User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 05/15/2002 Time: 14:55:15

Sample: AE09331
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 5/15/02 14:54 Ended: 5/15/02 14:54 Analyst: DLV
Page: 1

	Component Name	Vial	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		79		10.0

Comments for Sample AE09331 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 05-15-2002 Time: 14:56:04

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds.
Recoveries for 4 of the 43 compounds in the LCS Standard were low.

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\050902\9331.D Vial: 21
Acq On : 10 May 2002 8:24 am Operator: Mclean
Sample : 5/8 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 13 11:42:05 2002 Quant Results File: 050102.RES

Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Mon May 13 11:40:29 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

re-extracted batch 6

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.94	152	8071019	40.00	ug/L	0.01
10) Napthalene-d8	13.43	136	32105407	40.00	ug/L	0.00
17) Acenapthalene-d10	18.57	164	17444422	40.00	ug/L	0.00
29) Phenanthranene-d10	22.96	188	25280967	40.00	ug/L	0.00
37) Chrysene-d12	30.81	240	17133735	40.00	ug/L	0.00
43) Perylene-d12	34.71	264	9562502	40.00	ug/L	0.00
System Monitoring Compounds						
27) TCMX	20.55	209	3349971	31.63	ug/L	0.00
Spiked Amount 40.000			Recovery = 79.07%			
Target Compounds						Qvalue
2) n-nitros-dimethyl amine	0.00	74	0	N.D.	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	0.00	93	0	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	20.55	77	61193	N.D.		
30) Nnitrosodiphenylamine	20.55	169	64012	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.		
34) Anthracene	0.00	178	0	N.D.		
35) Di-n-butylphthalate	25.08	149	42937	N.D.		

(#) = qualifier out of range (m) = manual integration
9331.D 050102.M Mon May 13 11:42:23 2002 Page 1

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\050902\9331.D Vial: 21
Acq On : 10 May 2002 8:24 am Operator: Mclean
Sample : 5/8 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 13 11:42:05 2002 Quant Results File: 050102.RES

Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Mon May 13 11:40:29 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

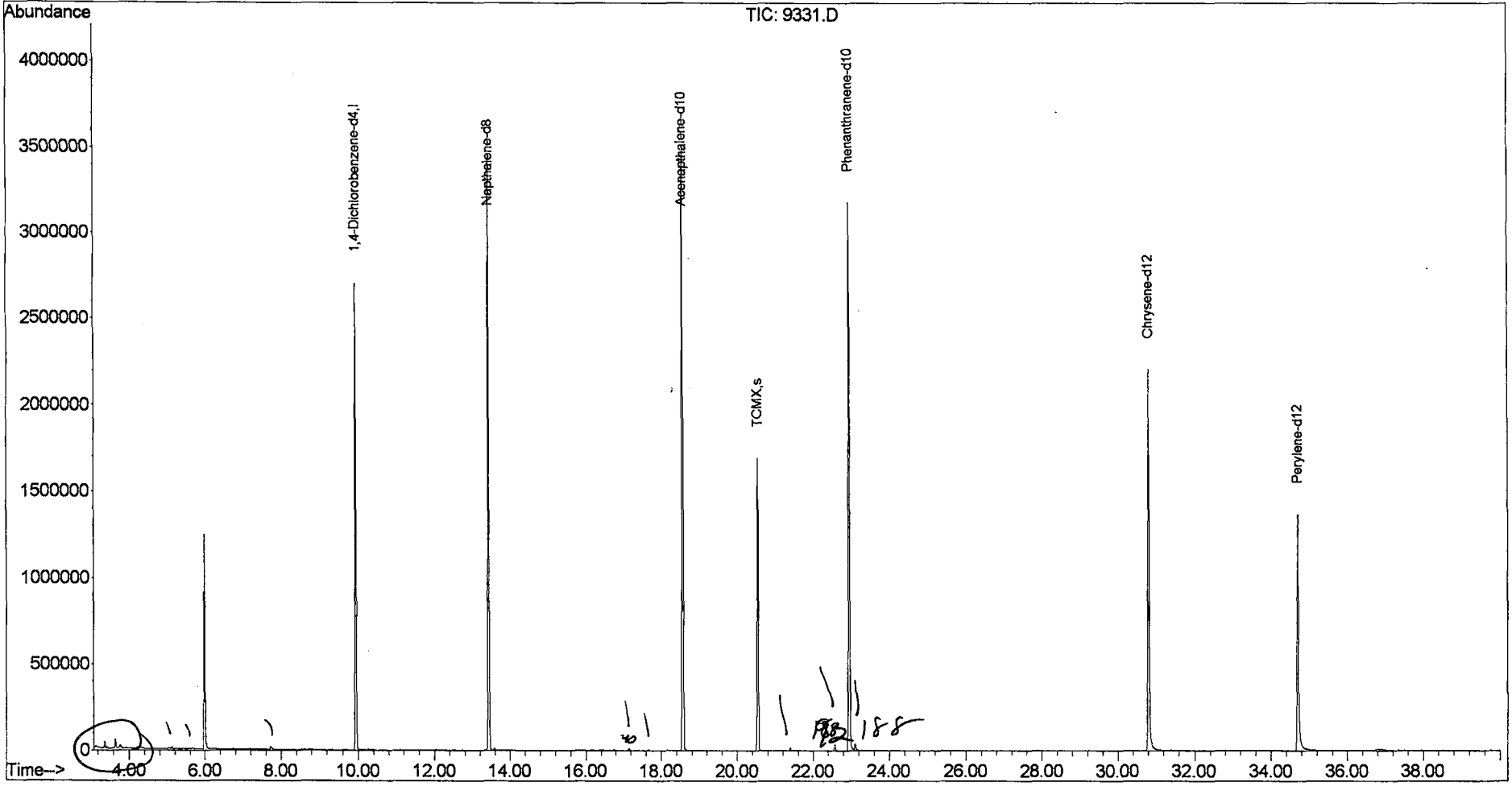
	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
36)	Fluoroanthene	0.00	202	0	N.D.	
38)	Pyrene	0.00	202	0	N.D.	
39)	Butylbenzylphthalate	0.00	149	0	N.D.	
40)	Benzo(a)anthracene	30.81	228	44599	N.D.	
41)	Bis(2-ethylhexyl)phthalate	0.00	149	0	N.D.	
42)	Chrysene	0.00	228	0	N.D.	
44)	Di-n-octylphthalate	0.00	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.	
50)	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\050902\9331.D Vial: 21
Acq On : 10 May 2002 8:24 am Operator: Mclean
Sample : 5/8 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 13 11:42 2002 Quant Results File: 050102.RES

Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Mon May 13 11:40:29 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\050902\9331.D Vial: 21
Acq On : 10 May 2002 8:24 am Operator: Mclean
Sample : 5/8 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

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

Signal : TIC

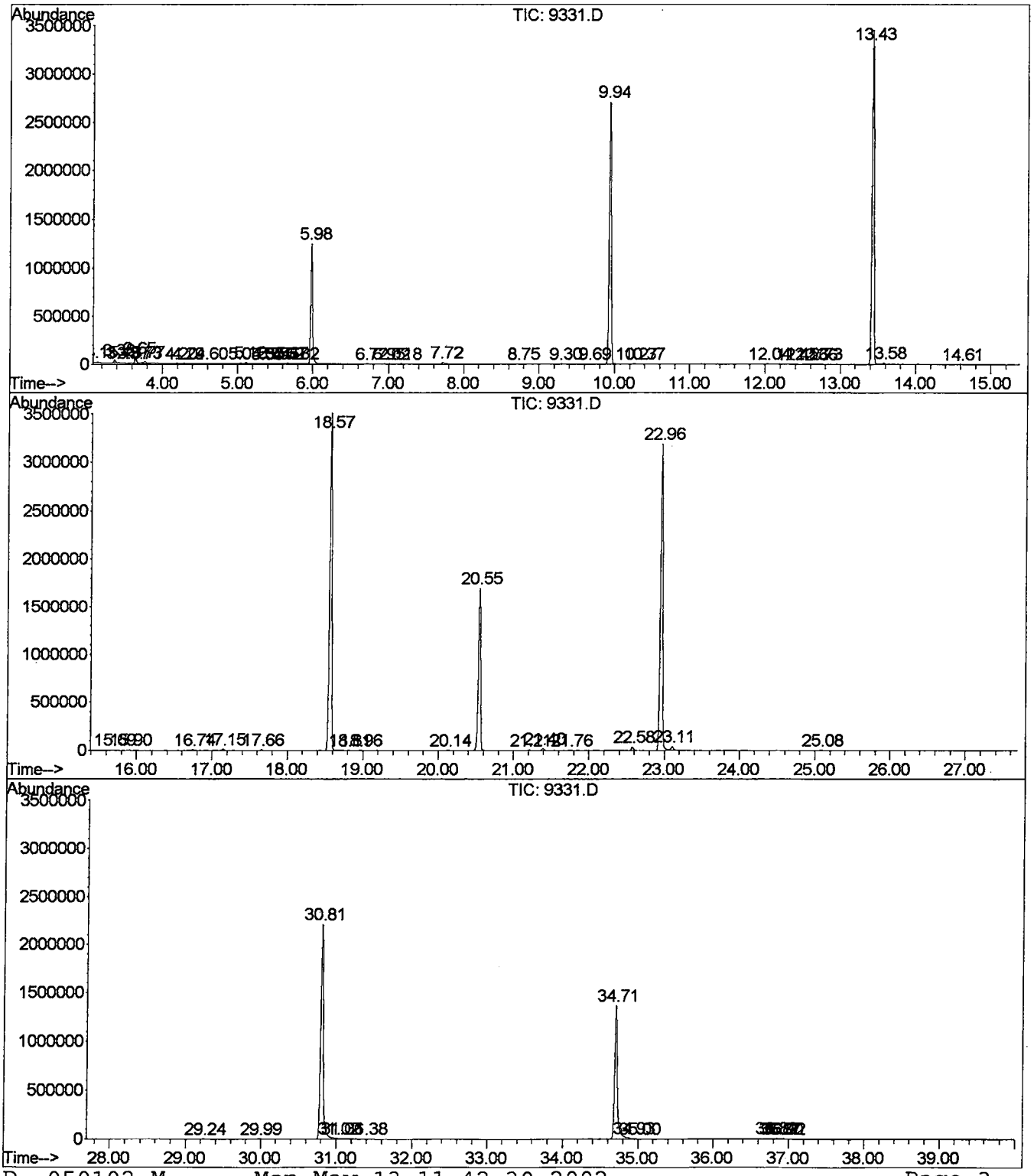
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.144	3	11	22	BV 3	6724	126846	0.18%	0.031%
2	3.379	45	51	56	PV	33692	523176	0.73%	0.128%
3	3.426	56	59	68	VV 6	5868	104342	0.15%	0.025%
4	3.567	76	83	91	PV 9	4503	116567	0.16%	0.028%
5	3.649	91	97	107	VV	53971	911669	1.27%	0.222%
6	3.732	107	111	114	VV 2	7922	135474	0.19%	0.033%
7	3.773	114	118	133	VV	18048	399038	0.56%	0.097%
8	4.196	181	190	196	VV 6	3849	89948	0.13%	0.022%
9	4.290	196	206	210	VV 4	4280	82192	0.11%	0.020%
10	4.601	255	259	267	PV 6	2002	45358	0.06%	0.011%
11	5.042	330	334	342	PV 4	6261	130483	0.18%	0.032%
12	5.118	342	347	356	VV 2	13907	241620	0.34%	0.059%
13	5.353	383	387	392	PV 6	1694	33131	0.05%	0.008%
14	5.435	397	401	406	VV 4	4078	70966	0.10%	0.017%
15	5.529	406	417	419	VV 8	2919	63542	0.09%	0.015%
16	5.547	419	420	424	VV 4	2614	41722	0.06%	0.010%
17	5.618	429	432	436	VV 5	3058	48611	0.07%	0.012%
18	5.676	436	442	450	VV 3	9105	188113	0.26%	0.046%
19	5.817	461	466	470	VV 4	2878	49377	0.07%	0.012%
20	5.982	480	494	521	VV	1207198	21372147	29.74%	5.212%
21	6.722	616	620	628	VV 7	2469	58646	0.08%	0.014%
22	6.951	655	659	665	VV 6	2114	42175	0.06%	0.010%
23	7.016	665	670	672	VV 4	2256	35556	0.05%	0.009%
24	7.175	688	697	705	PV 10	3522	85607	0.12%	0.021%
25	7.721	786	790	808	VV 2	16989	465662	0.65%	0.114%
26	8.749	960	965	974	VV 6	2911	59822	0.08%	0.015%
27	9.302	1055	1059	1069	VV 3	3861	88855	0.12%	0.022%
28	9.689	1122	1125	1127	VV 3	2192	23904	0.03%	0.006%
29	9.942	1158	1168	1189	VV	2660430	49412010	68.77%	12.051%
30	10.265	1217	1223	1231	VV 6	5443	145462	0.20%	0.035%
31	10.371	1234	1241	1253	VV 8	2983	93285	0.13%	0.023%
32	12.040	1519	1525	1537	VV 2	6041	117100	0.16%	0.029%
33	12.404	1583	1587	1593	VV 2	2039	28127	0.04%	0.007%
34	12.527	1593	1608	1619	PV 4	3496	75896	0.11%	0.019%
35	12.662	1627	1631	1635	PV 4	2025	24297	0.03%	0.006%
36	12.727	1635	1642	1647	VV	3798	65171	0.09%	0.016%
37	13.432	1751	1762	1781	PV	3403034	66326930	92.31%	16.176%

38	13.585	1781	1788	1801	VV		12544	225900	0.31%	0.055%
39	14.607	1958	1962	1972	VV	5	2344	45386	0.06%	0.011%
40	15.694	2140	2147	2155	VV	3	3883	74261	0.10%	0.018%
41	15.900	2176	2182	2188	VV	2	3710	66152	0.09%	0.016%
42	16.740	2321	2325	2336	VV	3	7704	143272	0.20%	0.035%
43	17.145	2378	2394	2400	PV	4	12686	226235	0.31%	0.055%
44	17.657	2476	2481	2487	PV	3	6195	88497	0.12%	0.022%
45	18.573	2624	2637	2659	PV		3496990	71855373	100.00%	17.525%
46	18.808	2675	2677	2687	VV	5	2388	49882	0.07%	0.012%
47	18.955	2696	2702	2714	VV	4	3332	81448	0.11%	0.020%
48	20.142	2901	2904	2911	VV	5	2984	51493	0.07%	0.013%
49	20.547	2960	2973	2987	PV		1692539	33655129	46.84%	8.208%
50	21.211	3069	3086	3093	VV	8	3335	88883	0.12%	0.022%
51	21.394	3111	3117	3130	VV	3	17817	327503	0.46%	0.080%
52	21.758	3173	3179	3189	VV	4	2670	65049	0.09%	0.016%
53	22.575	3309	3318	3331	PV		34441	622331	0.87%	0.152%
54	22.956	3370	3383	3402	PV	2	3143718	69426648	96.62%	16.932%
55	23.109	3402	3409	3424	VV		37380	891875	1.24%	0.218%
56	25.083	3738	3745	3753	PV	2	2846	55493	0.08%	0.014%
57	29.237	4440	4452	4460	PV	4	2394	52040	0.07%	0.013%
58	29.984	4567	4579	4590	PV	6	2541	81274	0.11%	0.020%
59	30.812	4704	4720	4754	PV	2	2201364	53837558	74.92%	13.130%
60	31.018	4754	4755	4760	VV	2	7483	124643	0.17%	0.030%
61	31.059	4760	4762	4767	VV	5	4888	113791	0.16%	0.028%
62	31.382	4810	4817	4823	PV	8	3334	72828	0.10%	0.018%
63	34.708	5367	5383	5419	PV	2	1368602	35160385	48.93%	8.575%
64	34.931	5419	5421	5432	VV	10	5675	167741	0.23%	0.041%
65	35.001	5432	5433	5440	VV	5	1919	26034	0.04%	0.006%
66	36.817	5721	5742	5749	PV	5	5598	279295	0.39%	0.068%
67	36.870	5749	5751	5753	VV	3	3478	46138	0.06%	0.011%
68	36.899	5753	5756	5759	VV	4	3480	57608	0.08%	0.014%
69	36.923	5759	5760	5765	VV	5	2691	41885	0.06%	0.010%

Sum of corrected areas: 410024853

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\050902\9331.D
Operator : Mclean
Acquired : 10 May 2002 8:24 am using AcqMethod 508SCAN
Instrument : Instrumen
Sample Name: 5/8 eh
Misc Info :
Vial Number: 21
Quant File :050102.RES (Chemstation Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050902\9331.D
Acq On : 10 May 2002 8:24 am
Sample : 5/8 eh
Misc :
MS Integration Params: LSCINT.e

Vial: 21
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

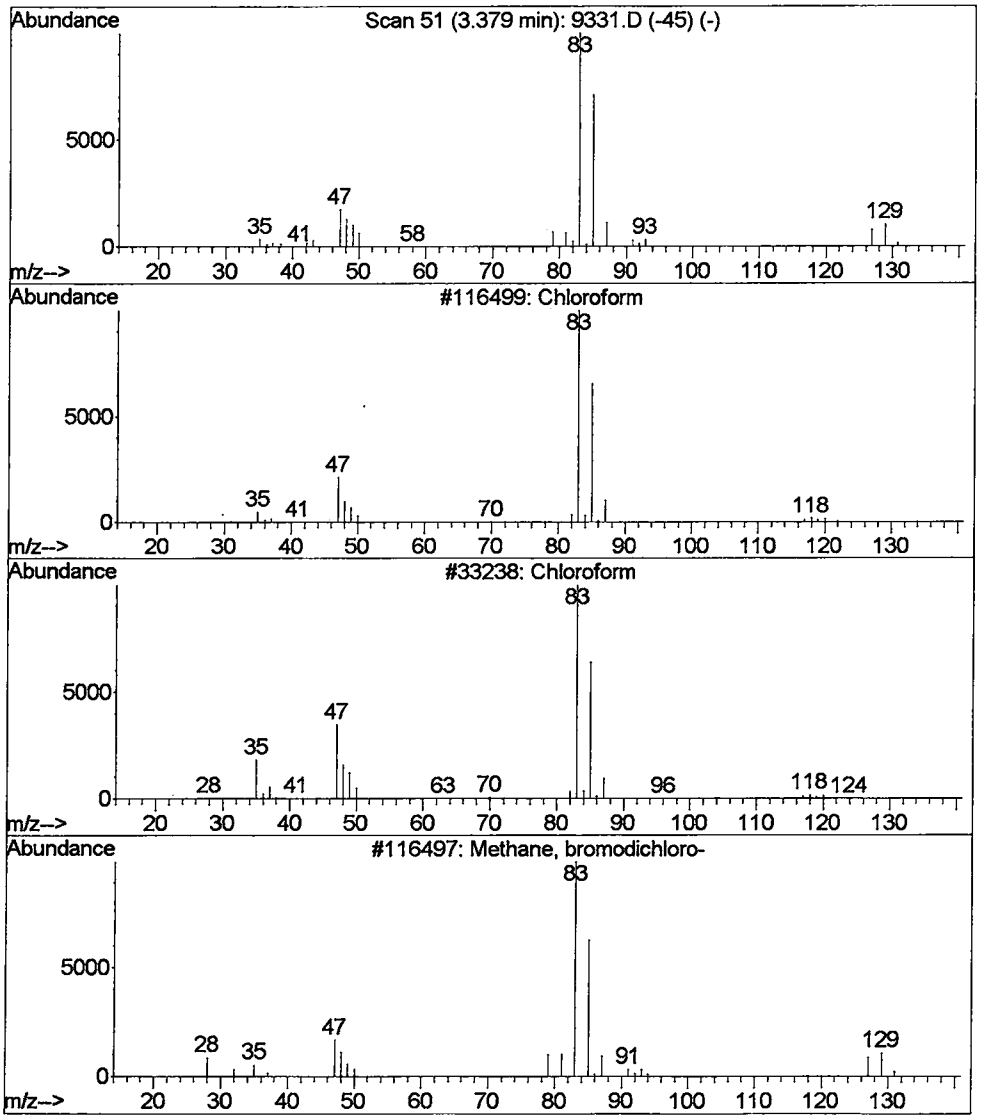
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

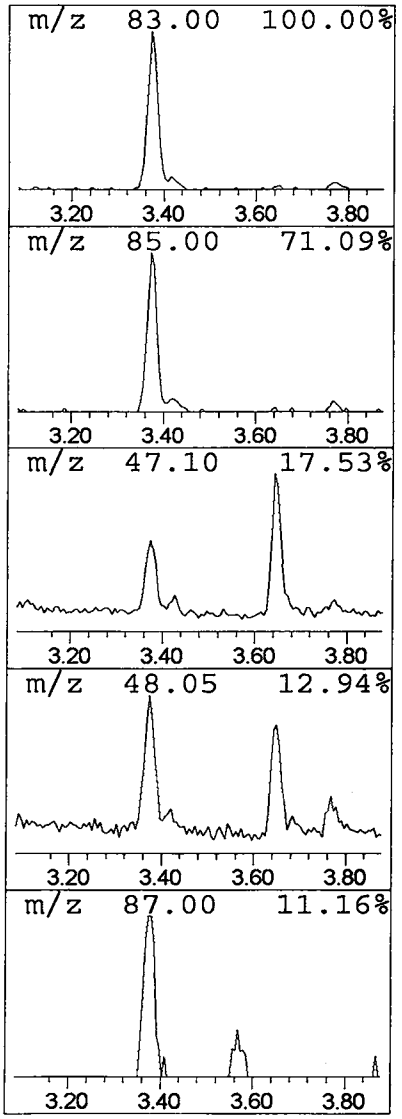
****

Peak Number 1 Chloroform Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.38	0.42 ug/L	523176	1,4-Dichlorobenzene-d4	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Chloroform	118	CHCl3	000067-66-3	78
2	Chloroform	118	CHCl3	000067-66-3	64
3	Methane, bromodichloro-	162	CHBrCl2	000075-27-4	59
4	Chloroform	118	CHCl3	000067-66-3	56





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050902\9331.D Vial: 21
Acq On : 10 May 2002 8:24 am Operator: Mclean
Sample : 5/8 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

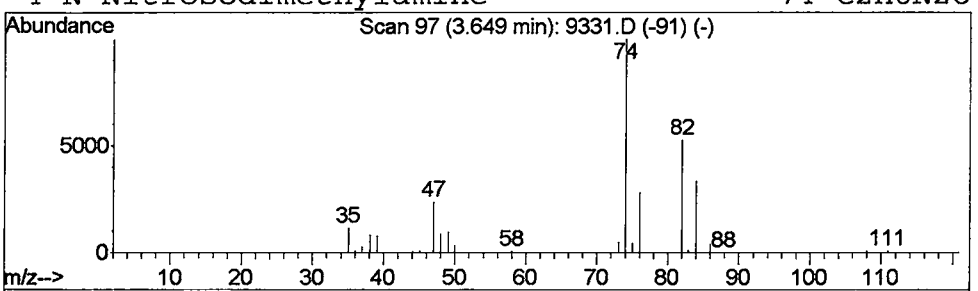
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

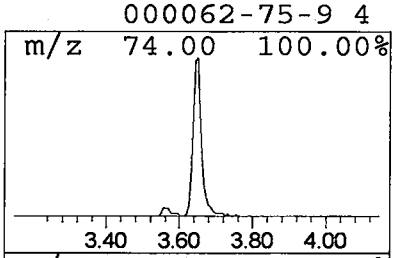
****

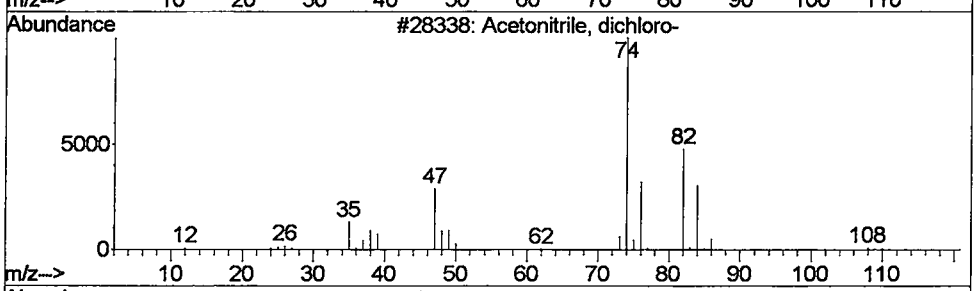
Peak Number 2 Acetonitrile, dichloro- Concentration Rank 2

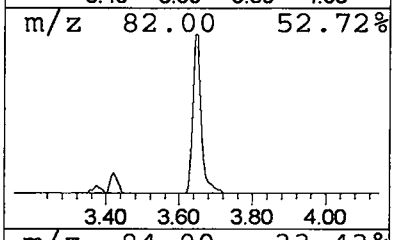
R.T.	EstConc	Area	Relative to ISTD	R.T.
3.65	0.74 ug/L	911669	1,4-Dichlorobenzene-d4	9.94

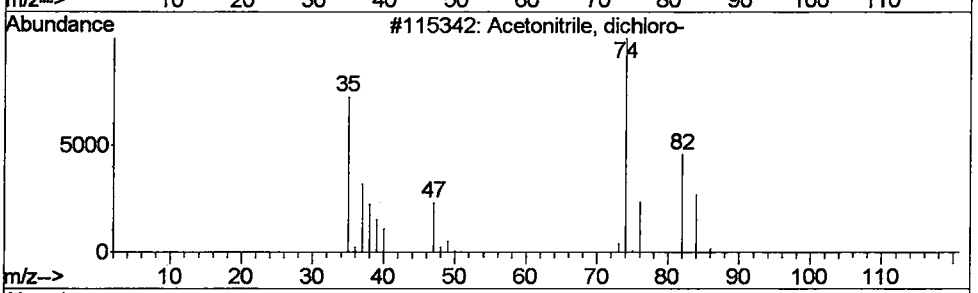
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	86
2	Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	64
3	3,3-Dichloropropyne	108	C3H2Cl2	1000222-23-9	38
4	N-Nitrosodimethylamine	74	C2H6N2O	000062-75-9	4

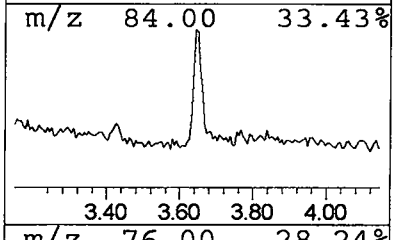


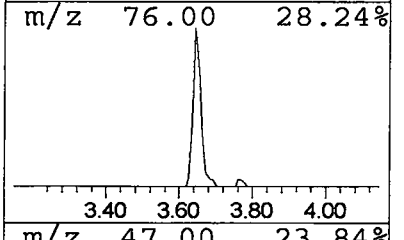


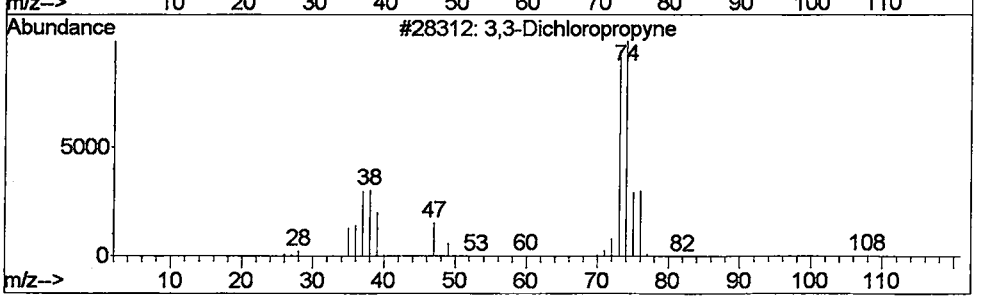


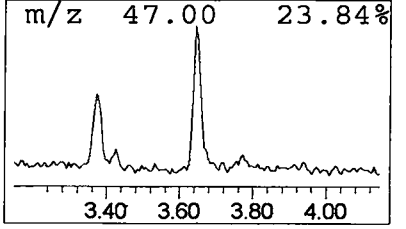












Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050902\9331.D
Acq On : 10 May 2002 8:24 am
Sample : 5/8 eh
Misc :
MS Integration Params: LSCINT.e

Vial: 21
Operator: Mclean
Inst : Instrumen
Multiplr: 1.00

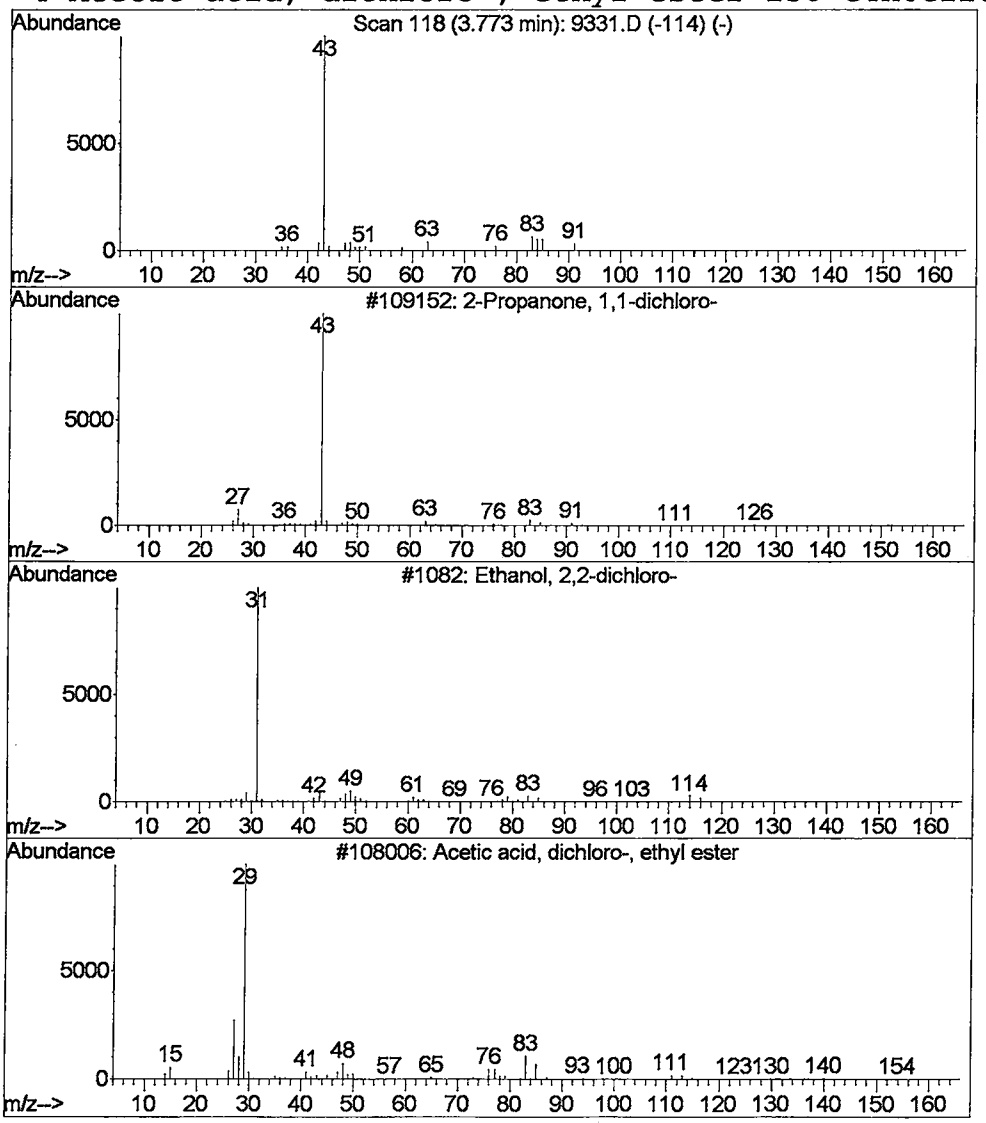
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

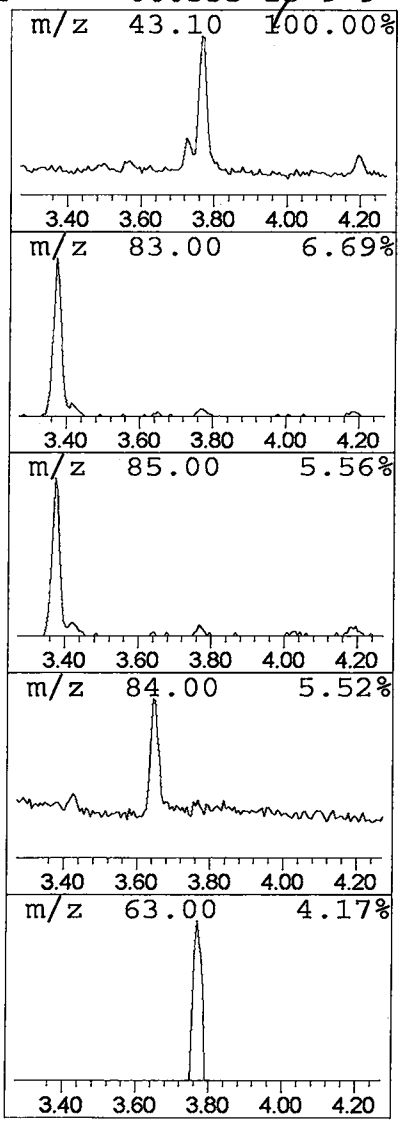
****

Peak Number 3 2-Propanone, 1,1-dichloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.77	0.32 ug/L	399038	1,4-Dichlorobenzene-d4	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanone, 1,1-dichloro-	126	C3H4Cl2O	000513-88-2	59
2	Ethanol, 2,2-dichloro-	114	C2H4Cl2O	000598-38-9	9
3	Acetic acid, dichloro-, ethyl ester	156	C4H6Cl2O2	000535-15-9	9
4	Acetic acid, dichloro-, ethyl ester	156	C4H6Cl2O2	000535-15-9	9





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050902\9331.D Vial: 21
Acq On : 10 May 2002 8:24 am Operator: Mclean
Sample : 5/8 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

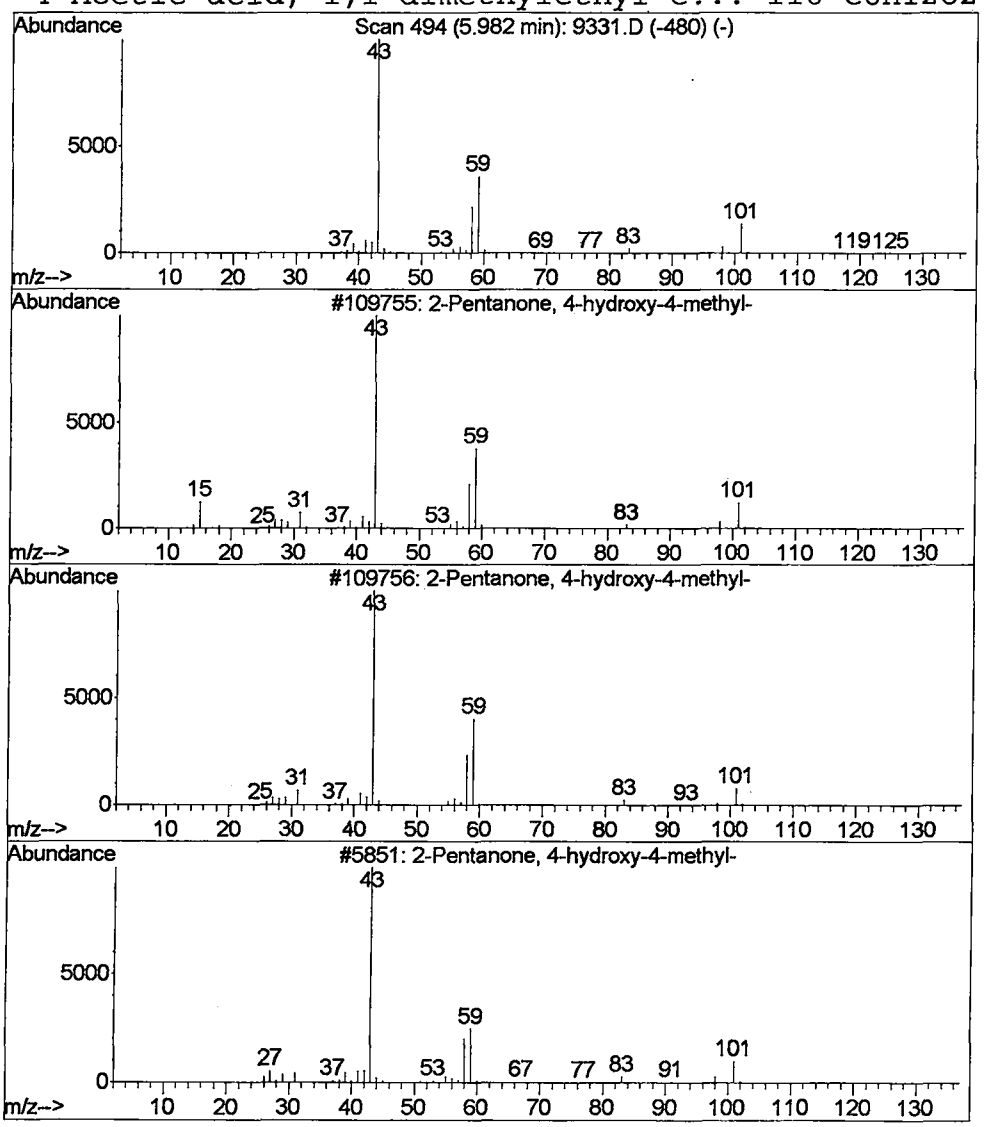
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

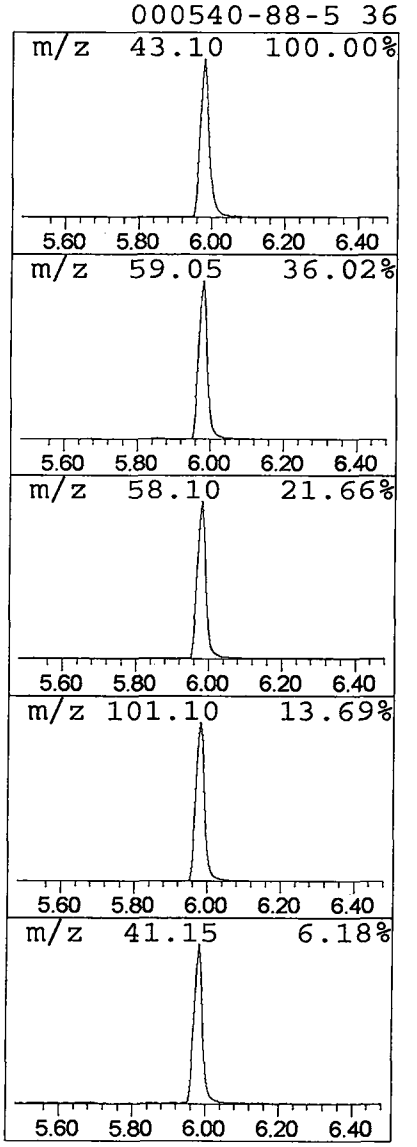
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.98	17.30 ug/L	21372100	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	90
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	36





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050902\9331.D Vial: 21
Acq On : 10 May 2002 8:24 am Operator: Mclean
Sample : 5/8 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

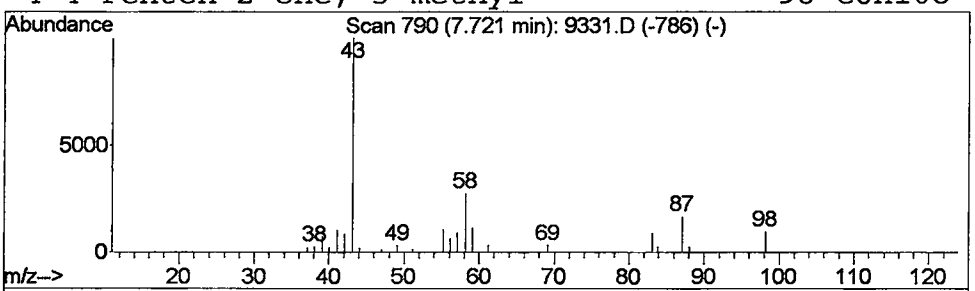
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

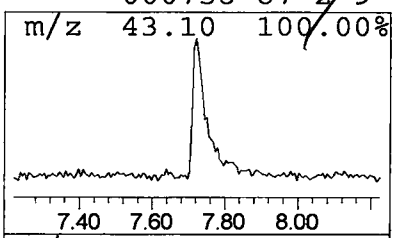
****

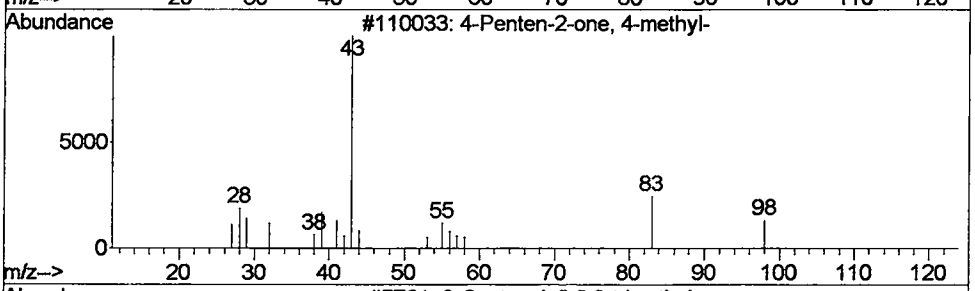
Peak Number 5 4-Penten-2-one, 4-methyl- Concentration Rank 5

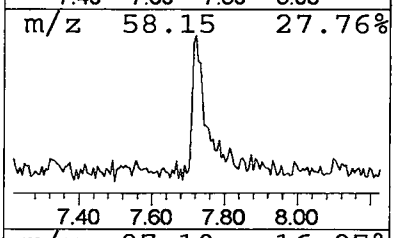
R.T.	EstConc	Area	Relative to ISTD	R.T.
7.72	0.38 ug/L	465662	1,4-Dichlorobenzene-d4	9.94

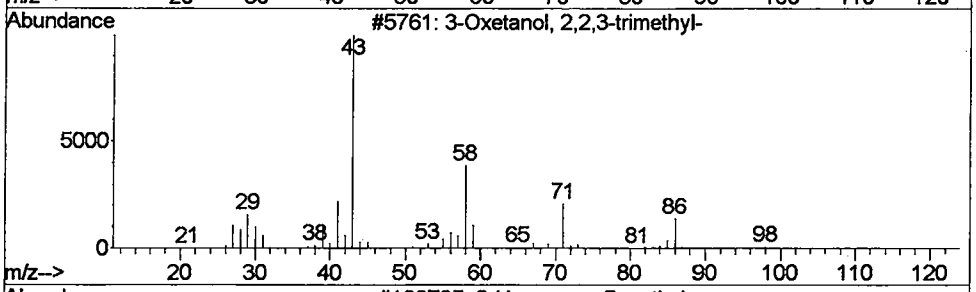
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	35
2	3-Oxetanol, 2,2,3-trimethyl-	116	C6H12O2	025910-96-7	28
3	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	28
4	4-Penten-2-one, 3-methyl-	98	C6H10O	000758-87-2	9

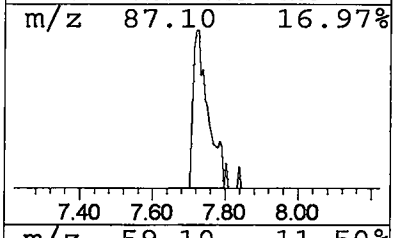


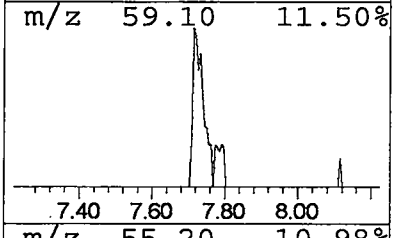


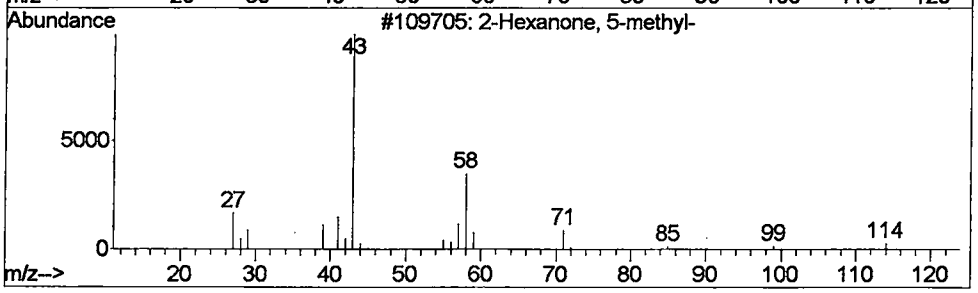


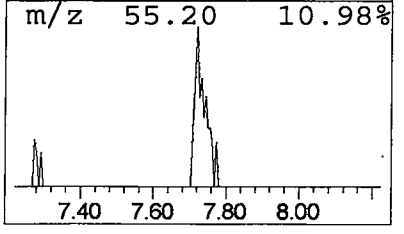












Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050902\9331.D Vial: 21
Acq On : 10 May 2002 8:24 am Operator: Mclean
Sample : 5/8 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

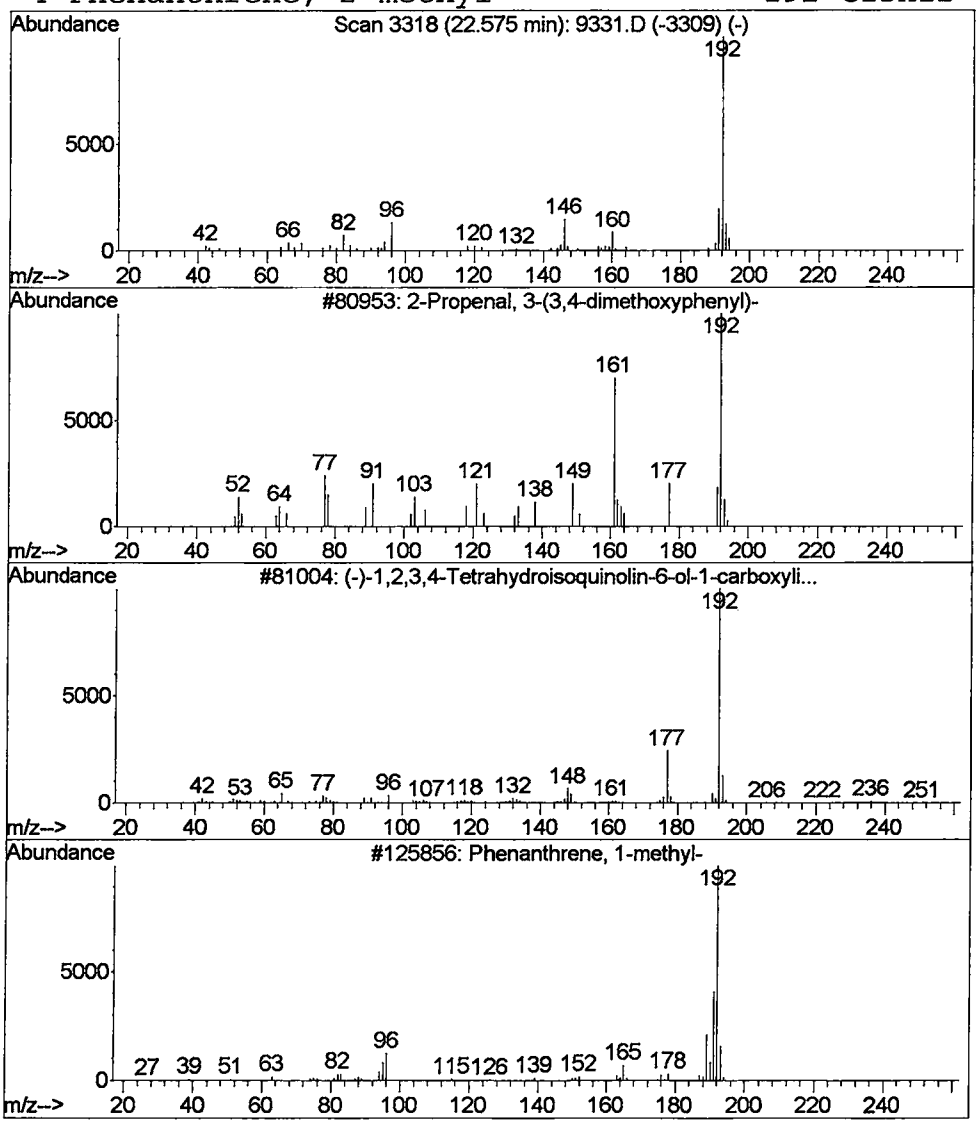
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

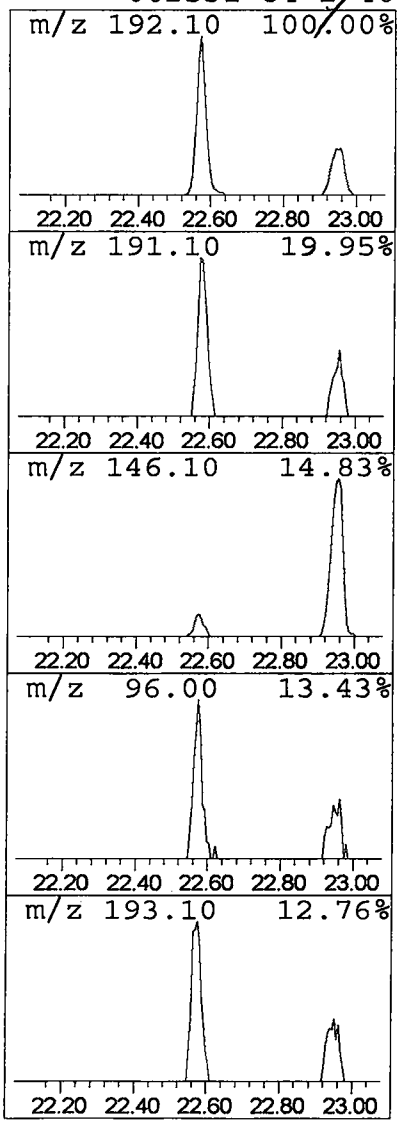
****

Peak Number 6 2-Propenal, 3-(3,4-dimethox... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.58	0.36 ug/L	622331	Phenanthranene-d10	22.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenal, 3-(3,4-dimethoxyphen...	192	C11H12O3	004497-40-9	45
2	(-)-1,2,3,4-Tetrahydroisoquinoli...	251	C13H17NO4	1000127-91-1	42
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	42
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	40





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050902\9331.D Vial: 21
Acq On : 10 May 2002 8:24 am Operator: Mclean
Sample : 5/8 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

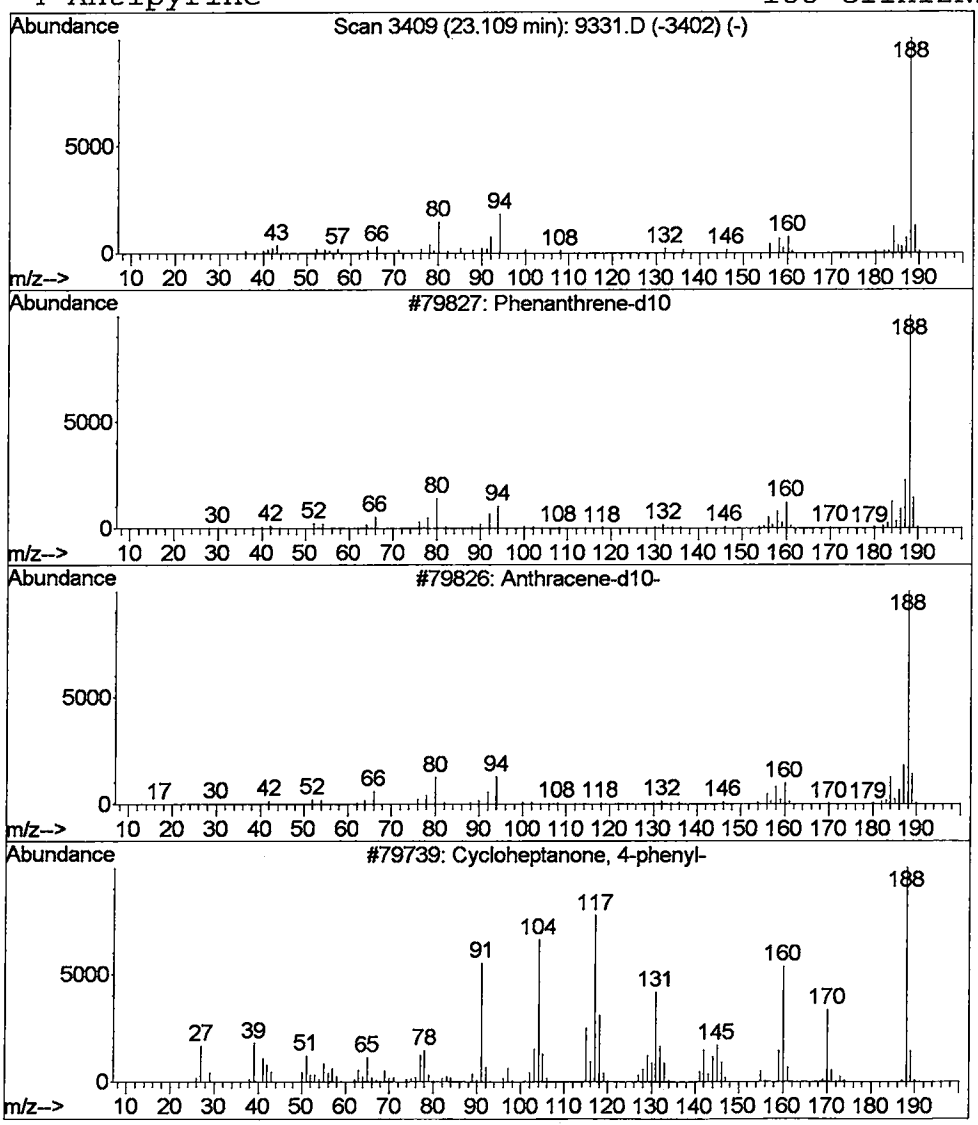
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

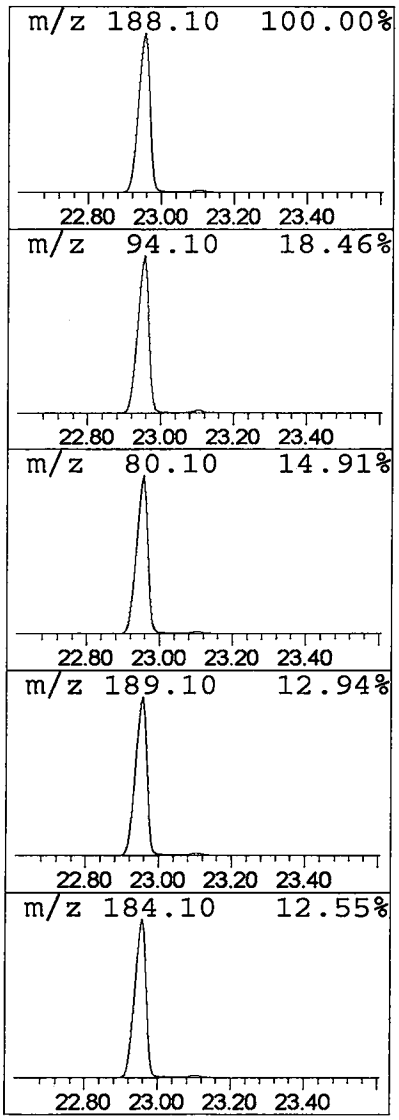
****

Peak Number 7 Phenanthrene-d10 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.11	0.51 ug/L	891875	Phenanthranene-d10	22.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene-d10	188	C14D10	001517-22-2	87
2	Anthracene-d10-	188	C14D10	001719-06-8	64
3	Cycloheptanone, 4-phenyl-	188	C13H16O	067688-28-2	50
4	Antipyrine	188	C11H12N2O	000060-80-0	50





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050902\9331.D Vial: 21
Acq On : 10 May 2002 8:24 am Operator: Mclean
Sample : 5/8 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

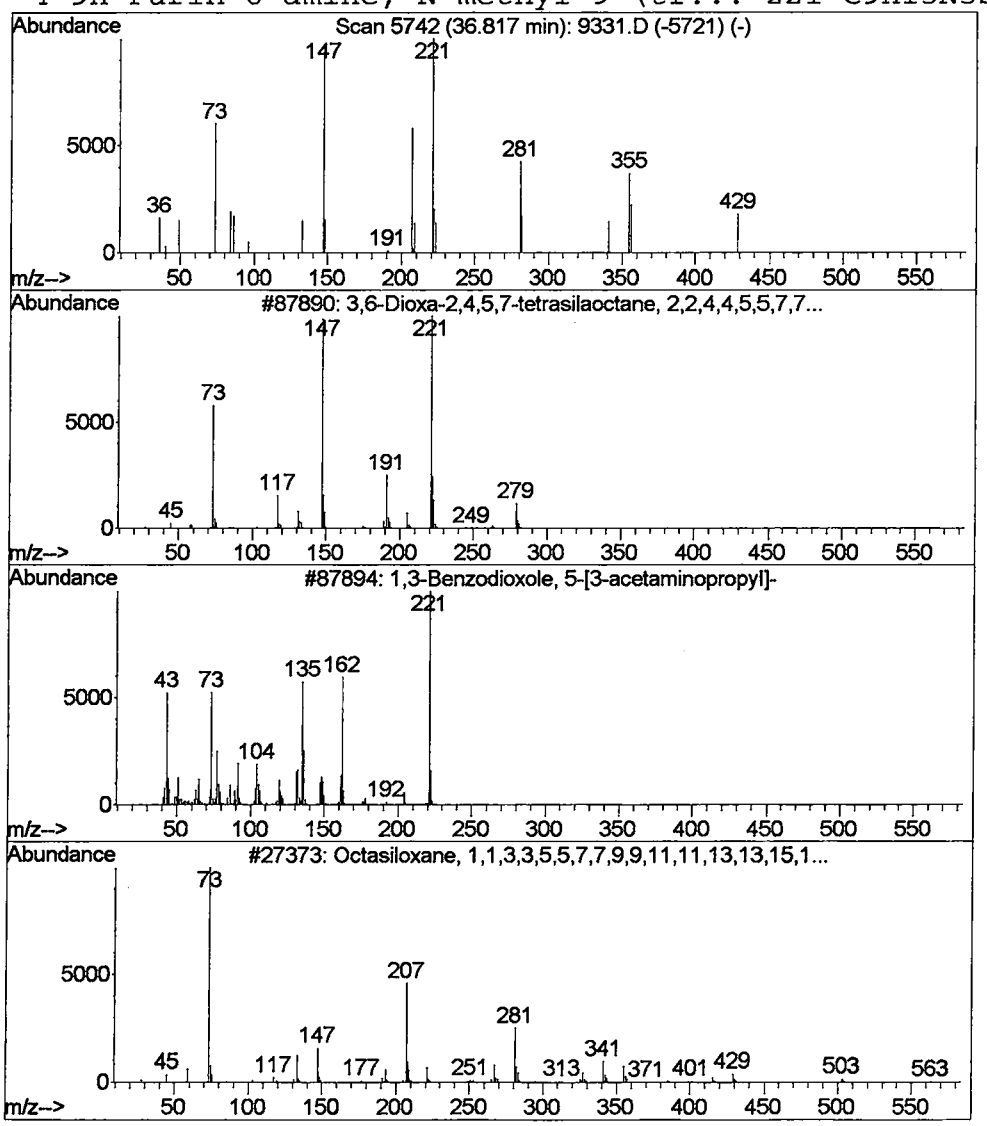
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

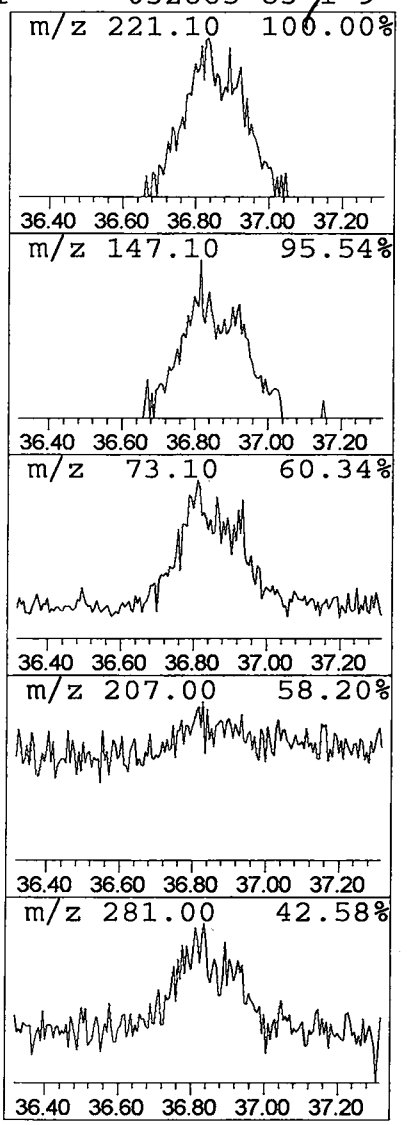
****

Peak Number 8 3,6-Dioxa-2,4,5,7-tetrasila... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
36.82	0.32 ug/L	279295	Perylene-d12	34.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	43
2	1,3-Benzodioxole, 5-[3-acetamino...	221	C12H15NO3	1000124-33-0	14
3	Octasiloxane, 1,1,3,3,5,5,7,7,9,...	578	C16H50O7Si8	019095-24-0	12
4	9H-Purin-6-amine, N-methyl-9-(tr...	221	C9H15N5Si	032865-83-1	9





Tentatively Identified Compound (LSC) summary

Operator ID: Mclean Date Acquired: 10 May 2002 8:24 am
Data File: C:\MSDCHEM\1\DATA\050902\9331.D
Name: 5/8 eh
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
Chloroform	3.38	0.4	ug/L	523176	1	9.94	49412000	40.0
Acetonitrile, dic...	3.65	0.7	ug/L	911669	1	9.94	49412000	40.0
2-Propanone, 1,1-...	3.77	0.3	ug/L	399038	1	9.94	49412000	40.0
2-Pentanone, 4-hy...	5.98	17.3	ug/L	21372100	1	9.94	49412000	40.0
4-Penten-2-one, 4...	7.72	0.4	ug/L	465662	1	9.94	49412000	40.0
2-Propenal, 3-(3,...	22.58	0.4	ug/L	622331	4	22.96	69426600	40.0
Phenanthrene-d10	23.11	0.5	ug/L	891875	4	22.96	69426600	40.0
3,6-Dioxa-2,4,5,7...	36.82	0.3	ug/L	279295	6	34.71	35160400	40.0