User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 05/10/2002 Time: 12:00:24

Sample: AE10203
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 5/10/02 11:59 Ended: 5/10/02 11:59 Analyst: DLV
Page: 1

	Component Name	Vial	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		88		10.0

Comments for Sample AE10203 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 05-10-2002 Time: 12:02:18

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds.
The recovery for 1 of the 43 compounds in the LCS Standard was low.

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\1\DATA\050802\10203.D Vial: 21
Acq On : 9 May 2002 7:42 am Operator: Mclean
Sample : 4/30 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 09 09:36:28 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	7511212	40.00	ug/L	0.00
10) Napthalene-d8	13.43	136	29922420	40.00	ug/L	0.00
17) Acenapthalene-d10	18.57	164	16315490	40.00	ug/L	0.00
29) Phenanthranene-d10	22.96	188	23738558	40.00	ug/L	0.00
37) Chrysene-d12	30.81	240	18467575	40.00	ug/L	0.00
43) Perylene-d12	34.71	264	12384289	40.00	ug/L	0.00
System Monitoring Compounds						
27) TCMX	20.54	209	828965	8.79	ug/L	0.00
Spiked Amount 10.000			Recovery = 87.90%			
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.		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	0.00	77	0	N.D.		
30) Nnitrosodiphenylamine	20.54	169	16612	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)

Data File : C:\MSDCHEM\1\DATA\050802\10203.D Vial: 21
Acq On : 9 May 2002 7:42 am Operator: Mclean
Sample : 4/30 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 09 09:36:28 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

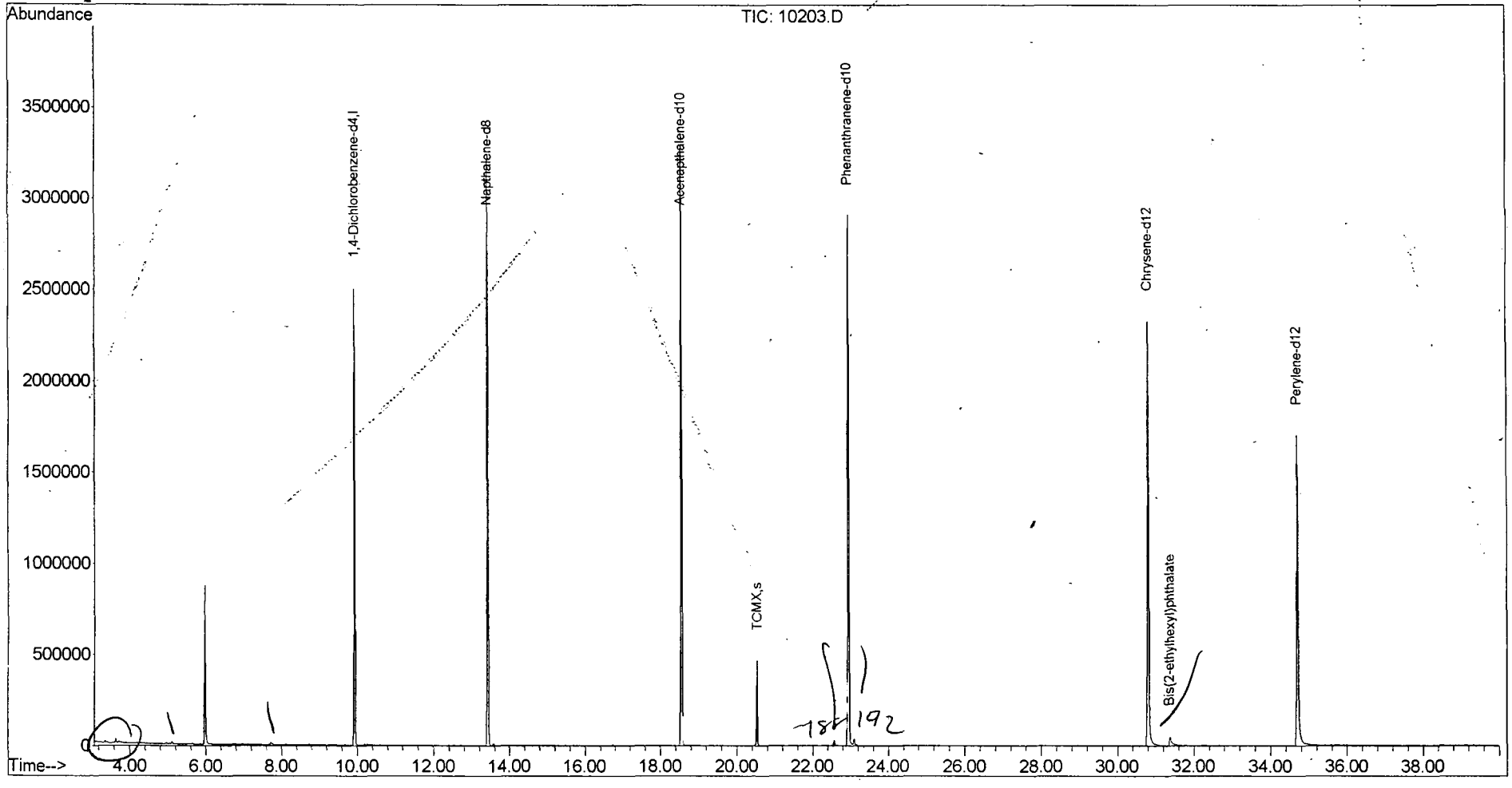
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
34) Anthracene	0.00	178	0	N.D.		
35) Di-n-butylphthalate	25.09	149	46783	N.D.		
36) Fluoroanthene	0.00	202	0	N.D.		
38) Pyrene	0.00	202	0	N.D.		
39) Butylbenzylphthalate	0.00	149	0	N.D.		
40) Benzo(a)anthracene	30.81	228	47246	N.D.		
41) Bis(2-ethylhexyl)phthalate	31.38	149	522584	1.14	ug/L	97
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\050802\10203.D Vial: 21
Acq On : 9 May 2002 7:42 am Operator: Mclean
Sample : 4/30 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 9 13:37 2002 Quant Results File: 050102.RES

Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Tue May 07 09:57:23 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\MSDCHEM\1\DATA\050802\10203.D
Acq On : 9 May 2002 7:42 am
Sample : 4/30 eh
Misc :
MS Integration Params: LSCINT.e

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

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

Signal : TIC

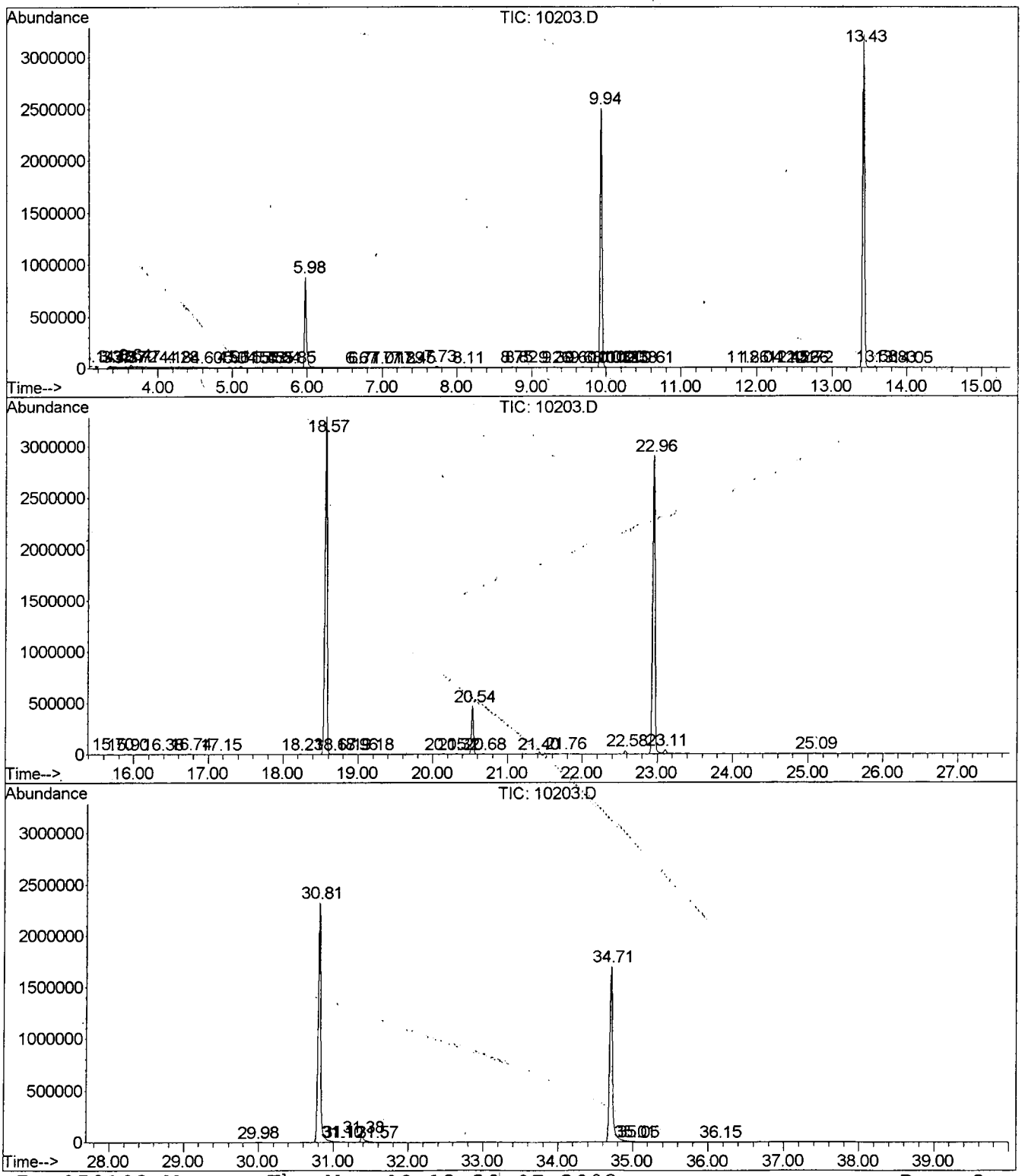
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.144	4	11	26	BV 3	5226	139002	0.21%	0.037%
2	3.367	42	49	53	PV 3	11329	168132	0.25%	0.045%
3	3.414	53	57	66	VV 8	4402	93930	0.14%	0.025%
4	3.567	77	83	87	PV 8	3513	66778	0.10%	0.018%
5	3.643	91	96	106	PV	23498	476773	0.71%	0.127%
6	3.725	106	110	114	VV 4	8826	174900	0.26%	0.046%
7	3.767	114	117	127	VV 2	8208	238984	0.36%	0.063%
8	4.178	183	187	201	VV 9	4588	163284	0.24%	0.043%
9	4.284	201	205	213	VV 5	5132	116071	0.17%	0.031%
10	4.601	254	259	262	VV 4	2512	46049	0.07%	0.012%
11	4.959	314	320	327	PV 4	7293	129365	0.19%	0.034%
12	5.042	331	334	341	VV 5	3186	59090	0.09%	0.016%
13	5.112	341	346	355	VV 5	12280	240949	0.36%	0.064%
14	5.429	394	400	405	VV 4	2797	49361	0.07%	0.013%
15	5.523	412	416	420	VV 5	3016	54329	0.08%	0.014%
16	5.553	420	421	423	VV 2	2333	24044	0.04%	0.006%
17	5.641	429	436	446	VV 8	4282	143538	0.21%	0.038%
18	5.852	467	472	476	VV 6	1986	30547	0.05%	0.008%
19	5.976	485	493	518	PV	845445	14923948	22.34%	3.965%
20	6.669	603	611	614	PV 6	1798	30650	0.05%	0.008%
21	6.716	614	619	622	VV 4	2219	39942	0.06%	0.011%
22	7.016	664	670	675	VV 7	1755	35700	0.05%	0.009%
23	7.180	694	698	709	VV 9	2233	47789	0.07%	0.013%
24	7.292	709	717	722	VV 7	1846	43189	0.06%	0.011%
25	7.445	736	743	752	VV 6	3360	69601	0.10%	0.018%
26	7.727	784	791	810	VV	14928	397846	0.60%	0.106%
27	8.109	850	856	860	VV 7	2243	40520	0.06%	0.011%
28	8.749	957	965	972	PV	7680	147032	0.22%	0.039%
29	8.814	972	976	984	VV 3	5089	104309	0.16%	0.028%
30	9.254	1045	1051	1054	VV 5	1920	33716	0.05%	0.009%
31	9.301	1054	1059	1071	VV 6	3331	88407	0.13%	0.023%
32	9.607	1106	1111	1119	VV 4	2523	71928	0.11%	0.019%
33	9.677	1119	1123	1131	VV 6	3539	102529	0.15%	0.027%
34	9.936	1158	1167	1181	VV	2455687	45837849	68.63%	12.178%
35	10.030	1181	1183	1187	VV 4	3138	46209	0.07%	0.012%
36	10.130	1197	1200	1206	VV 6	2548	46219	0.07%	0.012%
37	10.200	1206	1212	1217	VV	6695	121087	0.18%	0.032%

38	10.259	1217	1222	1232	VV	5	5569	135446	0.20%	0.036%
39	10.377	1236	1242	1248	VV	6	2968	66709	0.10%	0.018%
40	10.606	1277	1281	1287	VV	3	3228	52384	0.08%	0.014%
41	11.863	1487	1495	1507	PV	6	3868	80459	0.12%	0.021%
42	12.033	1513	1524	1533	VV	4	3746	84270	0.13%	0.022%
43	12.404	1578	1587	1592	PV	2	2436	40326	0.06%	0.011%
44	12.521	1599	1607	1615	BV	3	3989	72517	0.11%	0.019%
45	12.656	1626	1630	1636	VV	5	1894	34238	0.05%	0.009%
46	12.721	1636	1641	1648	VV		3327	75039	0.11%	0.020%
47	13.432	1751	1762	1780	PV		3194539	61390455	91.91%	16.309%
48	13.579	1780	1787	1798	VV		11321	242959	0.36%	0.065%
49	13.826	1825	1829	1836	VV	6	3093	51662	0.08%	0.014%
50	14.049	1858	1867	1881	VV	8	2003	50302	0.08%	0.013%
51	15.700	2143	2148	2156	VV	4	3386	65476	0.10%	0.017%
52	15.900	2178	2182	2192	VV	4	2680	47653	0.07%	0.013%
53	16.376	2257	2263	2273	VV	6	2388	53027	0.08%	0.014%
54	16.740	2316	2325	2337	VV	5	6781	153378	0.23%	0.041%
55	17.151	2390	2395	2398	VV	4	2065	31126	0.05%	0.008%
56	18.232	2574	2579	2582	PV	4	1741	27870	0.04%	0.007%
57	18.573	2615	2637	2652	PV		3263056	66793328	100.00%	17.745%
58	18.673	2652	2654	2667	VV	8	2127	83142	0.12%	0.022%
59	18.961	2696	2703	2712	VV	4	2916	71229	0.11%	0.019%
60	19.184	2733	2741	2747	VV	5	1674	24851	0.04%	0.007%
61	20.148	2901	2905	2910	VV	4	2520	45816	0.07%	0.012%
62	20.324	2929	2935	2942	VV	5	2812	55561	0.08%	0.015%
63	20.535	2961	2971	2988	VV		459959	8396425	12.57%	2.231%
64	20.682	2988	2996	3001	VV	5	2378	46216	0.07%	0.012%
65	21.393	3100	3117	3129	PV	6	3052	82417	0.12%	0.022%
66	21.758	3170	3179	3184	VV	4	4726	87677	0.13%	0.023%
67	22.580	3312	3319	3329	VV	3	31041	581119	0.87%	0.154%
68	22.956	3370	3383	3402	VV	2	2884327	65404589	97.92%	17.376%
69	23.109	3402	3409	3426	VV	2	39103	1071674	1.60%	0.285%
70	25.089	3737	3746	3757	PV	2	4296	90211	0.14%	0.024%
71	29.977	4571	4578	4586	PV	6	2224	55001	0.08%	0.015%
72	30.812	4708	4720	4766	PV	2	2328955	57815110	86.56%	15.360%
73	31.094	4766	4768	4771	VV	4	4335	76243	0.11%	0.020%
74	31.123	4771	4773	4780	VV	6	3163	74637	0.11%	0.020%
75	31.382	4807	4817	4848	PV		50234	1953951	2.93%	0.519%
76	31.576	4848	4850	4852	VV	3	1609	13786	0.02%	0.004%
77	34.713	5371	5384	5434	PV	2	1698065	45991970	68.86%	12.218%
78	35.013	5434	5435	5440	VV	3	4143	75525	0.11%	0.020%
79	35.054	5440	5442	5459	VV	7	2686	106863	0.16%	0.028%
80	36.147	5625	5628	5632	PV	4	1609	16396	0.02%	0.004%

Sum of corrected areas: 376412629

LSC Report - Integrated Chromatogram

File : C:\MSDCHEM\1\DATA\050802\10203.D
Operator : Mclean
Acquired : 9 May 2002 7:42 am using AcqMethod 508SCAN
Instrument : Instrumen
Sample Name: 4/30 eh
Misc Info :
Vial Number: 21
Quant File :050102.RES (Chemstation Integrator)



Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050802\10203.D Vial: 21
Acq On : 9 May 2002 7:42 am Operator: Mclean
Sample : 4/30 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

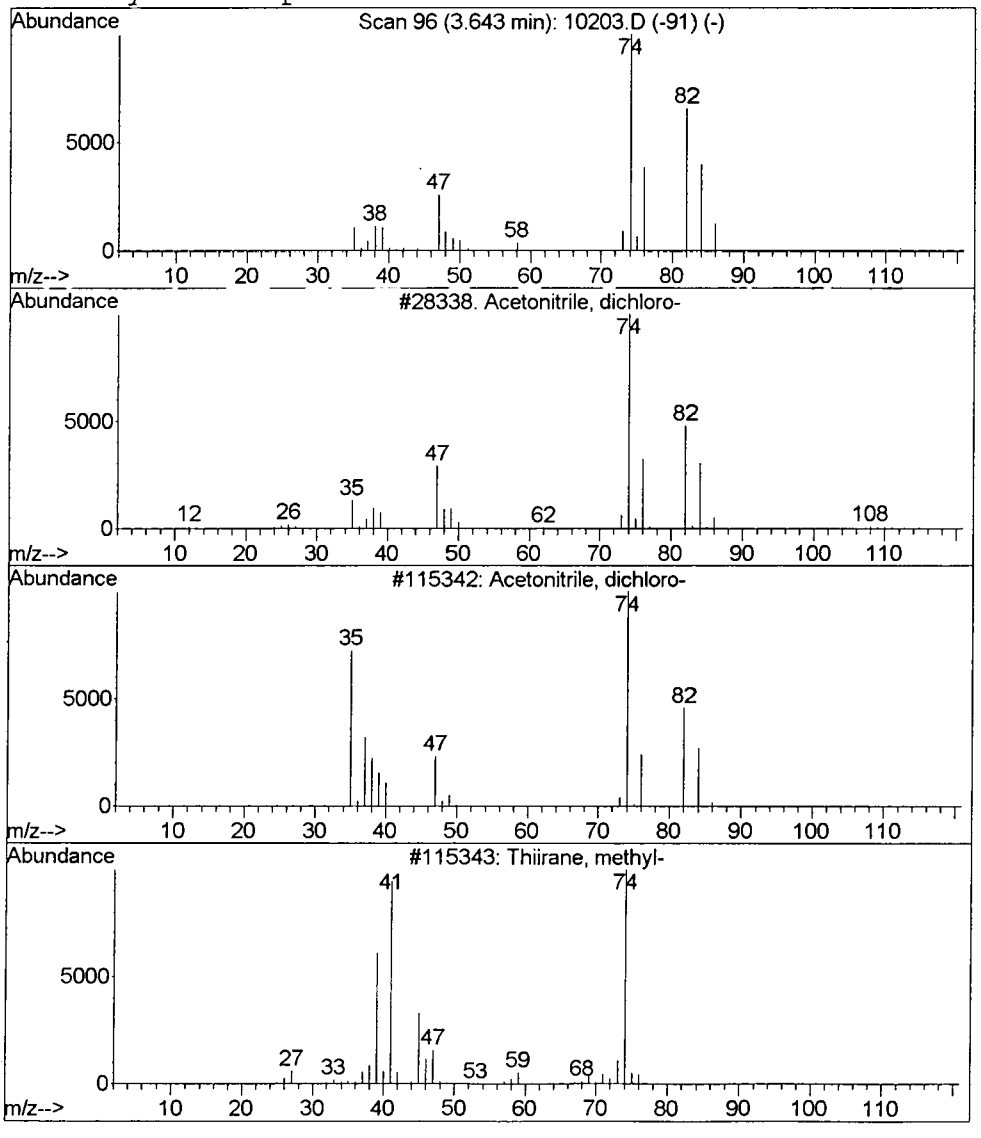
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

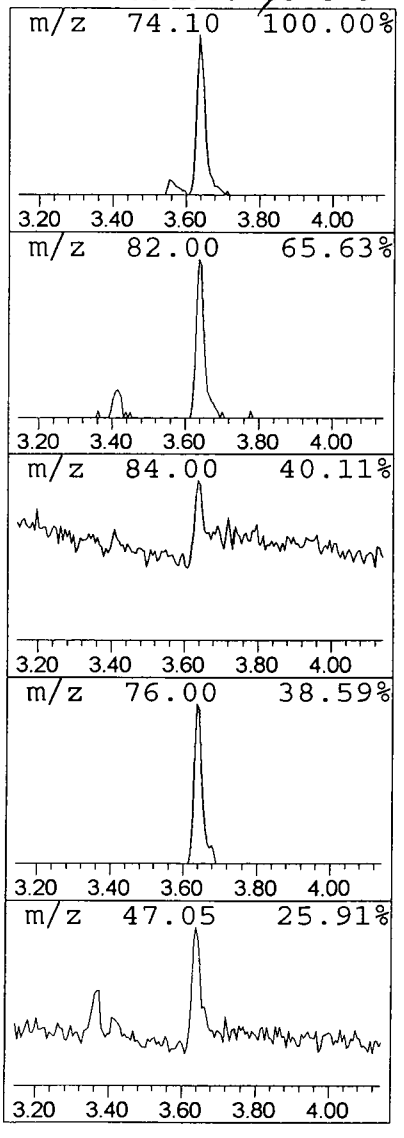
****

Peak Number 1 Acetonitrile, dichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.64	0.42 ug/L	476773	1,4-Dichlorobenzene-d4	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	72
2	Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	56
3	Thiirane, methyl-	74	C3H6S	001072-43-1	9
4	Allyl mercaptan	74	C3H6S	000870-23-5	9





Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050802\10203.D Vial: 21
Acq On : 9 May 2002 7:42 am Operator: Mclean
Sample : 4/30 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

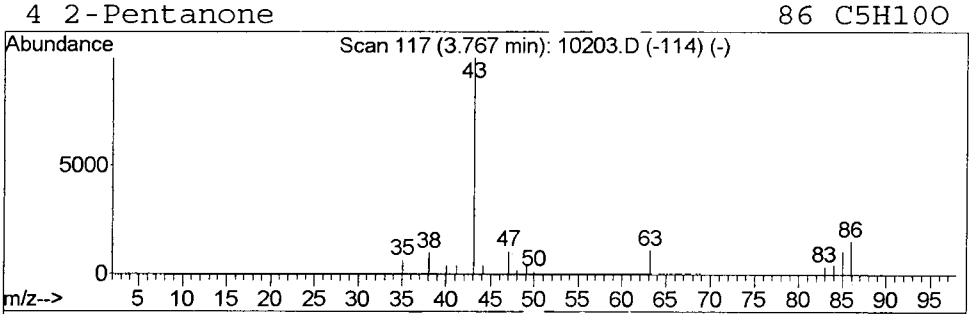
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

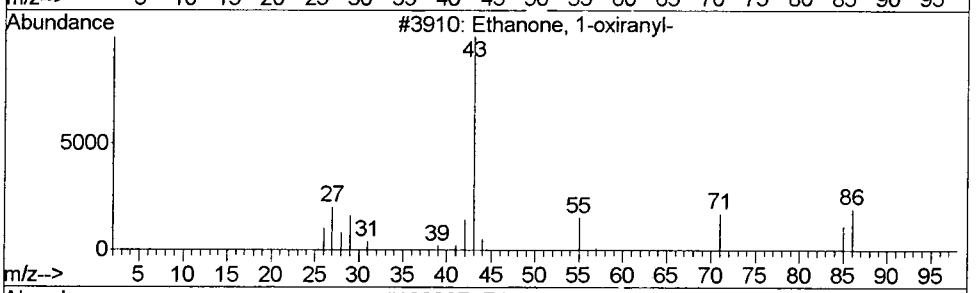
****

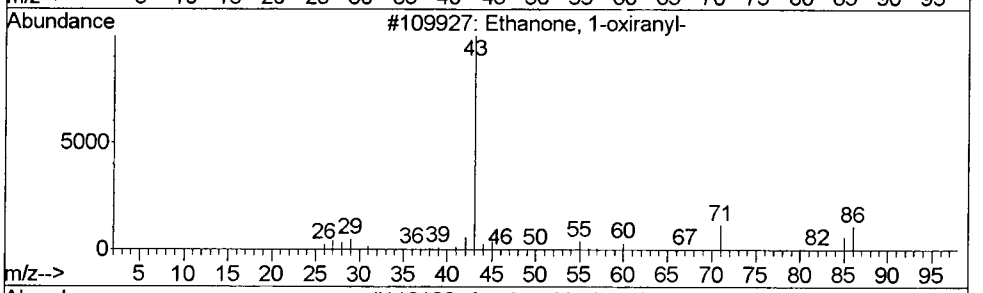
Peak Number 2 Ethanone, 1-oxiranyl- Concentration Rank 7

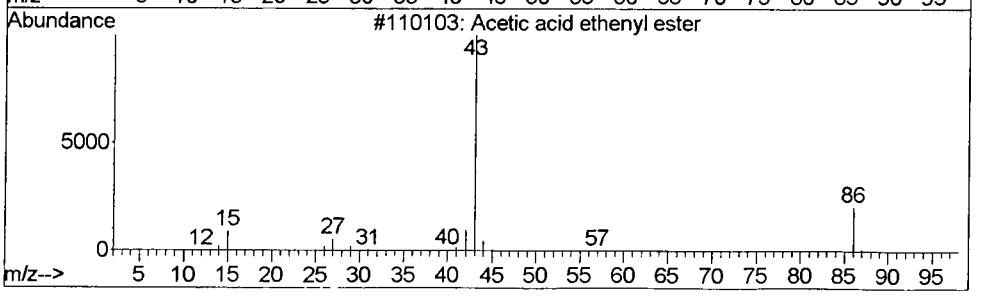
R.T.	EstConc	Area	Relative to ISTD	R.T.
3.77	0.21 ug/L	238984	1,4-Dichlorobenzene-d4	9.94

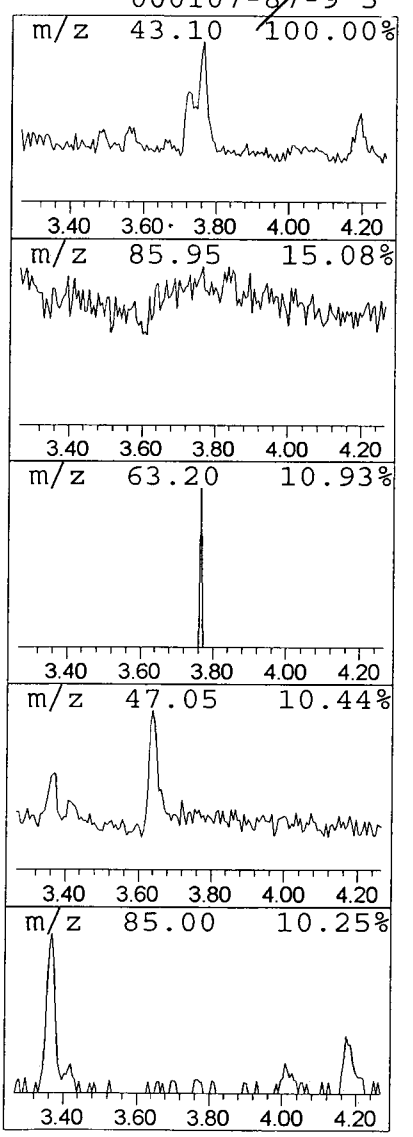
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanone, 1-oxiranyl-	86	C4H6O2	004401-11-0	9
2	Ethanone, 1-oxiranyl-	86	C4H6O2	004401-11-0	7
3	Acetic acid ethenyl ester	86	C4H6O2	000108-05-4	5
4	2-Pentanone	86	C5H10O	000107-87-9	5











Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050802\10203.D Vial: 21
Acq On : 9 May 2002 7:42 am Operator: Mclean
Sample : 4/30 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

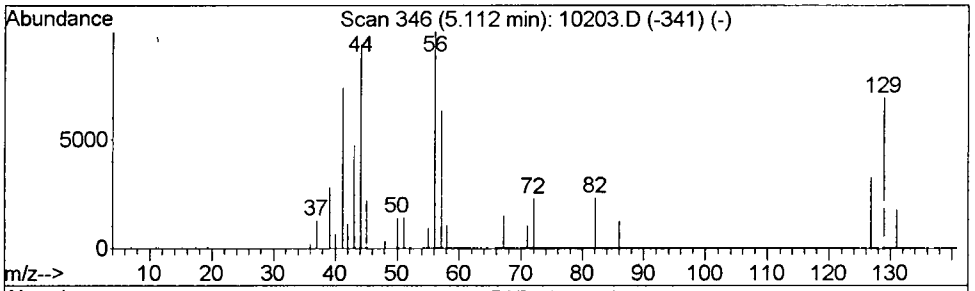
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

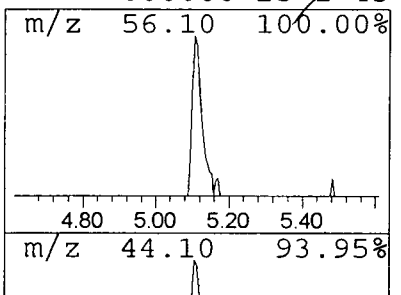
****

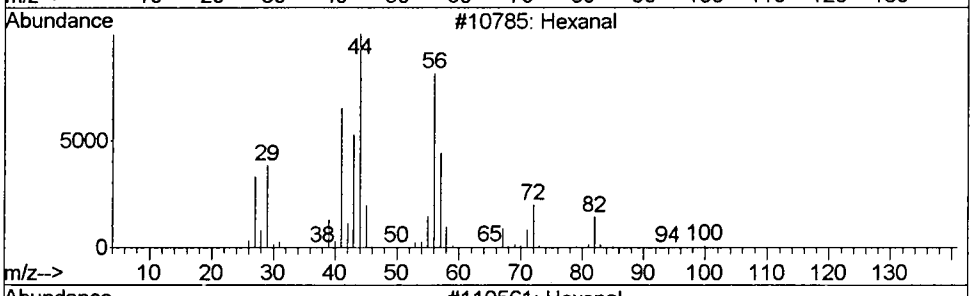
Peak Number 3 Hexanal Concentration Rank 6

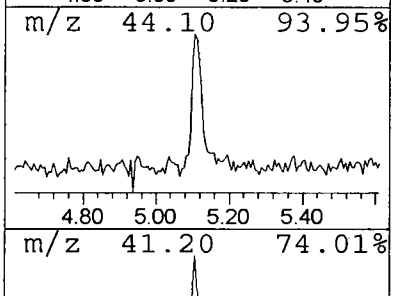
R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	0.21 ug/L	240949	1,4-Dichlorobenzene-d4	9.94

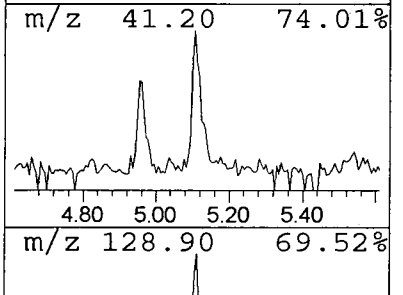
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanal	100	C6H12O	000066-25-1	46
2	Hexanal	100	C6H12O	000066-25-1	46
3	Hexanal	100	C6H12O	000066-25-1	46
4	Hexanal	100	C6H12O	000066-25-1	43

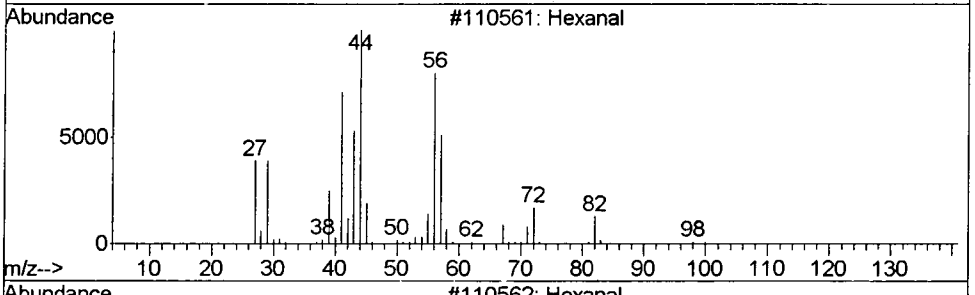


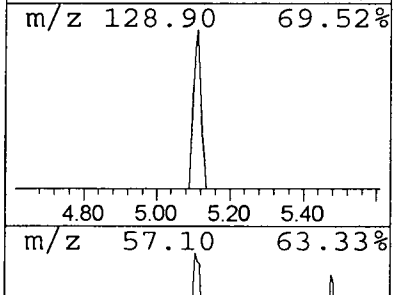


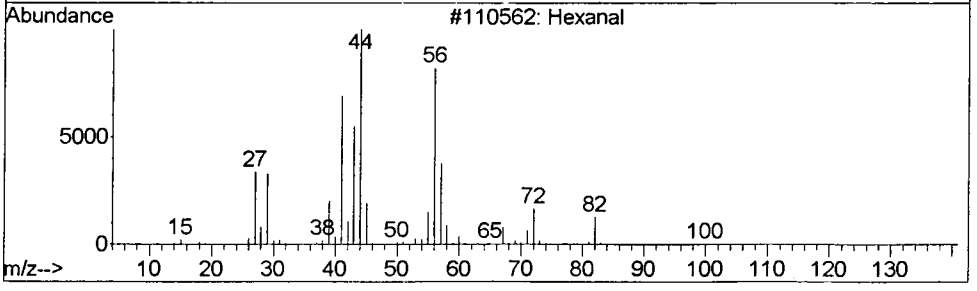


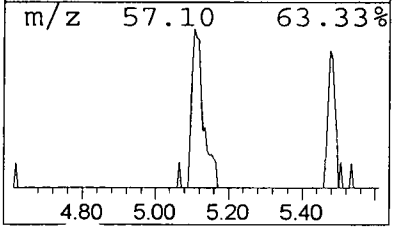












Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050802\10203.D Vial: 21
Acq On : 9 May 2002 7:42 am Operator: Mclean
Sample : 4/30 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

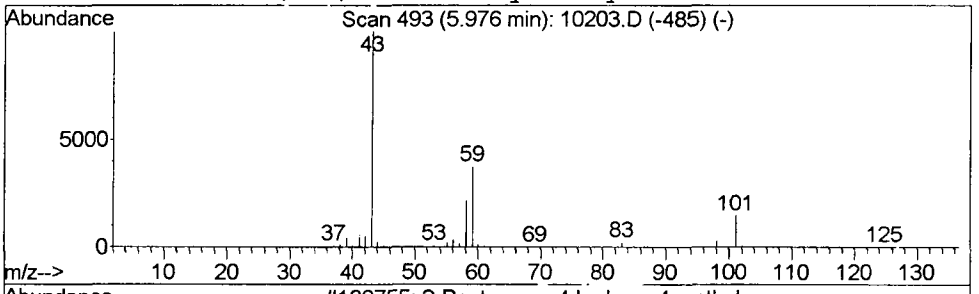
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

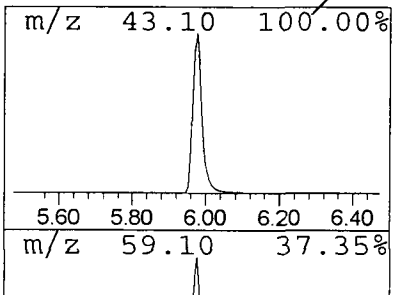
****

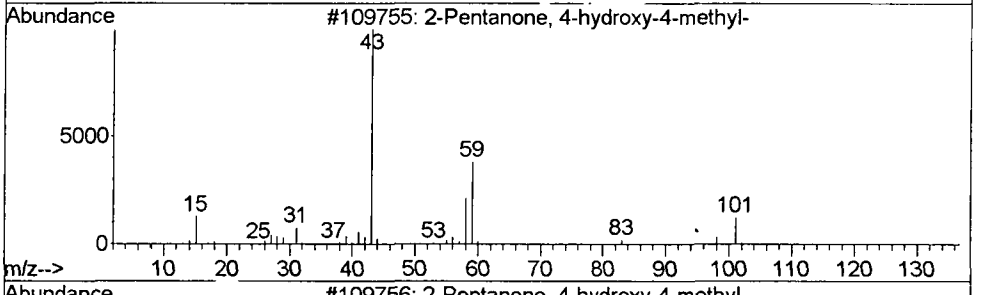
Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

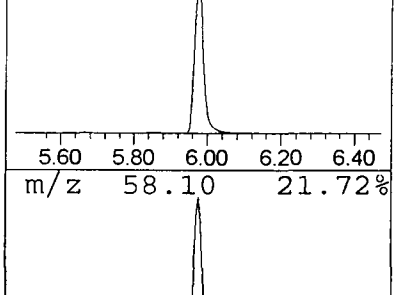
R.T.	EstConc	Area	Relative to ISTD	R.T.
5.98	13.02 ug/L	14923900	1,4-Dichlorobenzene-d4	9.94

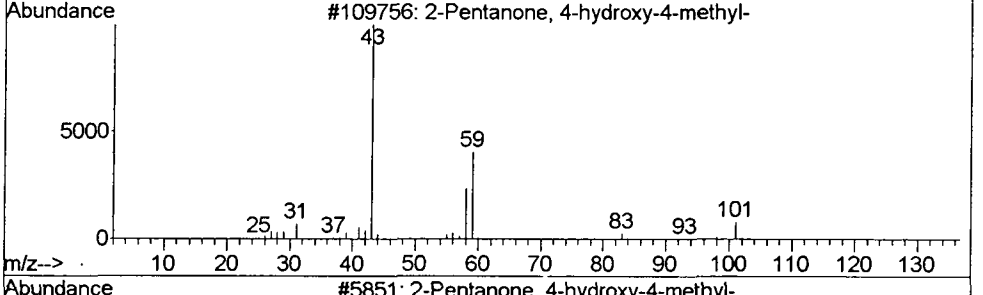
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	36

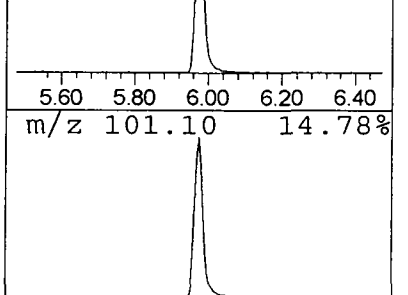


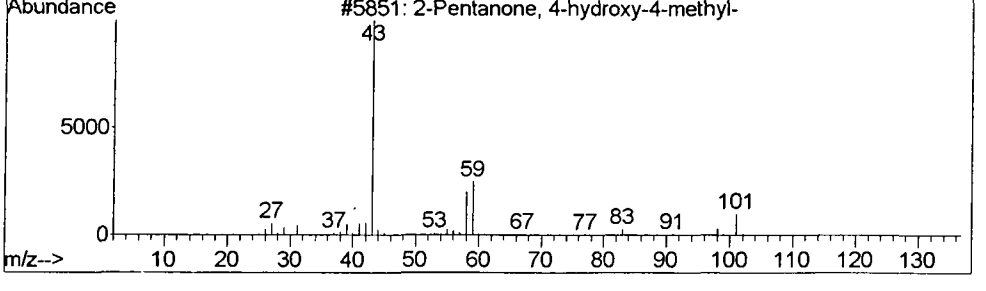


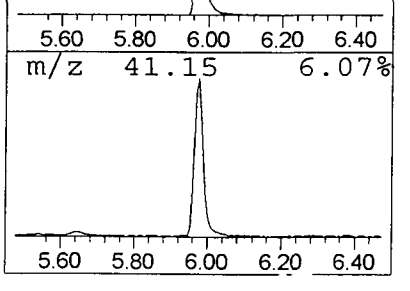












Library Search Compound Report

Data File : C:\MSDCHEM\1\DATA\050802\10203.D Vial: 21
Acq On : 9 May 2002 7:42 am Operator: Mclean
Sample : 4/30 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

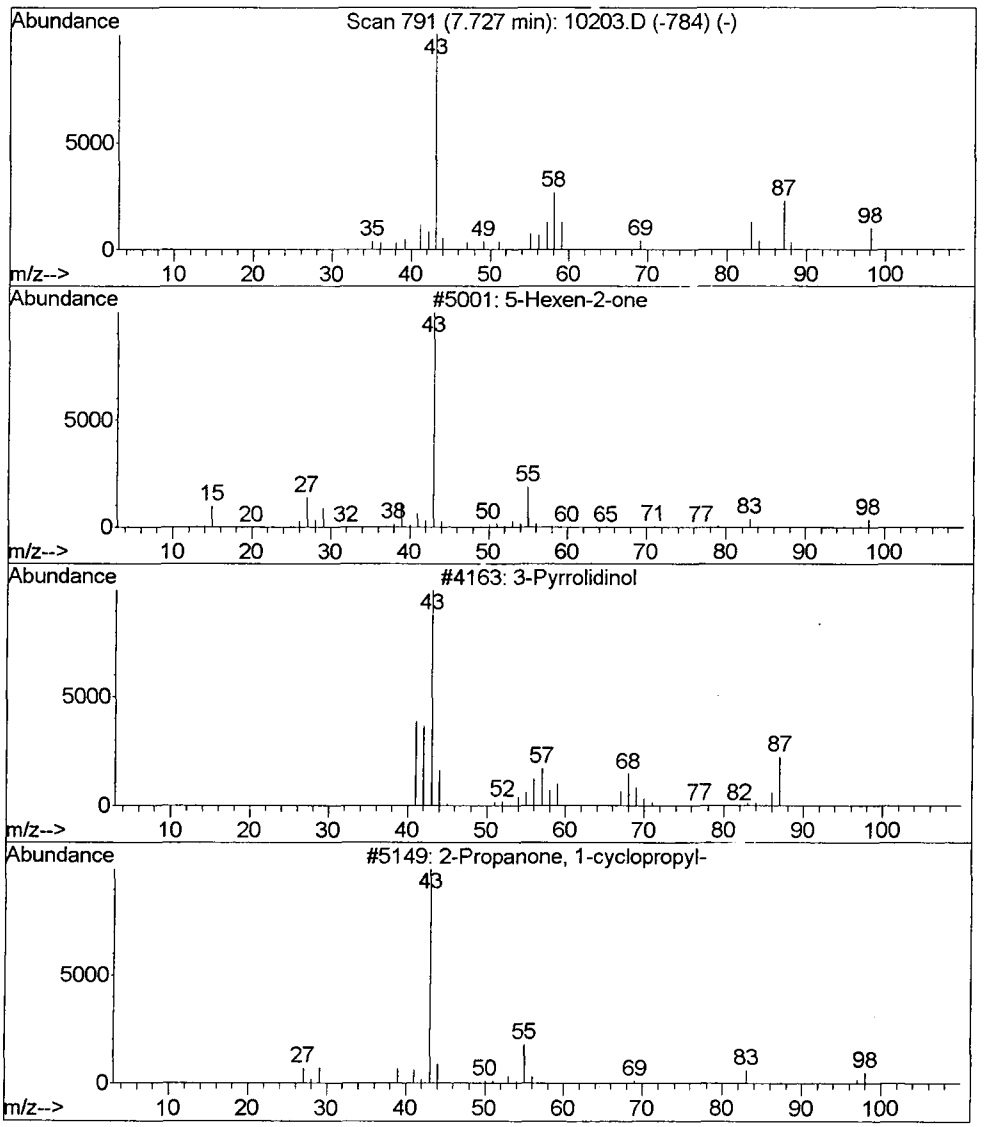
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

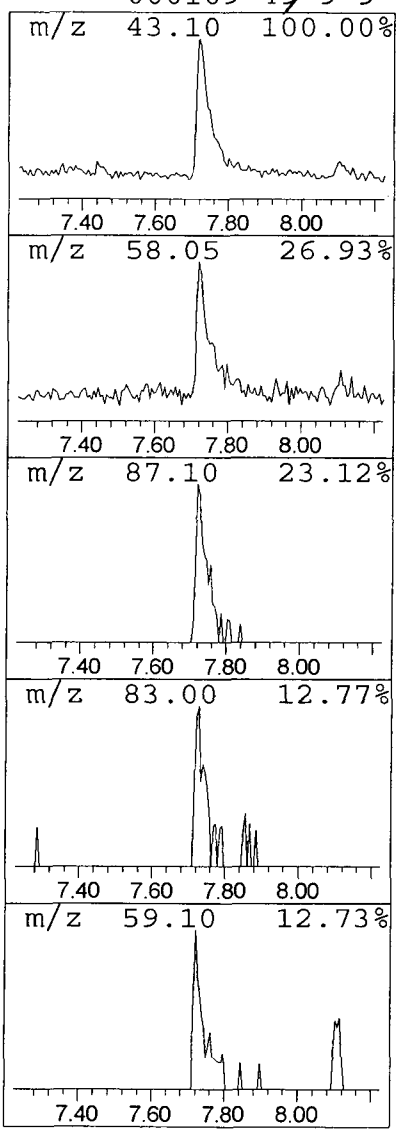
****

Peak Number 5 5-Hexen-2-one Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.73	0.35 ug/L	397846	1,4-Dichlorobenzene-d4	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Hexen-2-one	98	C6H10O	000109-49-9	10
2	3-Pyrrolidinol	87	C4H9NO	040499-83-0	9
3	2-Propanone, 1-cyclopropyl-	98	C6H10O	004160-75-2	9
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9





Library Search Compound Report

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Acq On : 9 May 2002 7:42 am Operator: Mclean
Sample : 4/30 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

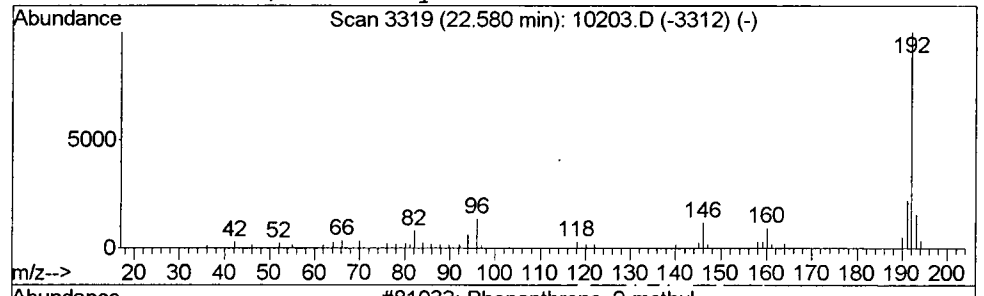
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

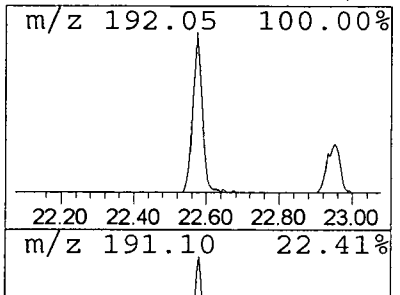
****

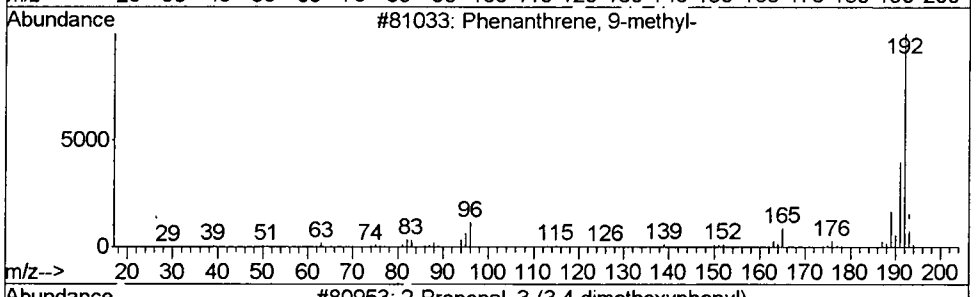
Peak Number 6 Phenanthrene, 9-methyl- Concentration Rank 4

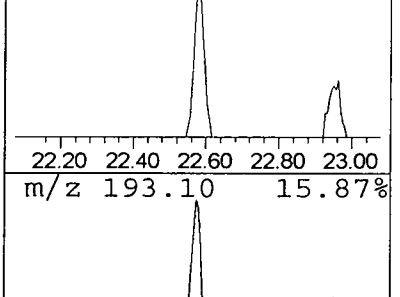
R.T.	EstConc	Area	Relative to ISTD	R.T.
22.58	0.36 ug/L	581119	Phenanthranene-d10	22.96

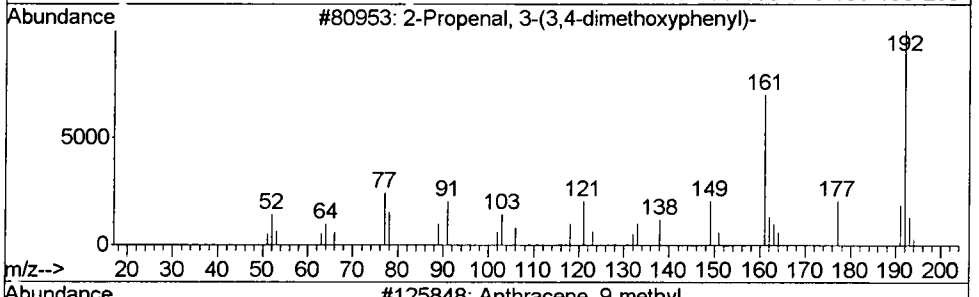
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 9-methyl-	192	C15H12	000883-20-5	53
2	2-Propenal, 3-(3,4-dimethoxyphen...	192	C11H12O3	004497-40-9	42
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	40
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	40

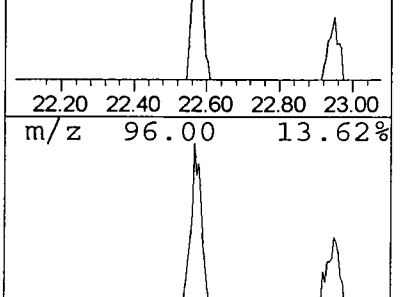


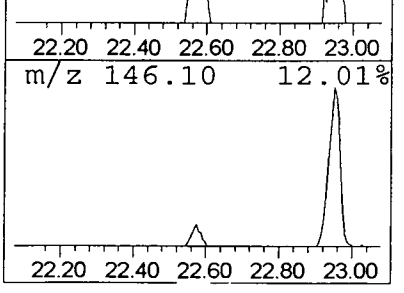


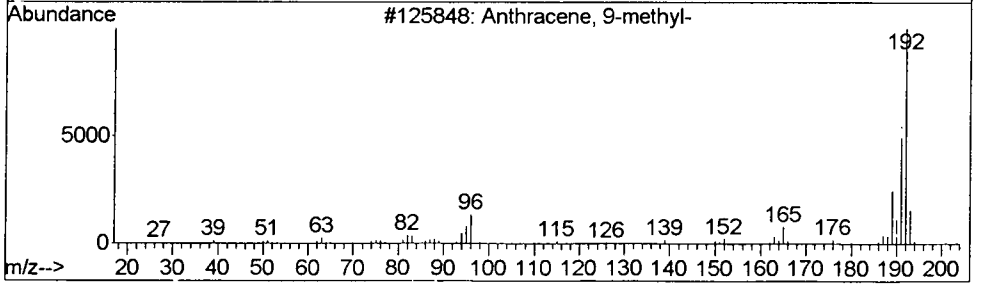












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Acq On : 9 May 2002 7:42 am Operator: Mclean
Sample : 4/30 eh Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

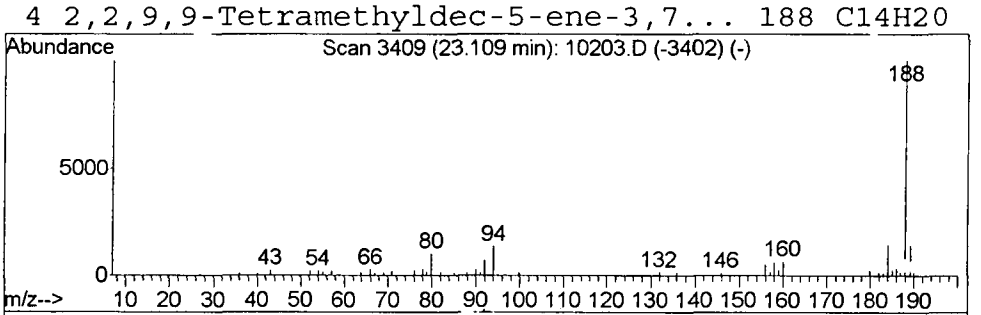
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

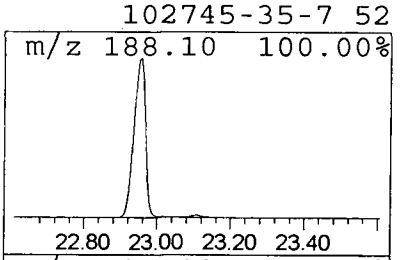
****

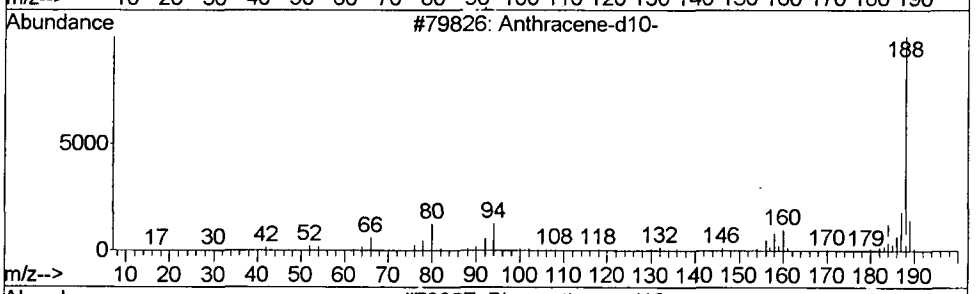
Peak Number 7 Anthracene-d10- Concentration Rank 2

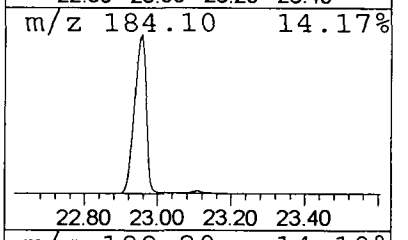
R.T.	EstConc	Area	Relative to ISTD	R.T.
23.11	0.66 ug/L	1071670	Phenanthranene-d10	22.96

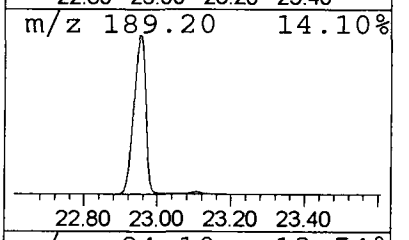
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene-d10-	188	C14D10	001719-06-8	83
2	Phenanthrene-d10	188	C14D10	001517-22-2	72
3	Selenide, ethyl 1-methyl-1-pente...	188	C8H12Se	025128-48-7	58
4	2,2,9,9-Tetramethyldec-5-ene-3,7...	188	C14H20	102745-35-7	52

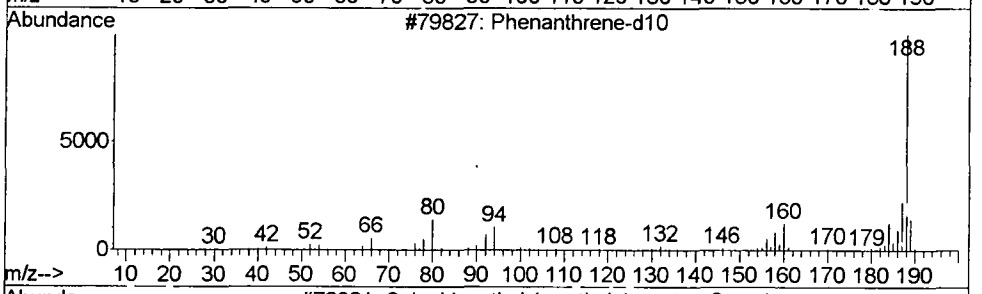


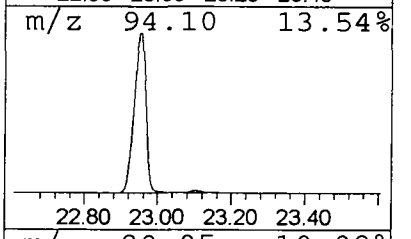


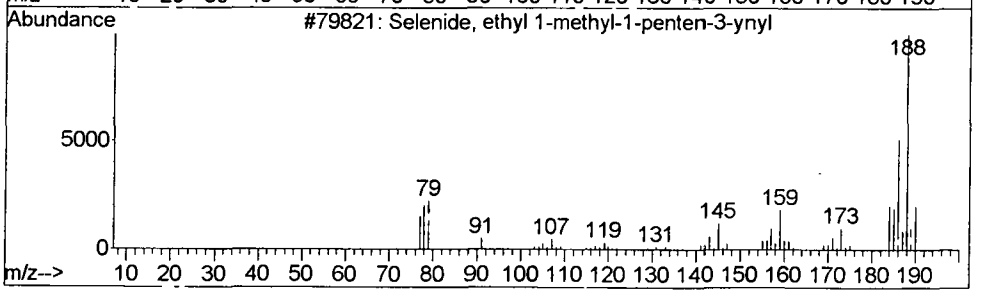


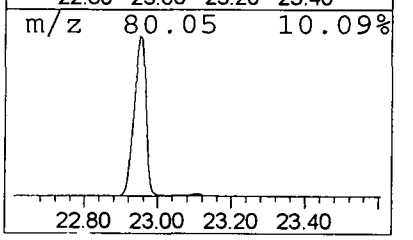












Tentatively Identified Compound (LSC) summary

Operator ID: Mclean Date Acquired: 9 May 2002 7:42 am
Data File: C:\MSDCHEM\1\DATA\050802\10203.D
Name: 4/30 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
Acetonitrile, dic...	3.64	0.4	ug/L	476773	1	9.94	45837800	40.0
Ethanone, 1-oxira...	3.77	0.2	ug/L	238984	1	9.94	45837800	40.0
Hexanal	5.11	0.2	ug/L	240949	1	9.94	45837800	40.0
2-Pentanone, 4-hy...	5.98	13.0	ug/L	14923900	1	9.94	45837800	40.0
5-Hexen-2-one	7.73	0.3	ug/L	397846	1	9.94	45837800	40.0
Phenanthrene, 9-m...	22.58	0.4	ug/L	581119	4	22.96	65404600	40.0
Anthracene-d10-	23.11	0.7	ug/L	1071670	4	22.96	65404600	40.0