User: Fakhouri, Morgan
Westchester Dept. of Labs & Research
Date: 05/09/2002 Time: 12:14:31

Sample: AE10603
Analysis: $B1GCMS Blank for GCMS Scan
Units: ug
Started: 5/9/02 12:14 Ended: 5/9/02 12:14 Analyst: MF
Result source: manual entry
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	n-Nitrosodimethylamine		< MDL		2.0
2	bis(2-Chloroethyl)ether		< MDL		2.0
3	1,3-Dichlorobenzene		< MDL		2.0
4	1,4-Dichlorobenzene		< MDL		2.0
5	1,2-Dichlorobenzene		< MDL		2.0
6	2,2'-oxybis(1-Chloropropane)		< MDL		2.0
7	n-Nitrosodi-n-propylamine		< MDL		2.0
8	Hexachloroethane		< MDL		2.0
9	Nitrobenzene		< MDL		2.0
10	Isophorone		< MDL		2.0
11	bis(2-Chloroethoxy)methane		< MDL		2.0
12	1,2,4-Trichlorobenzene		< MDL		2.0
13	Naphthalene		< MDL		2.0
14	Hexachlorobutadiene		< MDL		2.0
15	2-Chloronaphthalene		< MDL		2.0
16	Dimethylphthalate		< MDL		2.0
17	2,6-Dinitrotoluene		< MDL		2.0
18	Acenaphthylene		< MDL		2.0
19	Acenaphthene		< MDL		2.0
20	2,4-Dinitrotoluene		< MDL		2.0
21	Diethylphthalate		< MDL		2.0
22	4-Chlorophenylphenylether		< MDL		2.0
23	Fluorene		< MDL		2.0
24	Azobenzene		< MDL		2.0
25	n-Nitrosodiphenylamine		< MDL		2.0
26	4-Bromophenylphenylether		< MDL		2.0
27	Hexachlorobenzene		< MDL		2.0
28	Phenanthrene		< MDL		2.0
29	Anthracene		< MDL		2.0
30	Di-n-butylphthalate		< MDL		2.0
31	Fluoranthene		< MDL		2.0
32	Pyrene		< MDL		2.0
33	Butylbenzylphthalate		< MDL		2.0
34	Benzo(a)anthracene		< MDL		2.0
35	bis(2-Ethylhexyl)phthalate		< MDL		2.0
36	Chrysene		< MDL		2.0
37	Di-n-octylphthalate		< MDL		2.0
38	Benzo(b)fluoranthene		< MDL		2.0
39	Benzo(k)fluoranthene		< MDL		2.0
40	Benzo(a)pyrene		< MDL		2.0
41	Indeno(1,2,3-cd)pyrene		< MDL		2.0
42	Dibenz(a,h)anthracene		< MDL		2.0
43	Benzo(g,h,i)perylene		< MDL		2.0

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\050702\0503LB.D Vial: 5
Acq On : 7 May 2002 4:24 pm Operator: Mclean
Sample : 5/3 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 09 09:20:19 2002 Quant Results File: 050102.RES

Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Wed May 08 09:53:56 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.94	152	6023194	40.00	ug/L	0.01
10) Napthalene-d8	13.44	136	23909835	40.00	ug/L	0.00
17) Acenapthalene-d10	18.58	164	13075378	40.00	ug/L	0.00
29) Phenanthranene-d10	22.96	188	19252439	40.00	ug/L	0.00
37) Chrysene-d12	30.82	240	15038964	40.00	ug/L	0.00
43) Perylene-d12	34.72	264	10184272	40.00	ug/L	0.01
System Monitoring Compounds						
27) TCMX	20.54	209	655857	8.68	ug/L	0.00
Spiked Amount 10.000			Recovery = 86.80%			
50) 2,4-ddd	0.00	235	0	0.00	ug/L	
Spiked Amount 5.000			Recovery = 0.00%			
Target Compounds						Qvalue
2) n-nitros-dimethyl amine	0.00	74	0	N.D.		
3) bis(2-Chloroethyl)ether	0.00	93	0	N.D.		
4) 1,3-Dichlorobenzene	0.00	146	0	N.D.		
5) 1,4-Dichlorobenzene	0.00	146	0	N.D.		
6) 1,2-Dichlorobenzene	0.00	146	0	N.D.		
7) 2,2'oxybis(1-Chloropropane	0.00	45	0	N.D.		
8) N-nitroso-di-propylamine	0.00	70	0	N.D.		
9) Hexachloroethane	0.00	117	0	N.D.		
11) Nitrobenzene	0.00	77	0	N.D.		
12) Isophorone	0.00	82	0	N.D.		
13) bis(2-Chloroethoxy) methan	12.94	93	533	N.D.		
14) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.		
15) Napthalene	0.00	128	0	N.D.		
16) Hexachlorobutadiene	0.00	225	0	N.D.		
18) 2-Chloronaphthalene	0.00	162	0	N.D.		
19) Dimethylphthalate	0.00	163	0	N.D.		
20) 2,6-Dinitrotoluene	0.00	165	0	N.D.		
21) Acenapthylene	0.00	152	0	N.D.		
22) Acenapthalene	0.00	153	0	N.D.		
23) 2,4-Dinitrotoluene	0.00	165	0	N.D.		
24) Diethylphthalate	0.00	149	0	N.D.		
25) 4-Chlorophenyl-phenylether	0.00	204	0	N.D.		
26) Fluorene	0.00	166	0	N.D.		
28) Azobenzene	0.00	77	0	N.D.		
30) Nnitrosodiphenylamine	0.00	169	0	N.D.		
31) 4-Bromophenyl-phenylether	0.00	248	0	N.D.		
32) Hexachlorobenzene	0.00	284	0	N.D.		
33) Phenanthrene	0.00	178	0	N.D.		

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

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Sample : 5/3 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 09 09:20:19 2002 Quant Results File: 050102.RES

Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Wed May 08 09:53:56 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
34)	Anthracene	0.00	178	0	N.D.	
35)	Di-n-butylphthalate	25.09	149	20433	N.D.	
36)	Fluoroanthene	0.00	202	0	N.D.	
38)	Pyrene	0.00	202	0	N.D.	
39)	Butylbenzylphthalate	0.00	149	0	N.D.	
40)	Benzo(a)anthracene	30.82	228	44444	N.D.	
41)	Bis(2-ethylhexyl)phthalate	0.00	149	0	N.D.	
42)	Chrysene	0.00	228	0	N.D.	
44)	Di-n-octylphthalate	0.00	149	0	N.D.	
45)	Benzo(b)fluoranthene	0.00	252	0	N.D.	
46)	Benzo(k)fluoranthene	0.00	252	0	N.D.	
47)	Benzo(a)pyrene	0.00	252	0	N.D.	
48)	Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.	
49)	Dibenz(a,h)anthracene	0.00	278	0	N.D.	
51)	Benzo(g,h,i)perylene	0.00	276	0	N.D.	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\050702\0503LB.D Vial: 5
Acq On : 7 May 2002 4:24 pm Operator: Mclean
Sample : 5/3 kb Inst : Instrumen
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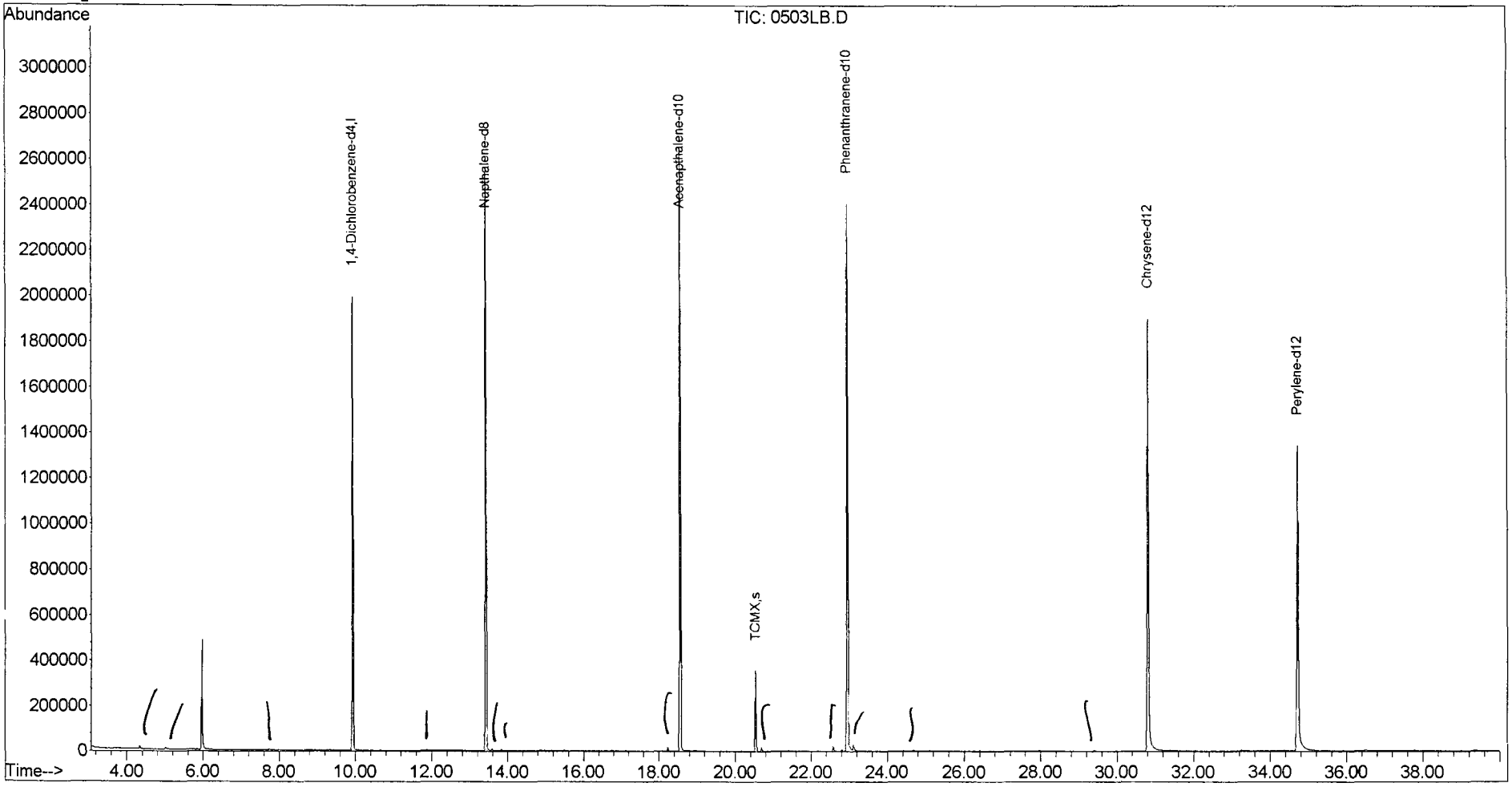
MS Integration Params: EVENTS.E
Quant Time: May 9 9:20 2002 Quant Results File: 050102.RES

Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Wed May 08 09:53:56 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\050702\0503LB.D
Acq On : 7 May 2002 4:24 pm
Sample : 5/3 kb
Misc :
MS Integration Params: LSCINT.e

Vial: 5
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan

Signal : TIC

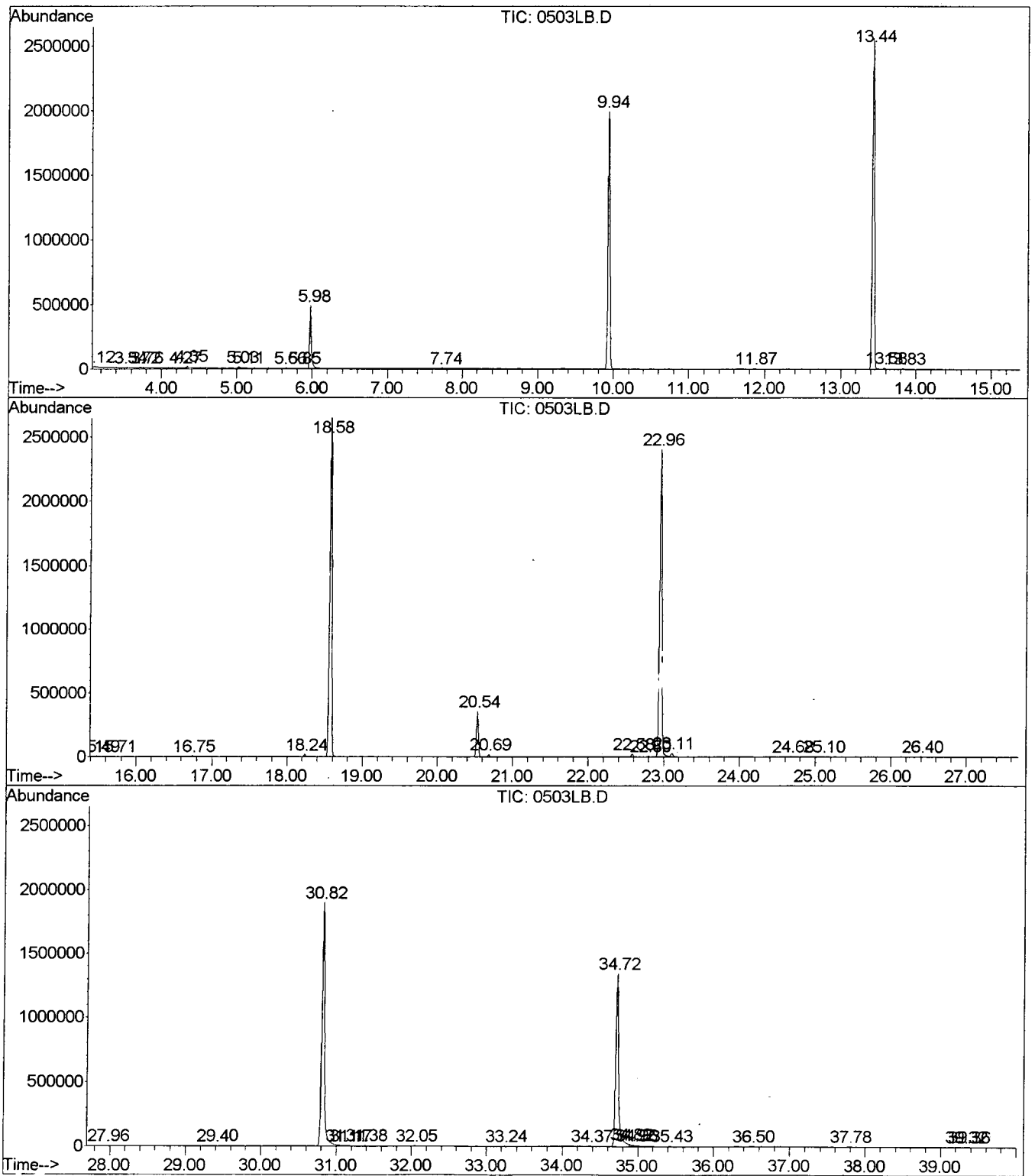
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.120	3	7	15	BV 3	3814	60344	0.11%	0.020%
2	3.543	76	79	87	PV 5	1896	25467	0.05%	0.009%
3	3.720	105	109	114	VV 5	2139	49545	0.09%	0.017%
4	3.761	114	116	119	VV 4	2072	27559	0.05%	0.009%
5	4.266	198	202	209	VV 6	2455	50896	0.09%	0.017%
6	4.348	209	216	223	VV	13882	228283	0.42%	0.077%
7	5.024	323	331	343	PV 2	8920	228597	0.42%	0.077%
8	5.106	343	345	354	VV 6	3095	60069	0.11%	0.020%
9	5.664	433	440	448	VV 8	2046	49864	0.09%	0.017%
10	5.846	465	471	487	PV 8	1668	38493	0.07%	0.013%
11	5.976	487	493	532	VV	471976	8128450	15.01%	2.732%
12	7.738	789	793	799	PV 2	4734	96542	0.18%	0.032%
13	9.942	1156	1168	1189	PV	1980076	36711361	67.80%	12.337%
14	11.869	1491	1496	1504	VV 2	4498	71334	0.13%	0.024%
15	13.438	1752	1763	1782	PV	2561100	49544645	91.50%	16.650%
16	13.585	1782	1788	1795	VV	8897	163833	0.30%	0.055%
17	13.831	1822	1830	1839	PV 3	5625	86533	0.16%	0.029%
18	15.488	2106	2112	2116	VV	3732	56724	0.10%	0.019%
19	15.706	2136	2149	2156	VV 5	2438	48186	0.09%	0.016%
20	16.746	2313	2326	2331	PV 3	3752	57103	0.11%	0.019%
21	18.238	2572	2580	2587	PV	15481	237526	0.44%	0.080%
22	18.579	2626	2638	2660	PV	2615665	54145077	100.00%	18.196%
23	20.541	2957	2972	2983	PV	350649	6475048	11.96%	2.176%
24	20.688	2983	2997	3010	VV	17796	286472	0.53%	0.096%
25	22.580	3310	3319	3331	PV 3	20581	390053	0.72%	0.131%
26	22.803	3351	3357	3367	VV 3	7560	135520	0.25%	0.046%
27	22.962	3371	3384	3404	PV 2	2404857	53033533	97.95%	17.823%
28	23.109	3404	3409	3424	VV	25794	611831	1.13%	0.206%
29	24.684	3666	3677	3686	PV 3	4397	94492	0.17%	0.032%
30	25.095	3733	3747	3758	PV 3	1714	37144	0.07%	0.012%
31	26.399	3963	3969	3974	PV 4	3331	59290	0.11%	0.020%
32	27.956	4226	4234	4246	VV 4	2570	66047	0.12%	0.022%
33	29.402	4473	4480	4488	PV 4	2872	56825	0.10%	0.019%
34	30.818	4700	4721	4770	PV 2	1881271	47672920	88.05%	16.021%
35	31.111	4770	4771	4778	VV 5	4280	99740	0.18%	0.034%
36	31.170	4778	4781	4788	VV 6	3049	86995	0.16%	0.029%
37	31.382	4809	4817	4831	VV 6	4762	165761	0.31%	0.056%

38	32.046	4913	4930	4936	VV	3	3218	79628	0.15%	0.027%
39	33.244	5126	5134	5144	PV	5	3309	68853	0.13%	0.023%
40	34.372	5319	5326	5335	VV	3	3296	86858	0.16%	0.029%
41	34.719	5365	5385	5413	VV	2	1331691	36911615	68.17%	12.405%
42	34.889	5413	5414	5419	VV	4	16746	284181	0.52%	0.096%
43	34.925	5419	5420	5425	VV	3	11558	215817	0.40%	0.073%
44	34.966	5425	5427	5433	VV	5	8307	189888	0.35%	0.064%
45	35.424	5498	5505	5514	VV	6	3234	74146	0.14%	0.025%
46	36.505	5673	5689	5696	PV	7	2687	75732	0.14%	0.025%
47	37.780	5892	5906	5914	BV	6	1762	63548	0.12%	0.021%
48	39.320	6153	6168	6170	PV	4	1416	30304	0.06%	0.010%
49	39.361	6170	6175	6183	VV	5	1738	44333	0.08%	0.015%

Sum of corrected areas: 297563005

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\050702\0503LB.D
Operator : Mclean
Acquired : 7 May 2002 4:24 pm using AcqMethod 508SCAN
Instrument : Instrumen
Sample Name: 5/3 kb
Misc Info :
Vial Number: 5
Quant File :050102.RES (Chemstation Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050702\0503LB.D Vial: 5
Acq On : 7 May 2002 4:24 pm Operator: Mclean
Sample : 5/3 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

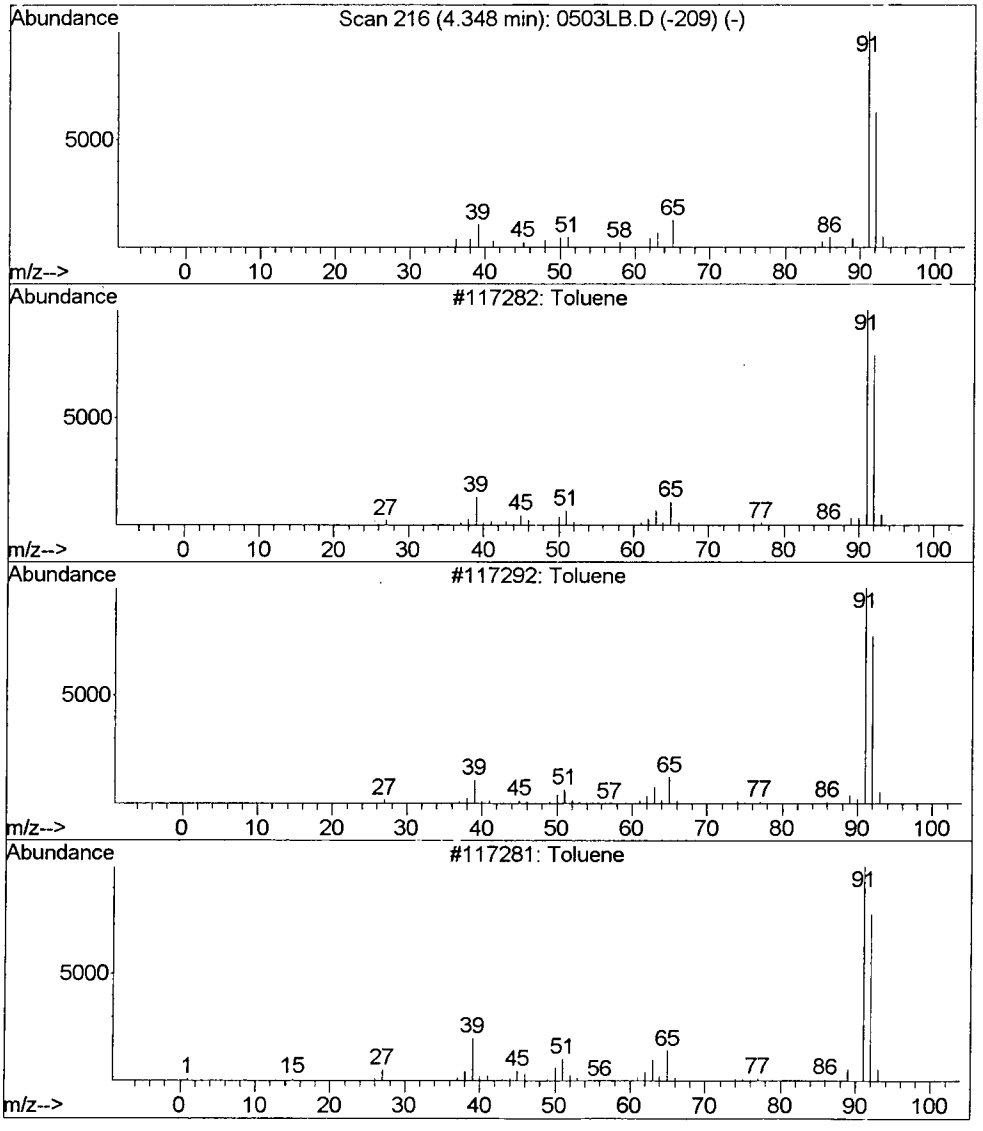
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

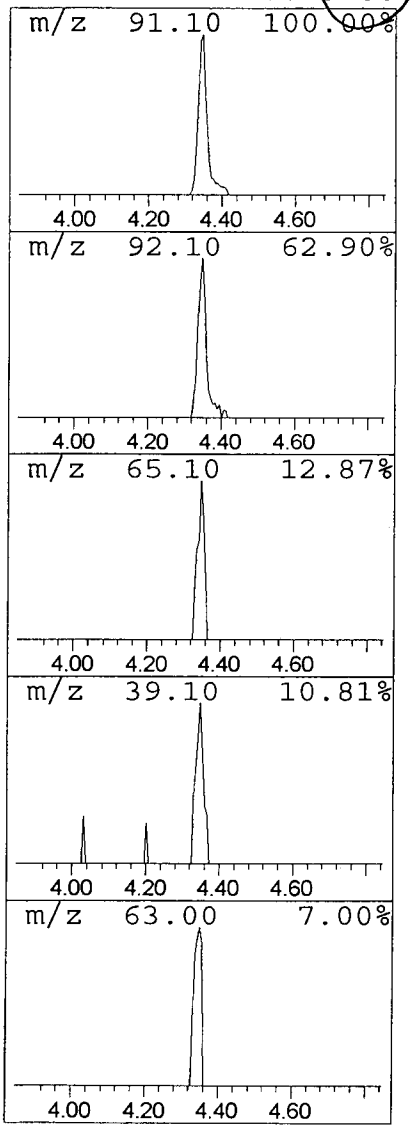
****

Peak Number 1 Toluene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.35	0.25 ug/L	228283	1,4-Dichlorobenzene-d4	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene	92	C7H8	000108-88-3	83
2	Toluene	92	C7H8	000108-88-3	83
3	Toluene	92	C7H8	000108-88-3	80
4	Toluene	92	C7H8	000108-88-3	80





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050702\0503LB.D Vial: 5
Acq On : 7 May 2002 4:24 pm Operator: Mclean
Sample : 5/3 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
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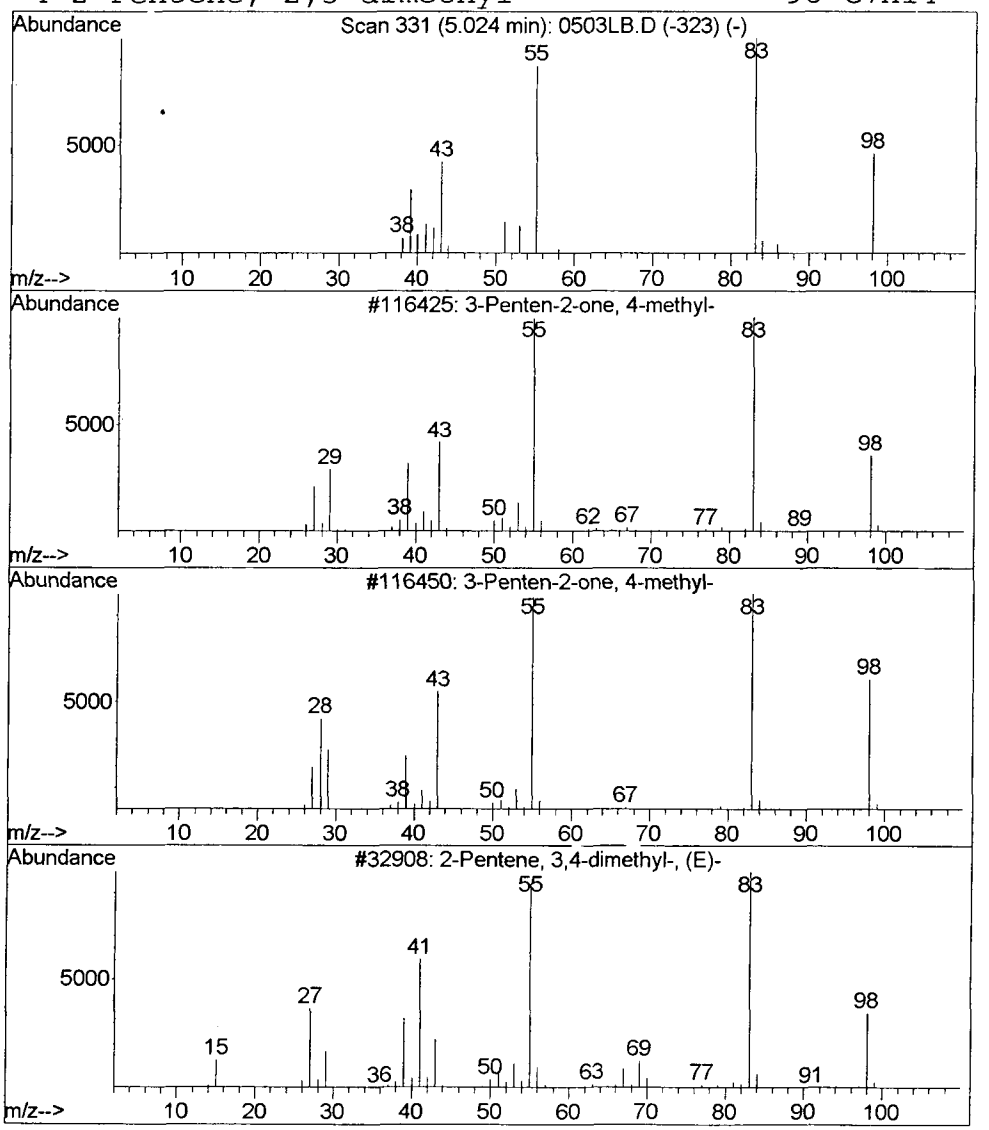
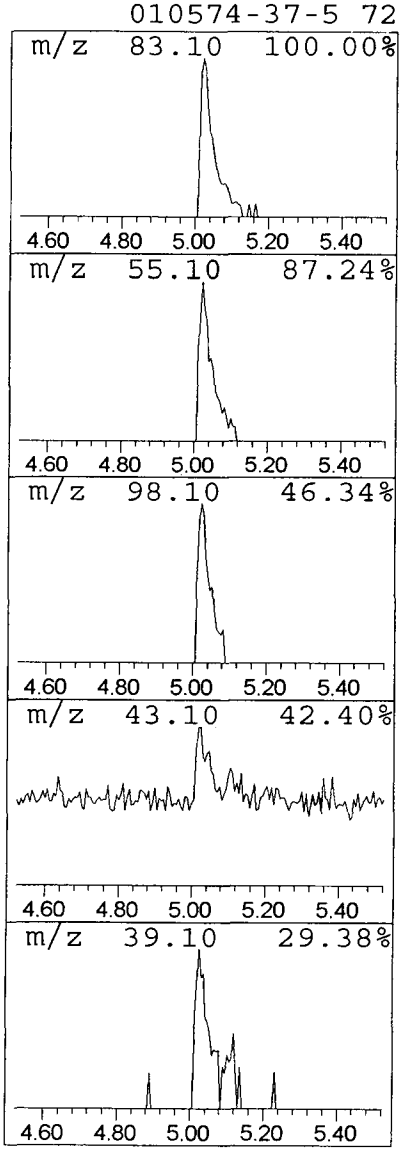
****

Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.03	0.25 ug/L	228597	1,4-Dichlorobenzene-d4	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	86
2	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	80
3	2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	72
4	2-Pentene, 2,3-dimethyl-	98	C7H14	010574-37-5	72





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050702\0503LB.D Vial: 5
Acq On : 7 May 2002 4:24 pm Operator: Mclean
Sample : 5/3 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

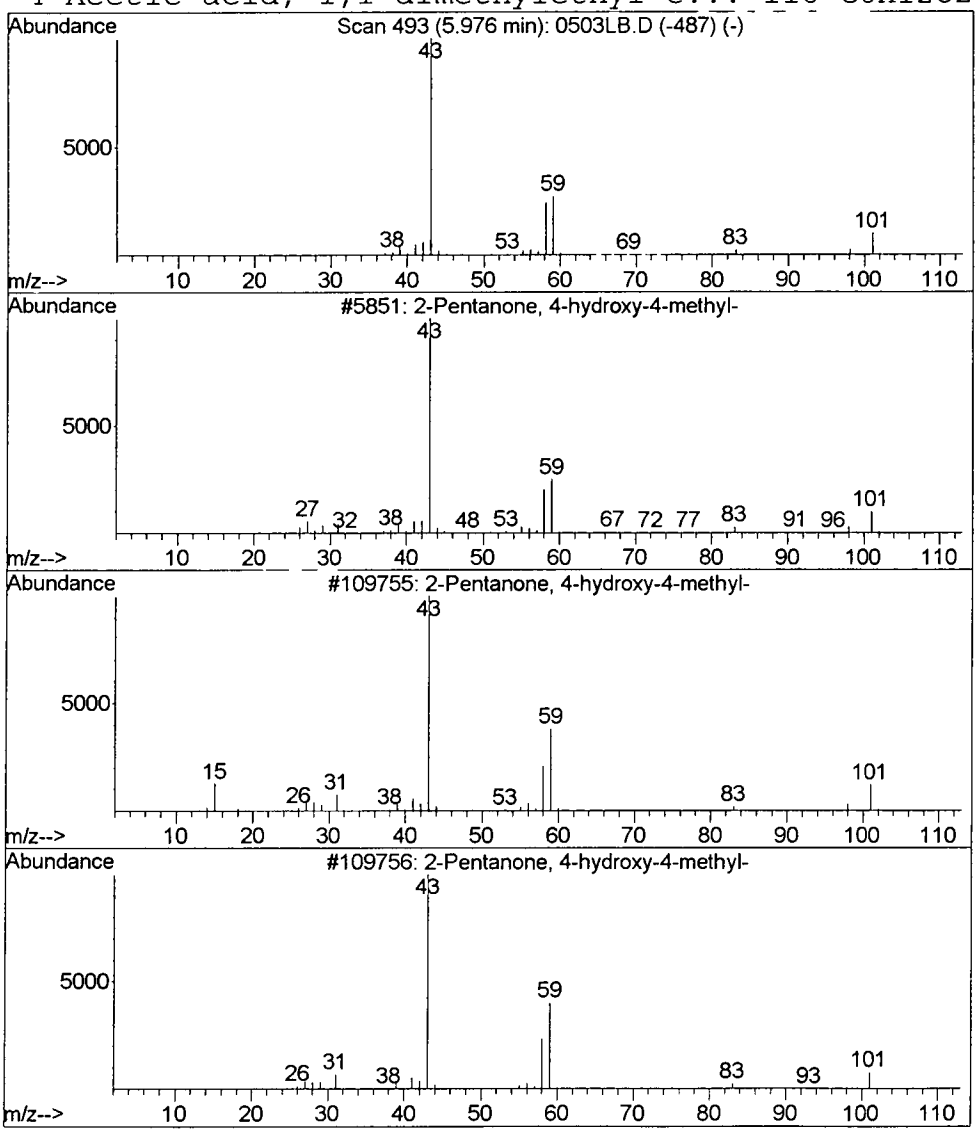
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

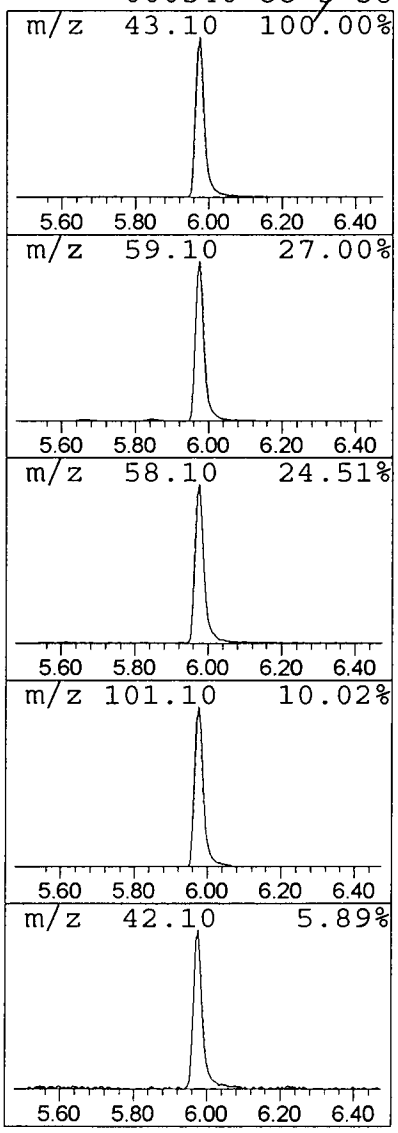
****

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.98	8.86 ug/L	8128450	1,4-Dichlorobenzene-d4	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38





Library Search Compound Report

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Acq On : 7 May 2002 4:24 pm Operator: Mclean
Sample : 5/3 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

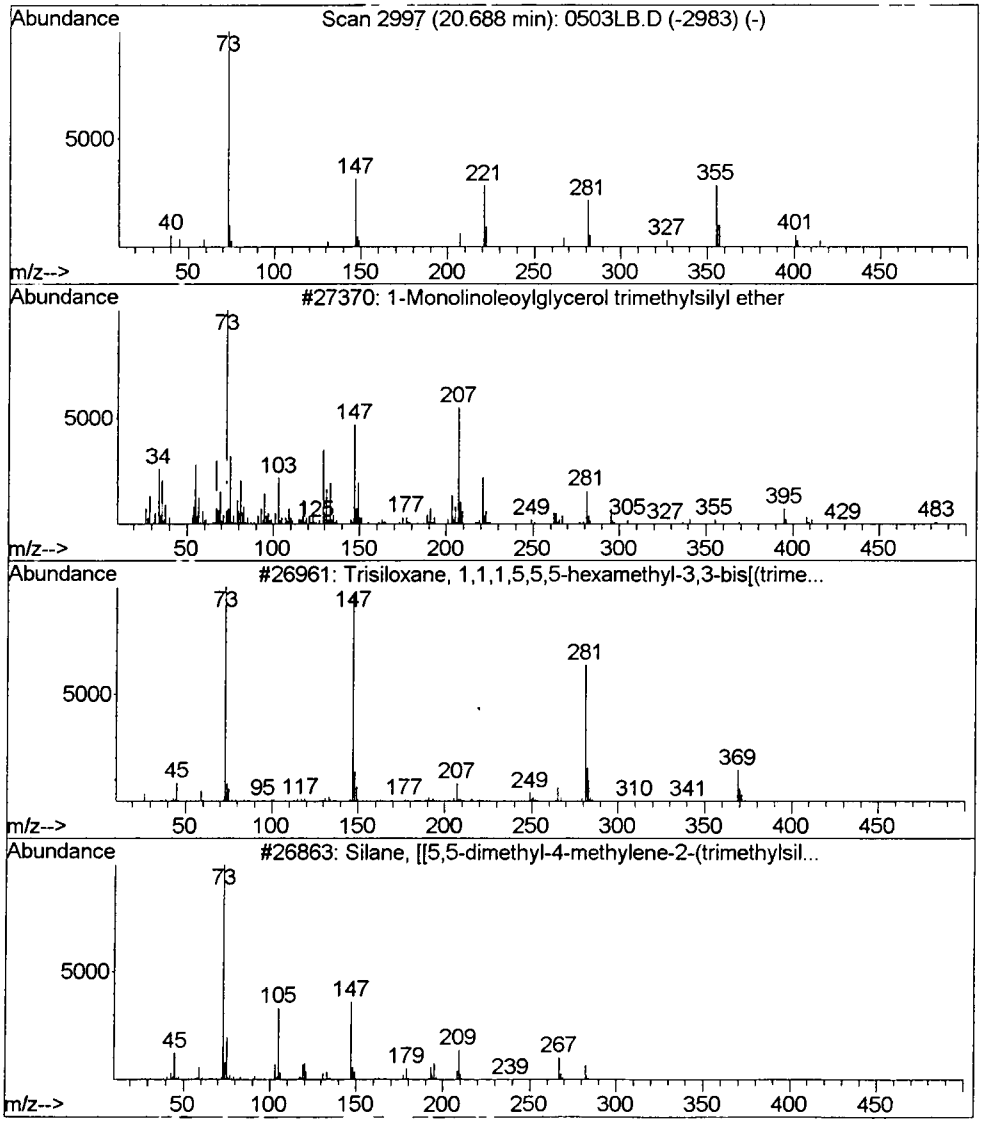
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

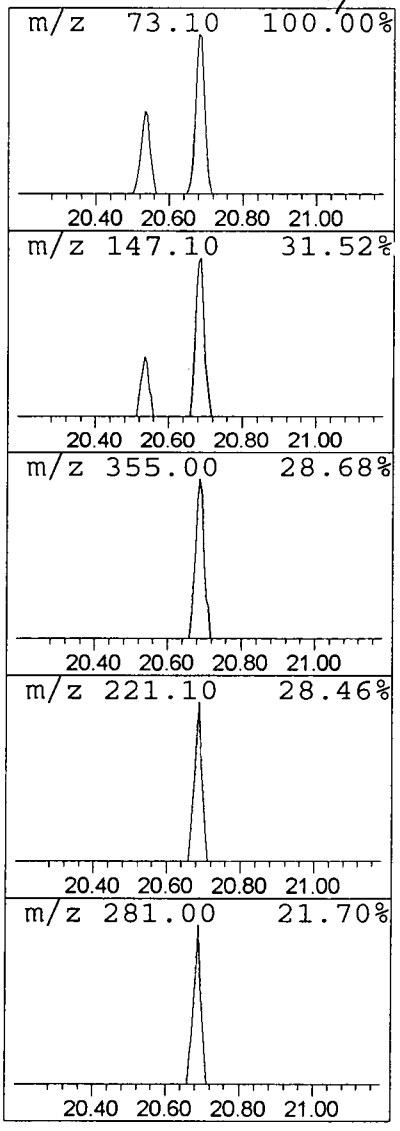
****

Peak Number 4 1-Monolinoleoylglycerol tri... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.69	0.21 ug/L	286472	Acenapthalene-d10	18.58

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Monolinoleoylglycerol trimethy...	498	C27H54O4Si2	054284-45-6	23
2	Trisiloxane, 1,1,1,5,5,5-hexamet...	384	C12H36O4Si5	003555-47-3	16
3	Silane, [[5,5-dimethyl-4-methyle...	282	C15H30OSi2	095798-15-5	14
4	Noradrenaline tetraTMS	457	C20H43NO3Si4	068595-65-3	12





Library Search Compound Report

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Acq On : 7 May 2002 4:24 pm Operator: Mclean
Sample : 5/3 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

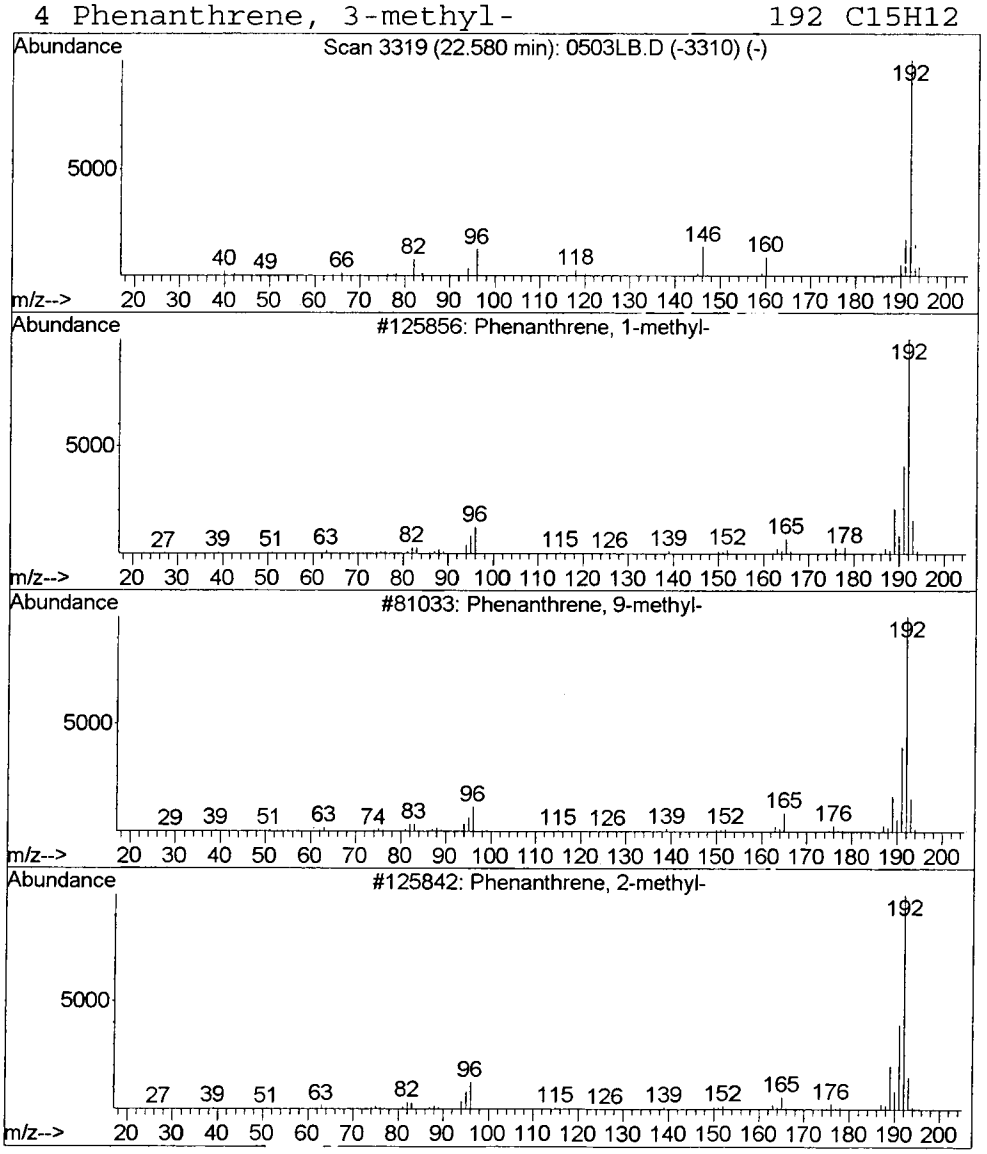
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

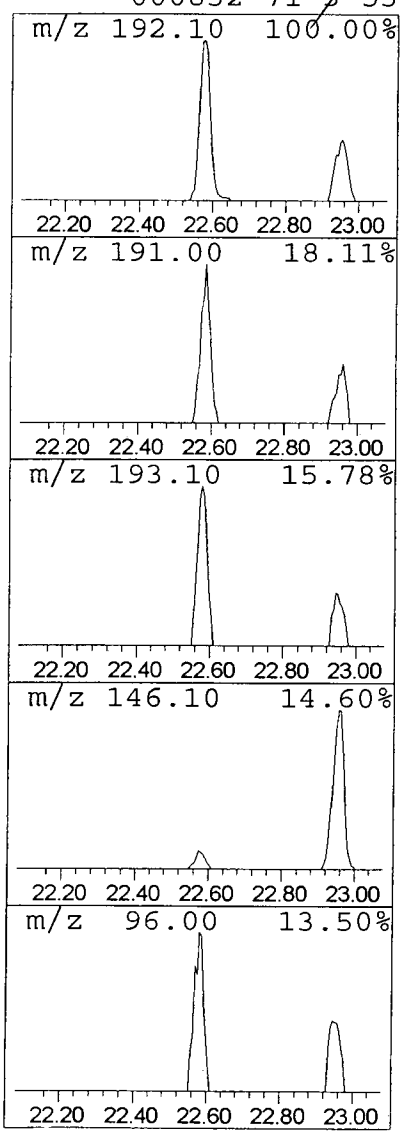
****

Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.58	0.29 ug/L	390053	Phenanthranene-d10	22.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	59
2	Phenanthrene, 9-methyl-	192	C15H12	000883-20-5	59
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	53
4	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	53





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050702\0503LB.D Vial: 5
Acq On : 7 May 2002 4:24 pm Operator: Mclean
Sample : 5/3 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

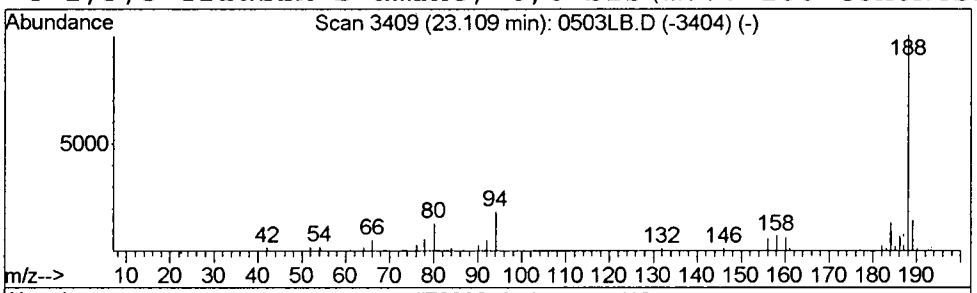
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

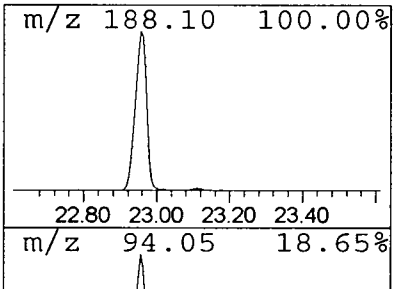
****

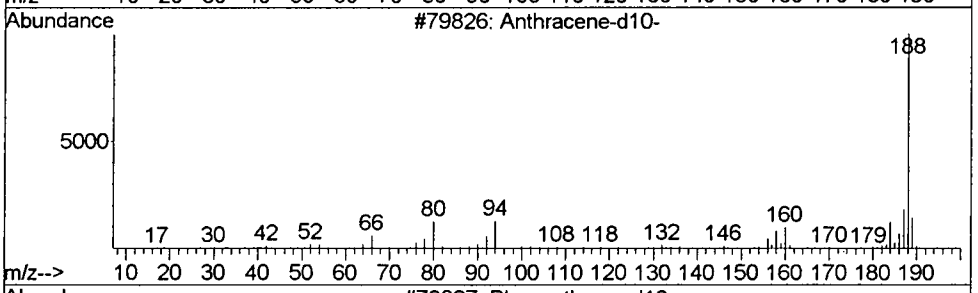
Peak Number 6 Anthracene-d10- Concentration Rank 2

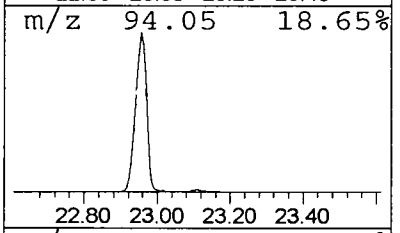
R.T.	EstConc	Area	Relative to ISTD	R.T.
23.11	0.46 ug/L	611831	Phenanthranene-d10	22.96

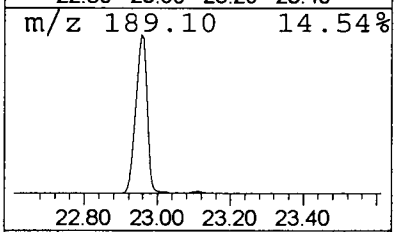
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene-d10-	188	C14D10	001719-06-8	93
2	Phenanthrene-d10	188	C14D10	001517-22-2	87
3	Antipyrine	188	C11H12N2O	000060-80-0	50
4	1,3,5-Triazin-2-amine, 4,6-bis(m...	188	C5H8N4S2	030358-19-1	38

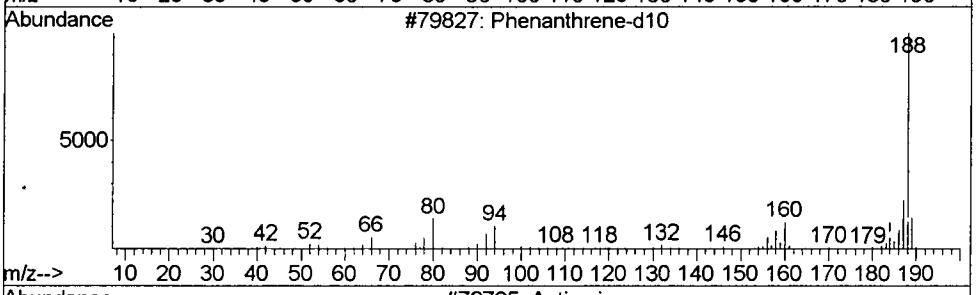


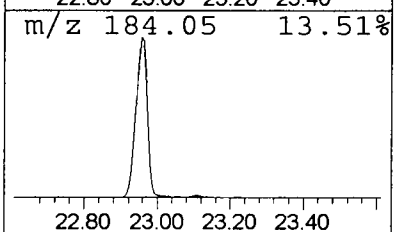


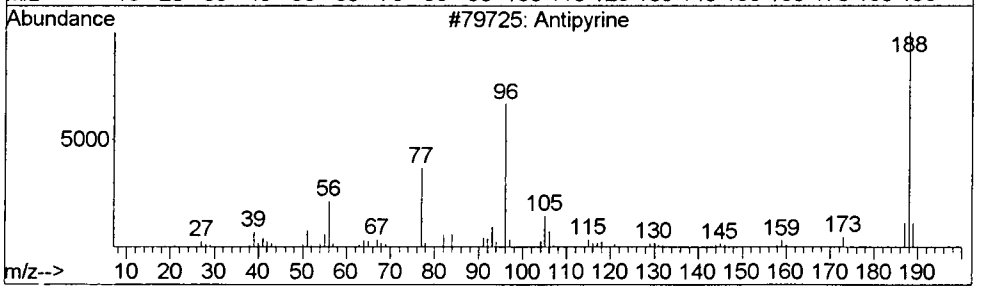


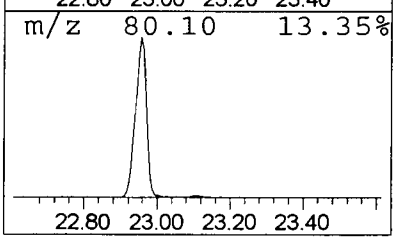












Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050702\0503LB.D Vial: 5
Acq On : 7 May 2002 4:24 pm Operator: Mclean
Sample : 5/3 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

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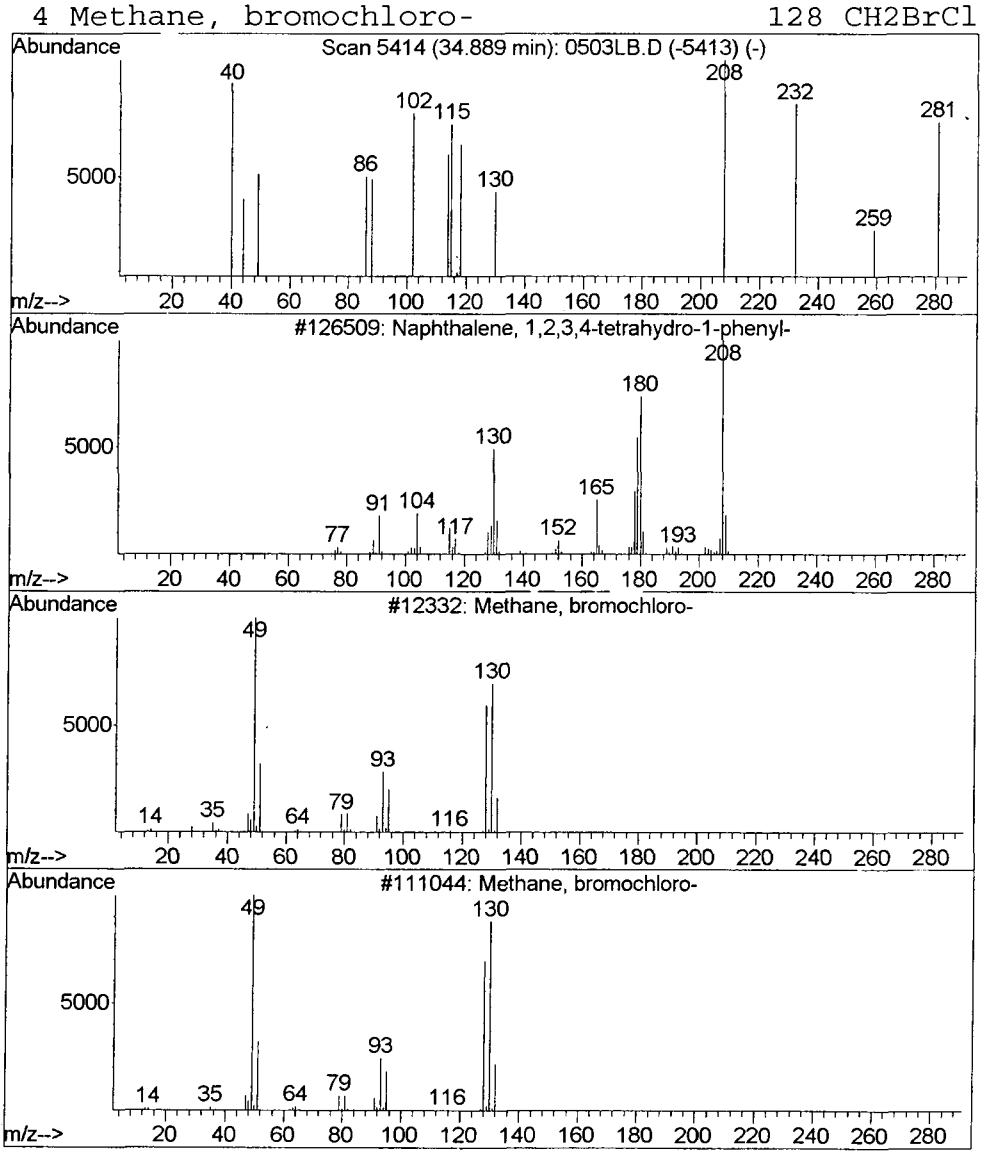
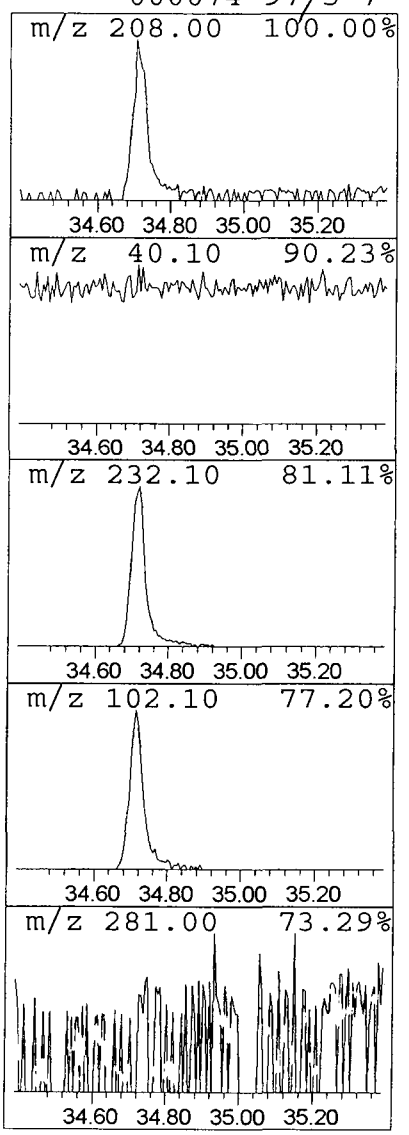
Peak Number 7 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.89	0.31 ug/L	284181	Perylene-d12	34.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	208	C16H16	003018-20-0	9
2	Methane, bromochloro-	128	CH2BrCl	000074-97-5	7
3	Methane, bromochloro-	128	CH2BrCl	000074-97-5	7
4	Methane, bromochloro-	128	CH2BrCl	000074-97-5	7





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050702\0503LB.D Vial: 5
Acq On : 7 May 2002 4:24 pm Operator: Mclean
Sample : 5/3 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

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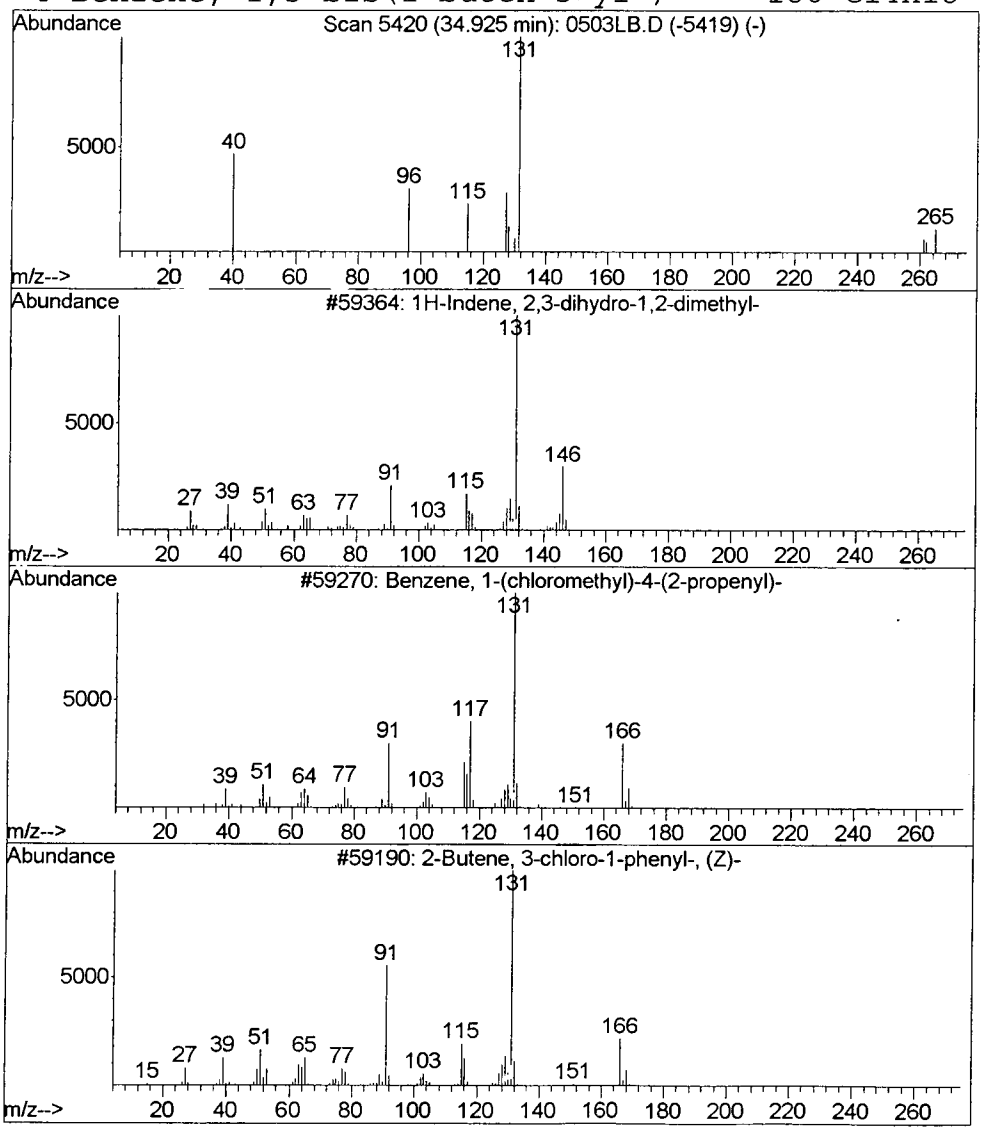
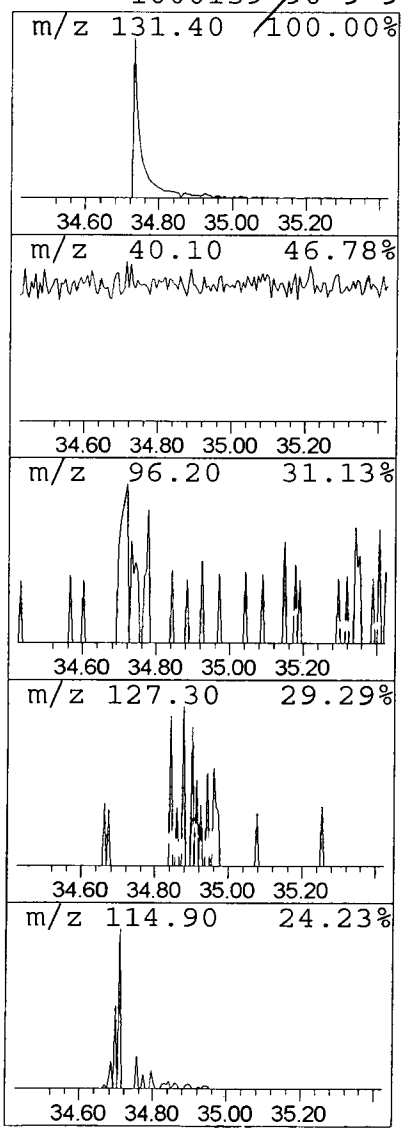
Peak Number 8 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.92	0.23 ug/L	215817	Perylene-d12	34.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	9
2	Benzene, 1-(chloromethyl)-4-(2-p...	166	C10H11Cl	036875-10-2	9
3	2-Butene, 3-chloro-1-phenyl-, (Z)-	166	C10H11Cl	016608-68-7	9
4	Benzene, 1,3-bis(1-buten-3-yl-)-	186	C14H18	1000159-90-9	9





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050702\0503LB.D Vial: 5
Acq On : 7 May 2002 4:24 pm Operator: Mclean
Sample : 5/3 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

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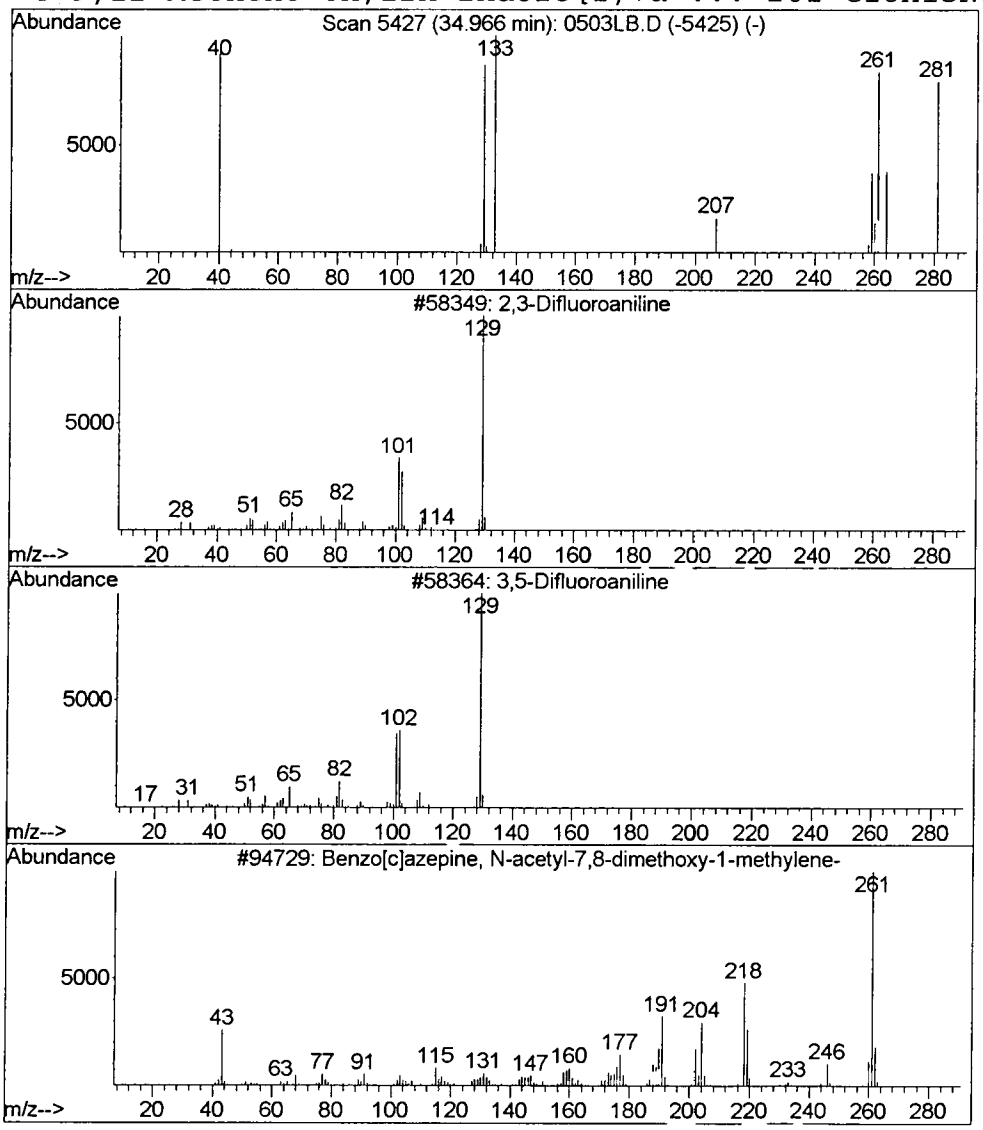
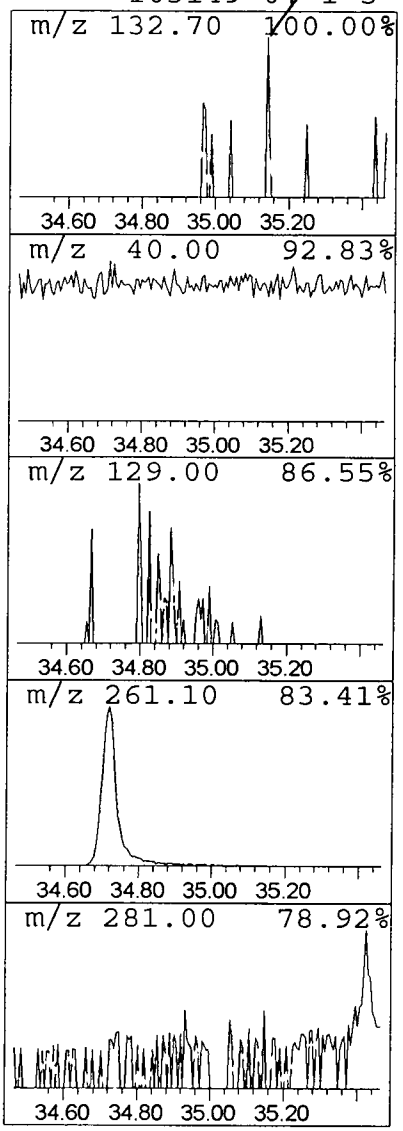
Peak Number 9 2,3-Difluoroaniline Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.96	0.21 ug/L	189888	Perylene-d12	34.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-Difluoroaniline	129	C6H5F2N	004519-40-8	7
2	3,5-Difluoroaniline	129	C6H5F2N	000372-39-4	7
3	Benzo[c]azepine, N-acetyl-7,8-di...	261	C15H19NO3	113193-86-5	5
4	9,12-Metheno-6H,12H-indolo[1,7a-...	261	C18H15NO	103149-07-1	5





Tentatively Identified Compound (LSC) summary

Operator ID: Mclean Date Acquired: 7 May 2002 4:24 pm
Data File: C:\MSDCHEM\1\DATA\050702\0503LB.D
Name: 5/3 kb
Misc:
Method: C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title: 508 Modified GC/MS scan
Library Searched: C:\DATABASE\NIST98.L

TIC Top Hit name	RT	EstConc	Units	Response	Internal Standard #	Internal Standard RT	Internal Standard Resp	Internal Standard Conc
Toluene	4.35	0.2	ug/L	228283	1	9.94	36711400	40.0
3-Penten-2-one, 4...	5.03	0.2	ug/L	228597	1	9.94	36711400	40.0
2-Pentanone, 4-hy...	5.98	8.9	ug/L	8128450	1	9.94	36711400	40.0
1-Monolinoleoylgl...	20.69	0.2	ug/L	286472	3	18.58	54145100	40.0
Phenanthrene, 1-m...	22.58	0.3	ug/L	390053	4	22.96	53033500	40.0
Anthracene-d10-	23.11	0.5	ug/L	611831	4	22.96	53033500	40.0
Naphthalene, 1,2,...	34.89	0.3	ug/L	284181	6	34.72	36911600	40.0
1H-Indene, 2,3-di...	34.92	0.2	ug/L	215817	6	34.72	36911600	40.0
2,3-Difluoroaniline	34.96	0.2	ug/L	189888	6	34.72	36911600	40.0