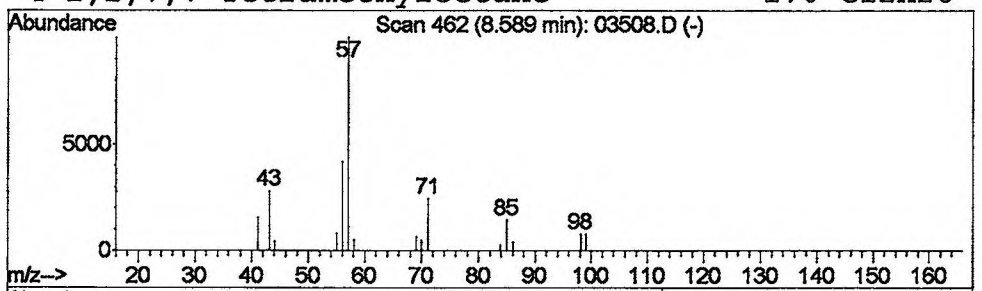
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

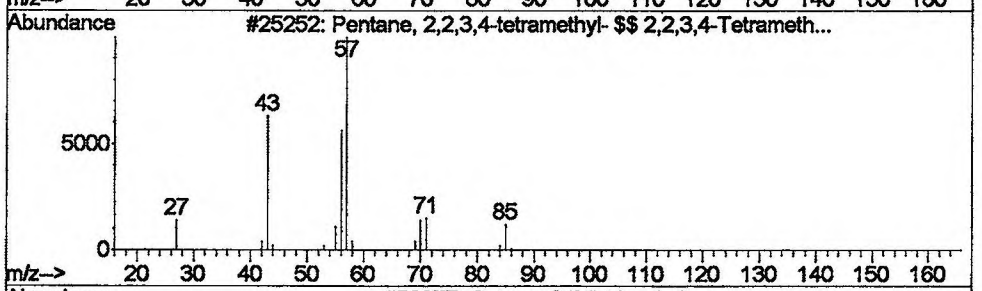
****

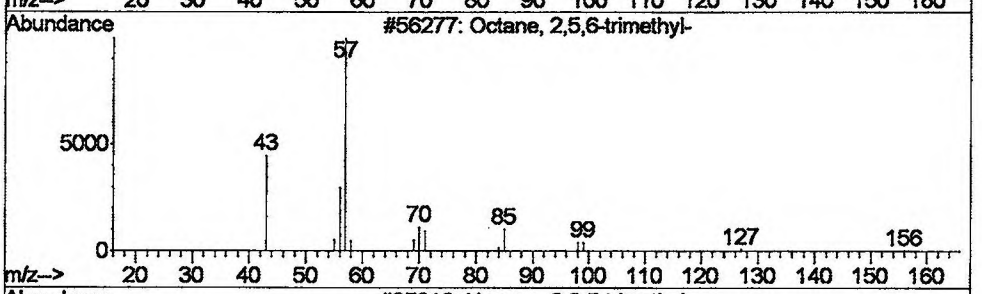
Peak Number 9 Pentane, 2,2,3,4-tetramethy... Concentration Rank 7

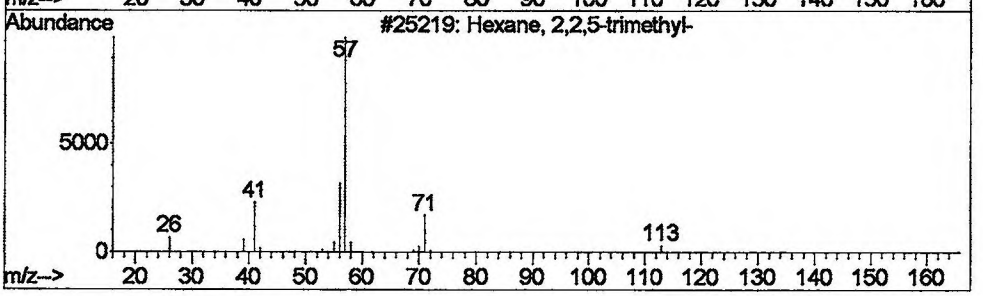
R.T.	EstConc	Area	Relative to ISTD	R.T.
8.59	0.12 ug/L	58135	1,4-Dichlorobenzene-d4	9.72

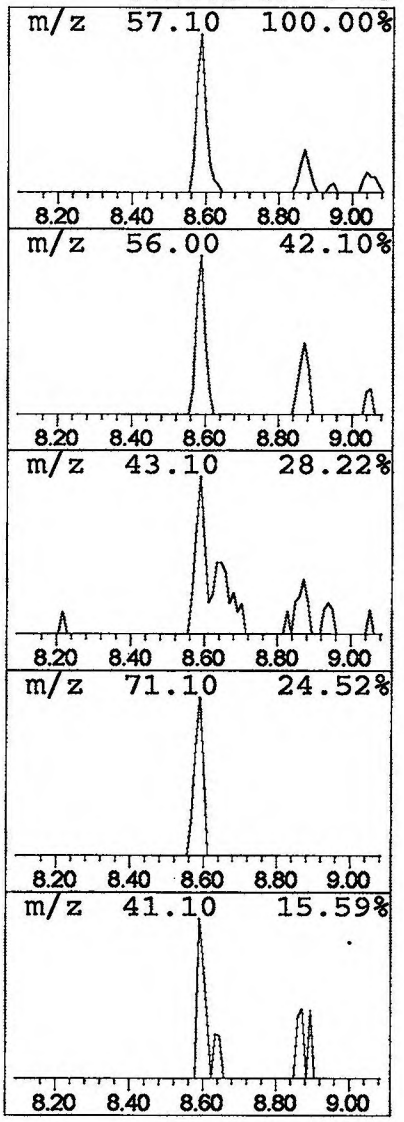
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,2,3,4-tetramethyl- $$...	128	C9H20	001186-53-4	59
2	Octane, 2,5,6-trimethyl-	156	C11H24	062016-14-2	50
3	Hexane, 2,2,5-trimethyl-	128	C9H20	003522-94-9	45
4	2,2,7,7-Tetramethyloctane	170	C12H26	001071-31-4	45











Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

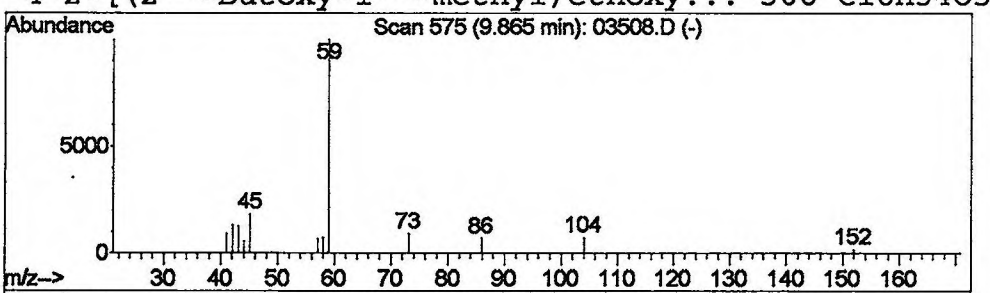
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

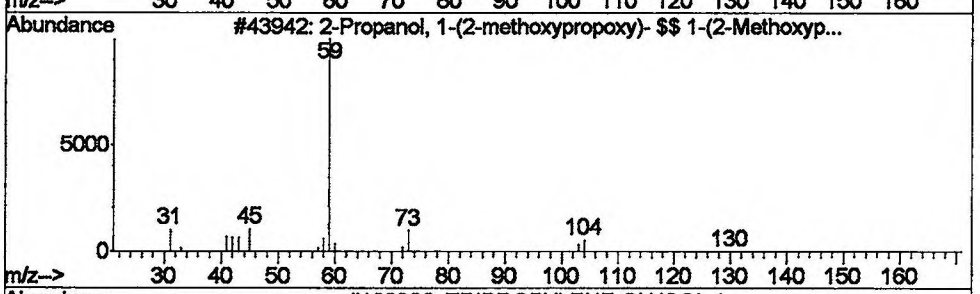
****

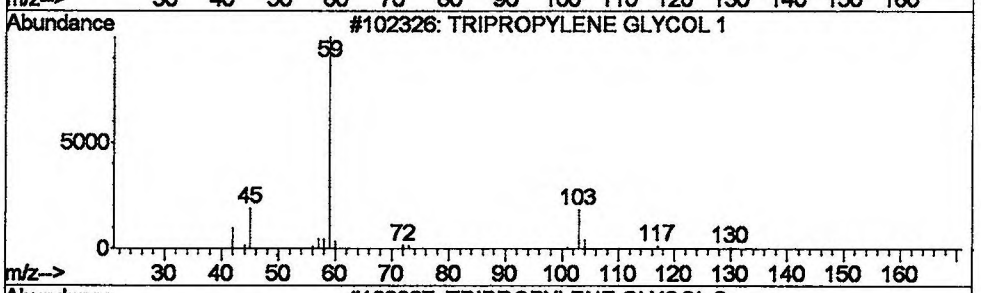
Peak Number 10 2-Propanol, 1-(2-methoxypro... Concentration Rank 12

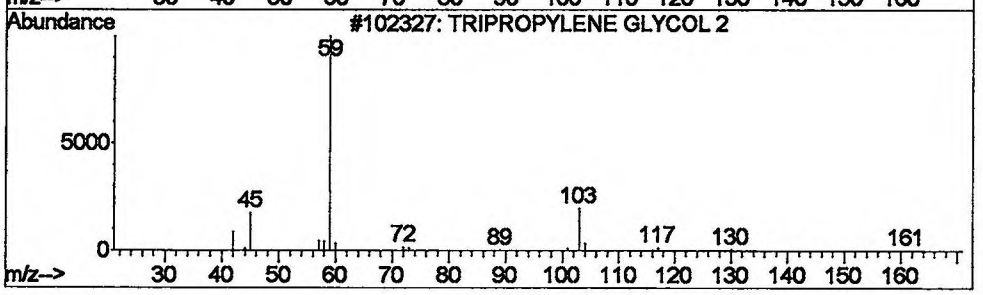
R.T.	EstConc	Area	Relative to ISTD	R.T.
9.86	0.08 ug/L	39077	1,4-Dichlorobenzene-d4	9.72

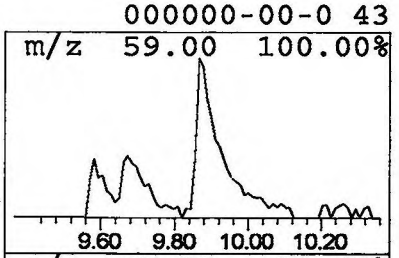
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-(2-methoxypropoxy)...	148	C7H16O3	013429-07-7	74
2	TRIPROPYLENE GLYCOL 1	192	C9H20O4	000000-00-0	56
3	TRIPROPYLENE GLYCOL 2	192	C9H20O4	000000-00-0	56
4	2-[(2''-Butoxy-1''-methyl)ethoxy...	306	C16H34O5	000000-00-0	43

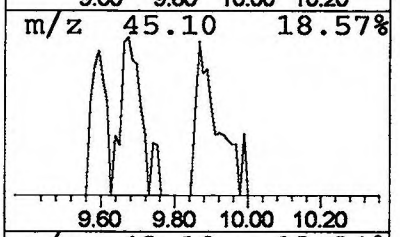


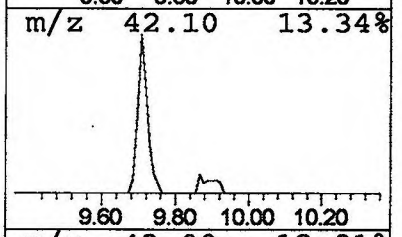


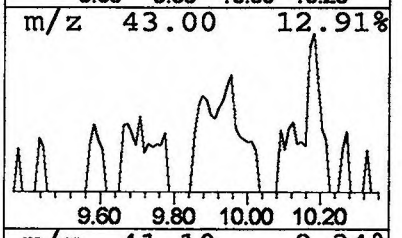


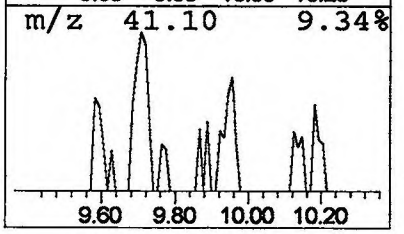












Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

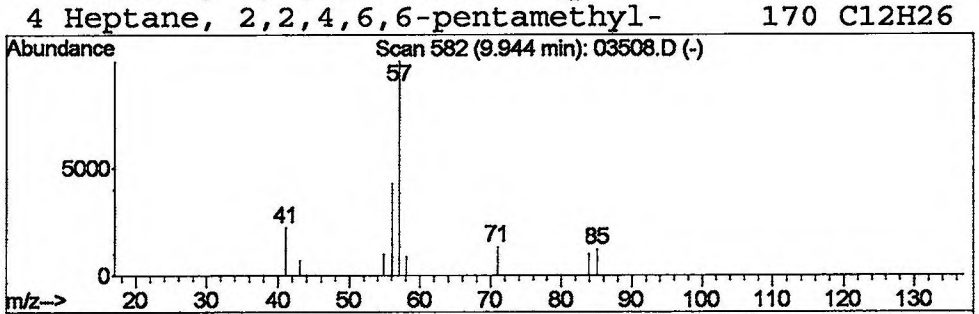
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

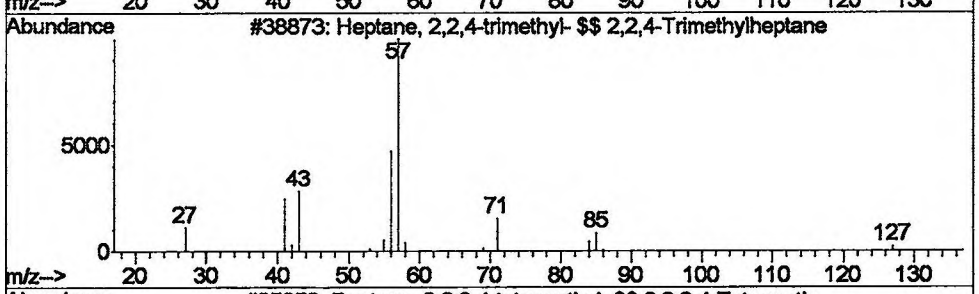
****

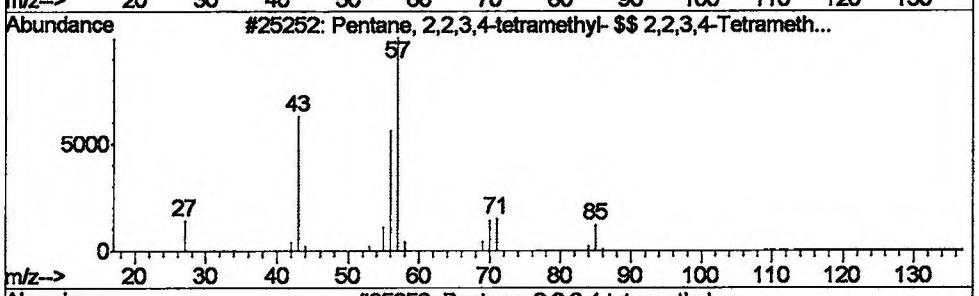
Peak Number 11 Heptane, 2,2,4-trimethyl- $... Concentration Rank 10

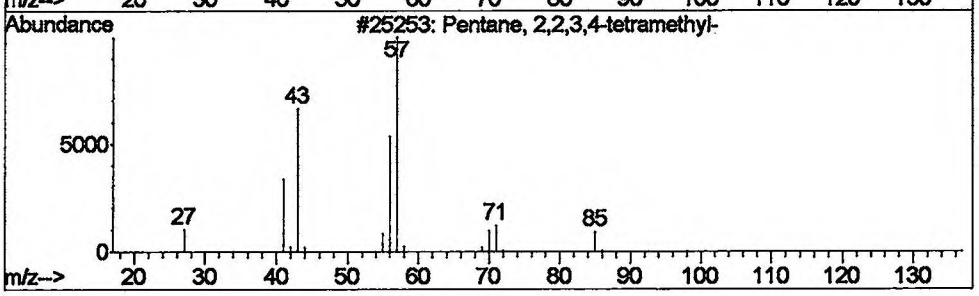
R.T.	EstConc	Area	Relative to ISTD	R.T.
9.94	0.10 ug/L	50278	1,4-Dichlorobenzene-d4	9.72

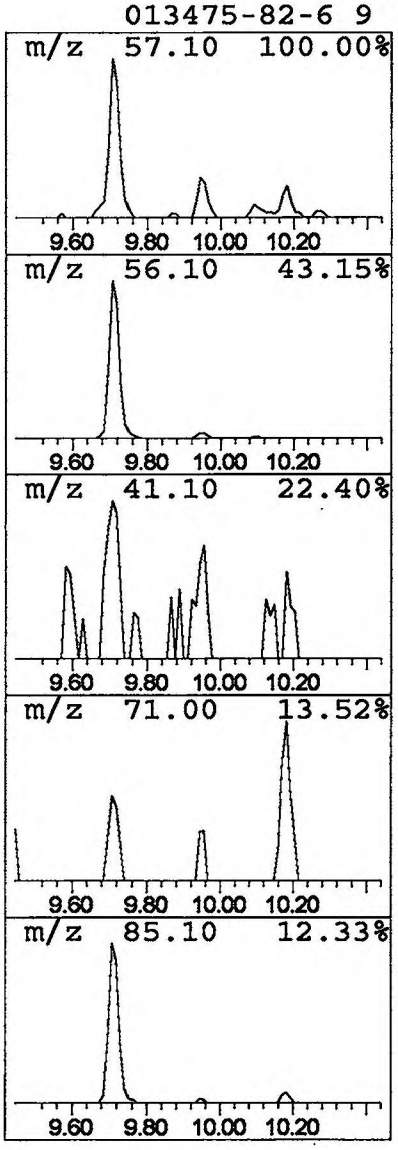
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,2,4-trimethyl- $$ 2,2...	142	C10H22	014720-74-2	64
2	Pentane, 2,2,3,4-tetramethyl- $$...	128	C9H20	001186-53-4	43
3	Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	40
4	Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	9











Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

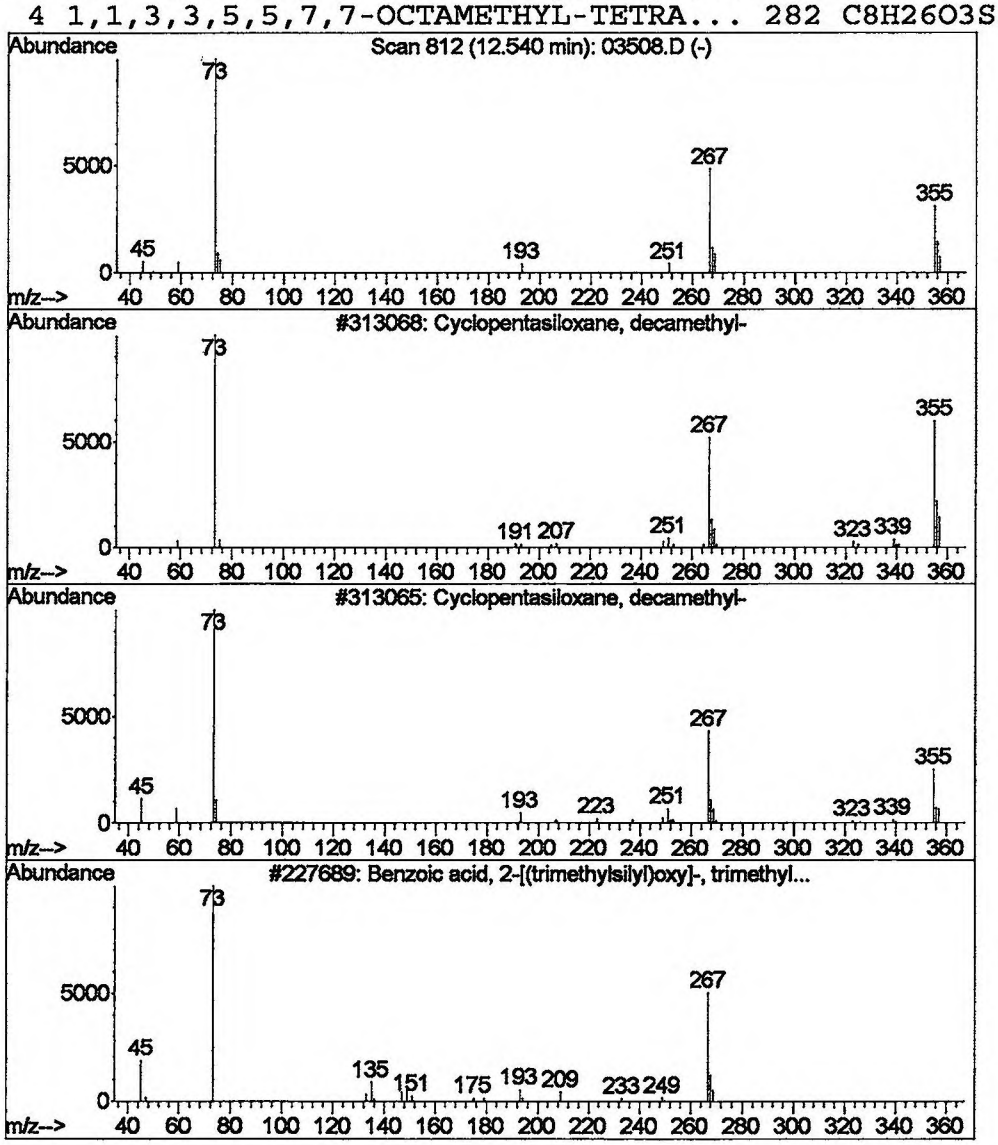
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

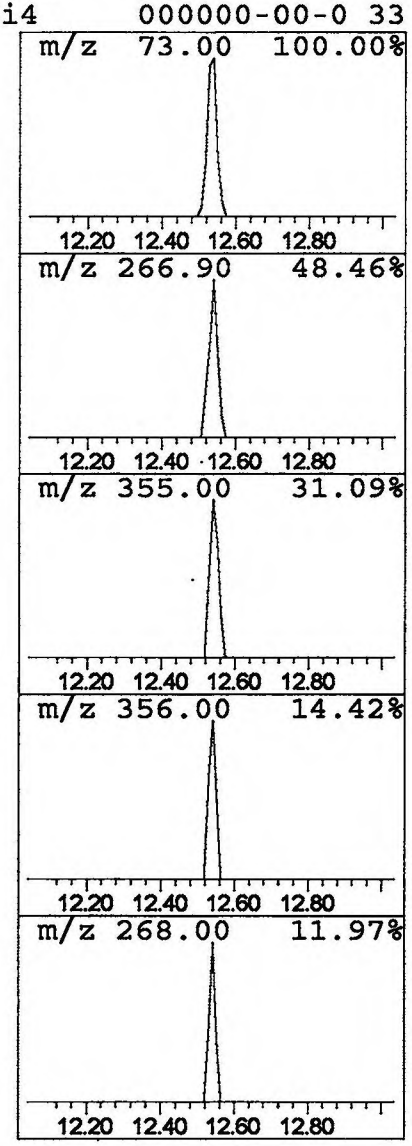
****

Peak Number 12 Cyclopentasiloxane, decamet... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	0.08 ug/L	59189	Naphthalene-d8	13.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	53
2	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	45
3	Benzoic acid, 2-[(trimethylsilyl...	282	C13H22O3Si2	003789-85-3	33
4	1,1,3,3,5,5,7,7-OCTAMETHYL-TETRA...	282	C8H26O3Si4	000000-00-0	33





Library Search Compound Report

Data File : C:\MSDCHEM\1\5973N\02MAR08\03508.D Vial: 4
Acq On : 8 Mar 2002 1:07 pm Operator:
Sample : 508 scan Inst : Instrumen
Misc : March 8, 2002 Multiplr: 1.00
MS Integration Params: LSCINT.P

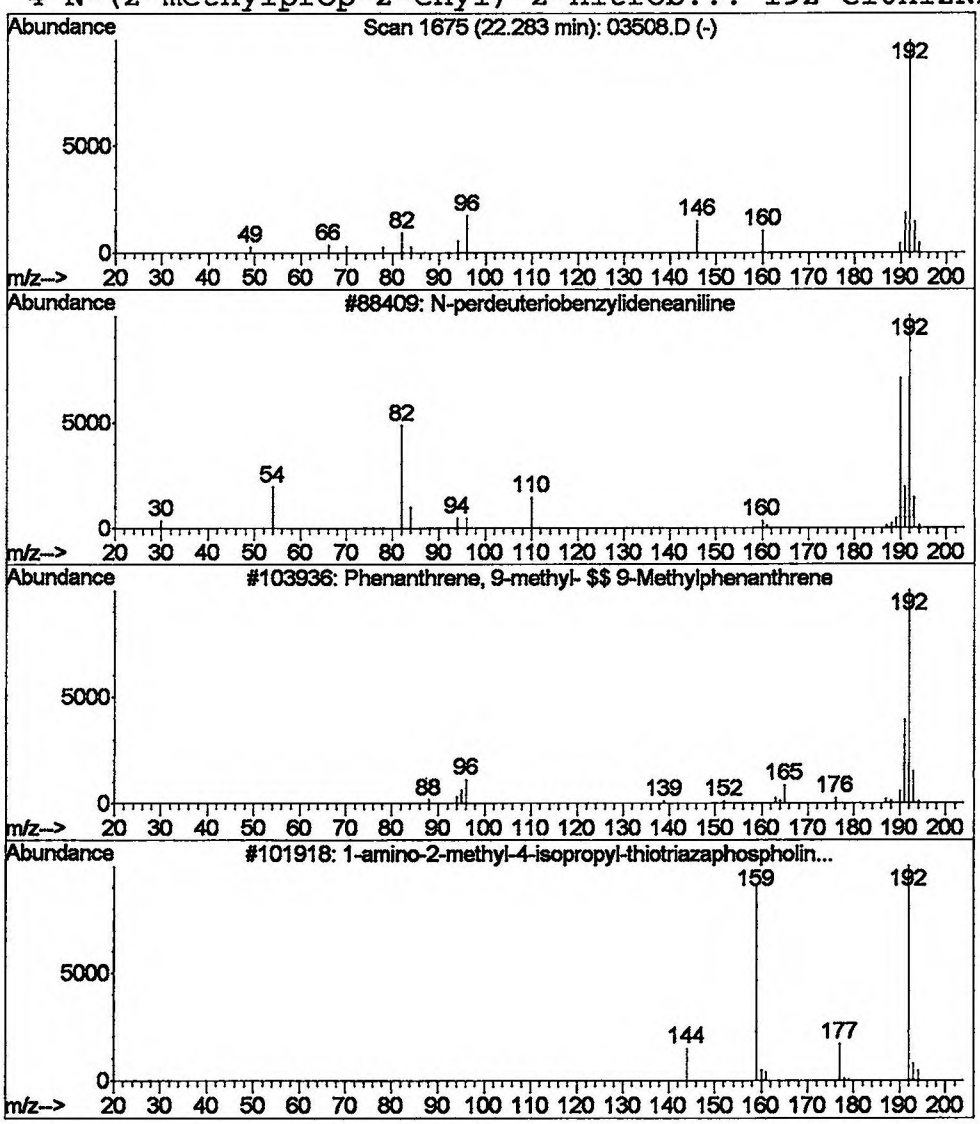
Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)
Title : 508 scan method
Library : C:\DATABASE\WILEY7N.L

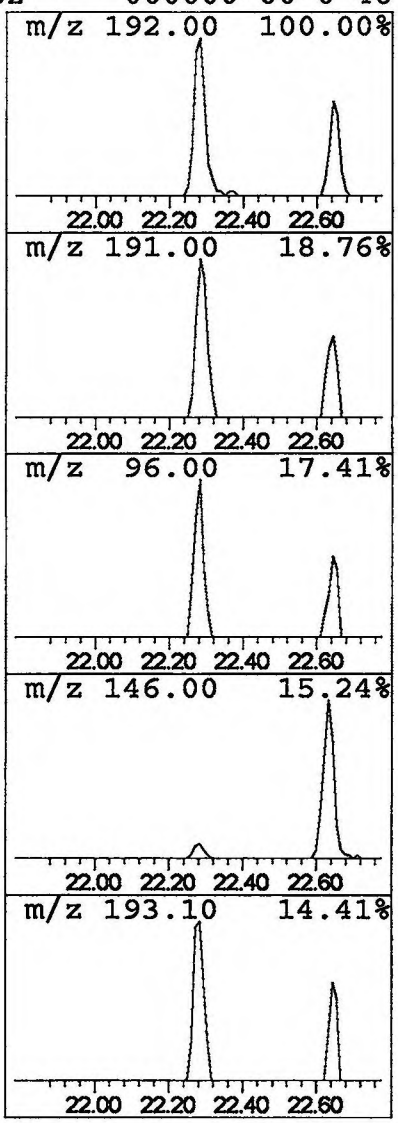
****

Peak Number 13 N-perdeuteriobenzylideneani... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	0.11 ug/L	84814	Phenanthrene-d10	22.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N-perdeuteriobenzylideneaniline	192	C13D11N	000538-51-2	64
2	Phenanthrene, 9-methyl- $$ 9-Met...	192	C15H12	000883-20-5	53
3	1-amino-2-methyl-4-isopropyl-thi...	192	C5H13N4PS	107264-54-0	50
4	N-(2-methylprop-2-enyl)-2-nitrob...	192	C10H12N2O2	000000-00-0	46





Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 8 Mar 2002 1:07 pm
Data File: C:\MSDCHEM\1\5973N\02MAR08\03508.D
Name. 508 scan
Misc: March 8, 2002
Method: C:\MSDCHEM\1\METHODS\METHODS\HP022102.M (RTE Integrator)
Title: 508 scan method
Library Searched: C:\DATABASE\WILEY7N.L

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--- #	RT	Resp	Conc
Butanoic acid, me...	3.61	0.2	ug/L	72967	1	9.72	9633320	20.0
ISOBUTYRALDEHYDE,...	3.93	0.1	ug/L	37878	1	9.72	9633320	20.0
2,2-Dimethoxybutane	4.24	0.1	ug/L	70114	1	9.72	9633320	20.0
Butanoic acid, 2-...	4.64	0.1	ug/L	53206	1	9.72	9633320	20.0
3,4-Dihydropyran ...	4.88	0.1	ug/L	62723	1	9.72	9633320	20.0
Butenol, methyl-	5.01	0.3	ug/L	127911	1	9.72	9633320	20.0
3-Penten-2-ol (CA...	5.71	0.1	ug/L	61888	1	9.72	9633320	20.0
Acetic acid, 3-[1...	5.89	3.9	ug/L	1874100	1	9.72	9633320	20.0
Pentane, 2,2,3,4-...	8.59	0.1	ug/L	58135	1	9.72	9633320	20.0
2-Propanol, 1-(2-...	9.86	0.1	ug/L	39077	1	9.72	9633320	20.0
Heptane, 2,2,4-tr...	9.94	0.1	ug/L	50278	1	9.72	9633320	20.0
Cyclopentasiloxan...	12.54	0.1	ug/L	59189	2	13.20	13944400	20.0
N-perdeuteriobenz...	22.28	0.1	ug/L	84814	4	22.63	15671400	20.0