User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 05/20/2002 Time: 14:46:10

Sample: AE10815
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 5/20/02 14:34 Ended: 5/20/02 14:34 Analyst: DLV
Page: 1

	Component Name	Vial	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		99		10.0

Comments for Sample AE10815 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 05-20-2002 Time: 14:46:34

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds.

Data File : C:\MSDCHEM\1\DATA\051402\10815.D Vial: 25
Acq On : 15 May 2002 9:55 am Operator: Mclean
Sample : 5/7 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 16 14:16:51 2002 Quant Results File: 050102.RES

Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Thu May 16 12:34:41 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	6880723	40.00	ug/L	0.00
10) Napthalene-d8	13.43	136	27369097	40.00	ug/L	0.00
17) Acenapthalene-d10	18.57	164	14790774	40.00	ug/L	0.00
29) Phenanthranene-d10	22.95	188	21438990	40.00	ug/L	0.00
37) Chrysene-d12	30.81	240	15281340	40.00	ug/L	0.00
43) Perylene-d12	34.71	264	9761242	40.00	ug/L	0.00
System Monitoring Compounds						
27) TCMX	20.54	209	891664	9.93	ug/L	0.00
Spiked Amount 10.000			Recovery = 99.30%			
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	12.94	93	3097	N.D.		
14) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.		
15) Napthalene	13.48	128	36564	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	20.14	149	14318	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	17359	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	25.09	149	38402	N.D.		

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

Data File : C:\MSDCHEM\1\DATA\051402\10815.D Vial: 25
Acq On : 15 May 2002 9:55 am Operator: Mclean
Sample : 5/7 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 16 14:16:51 2002 Quant Results File: 050102.RES

Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Thu May 16 12:34:41 2002
Response via : Initial Calibration
DataAcq Meth : 508SCAN

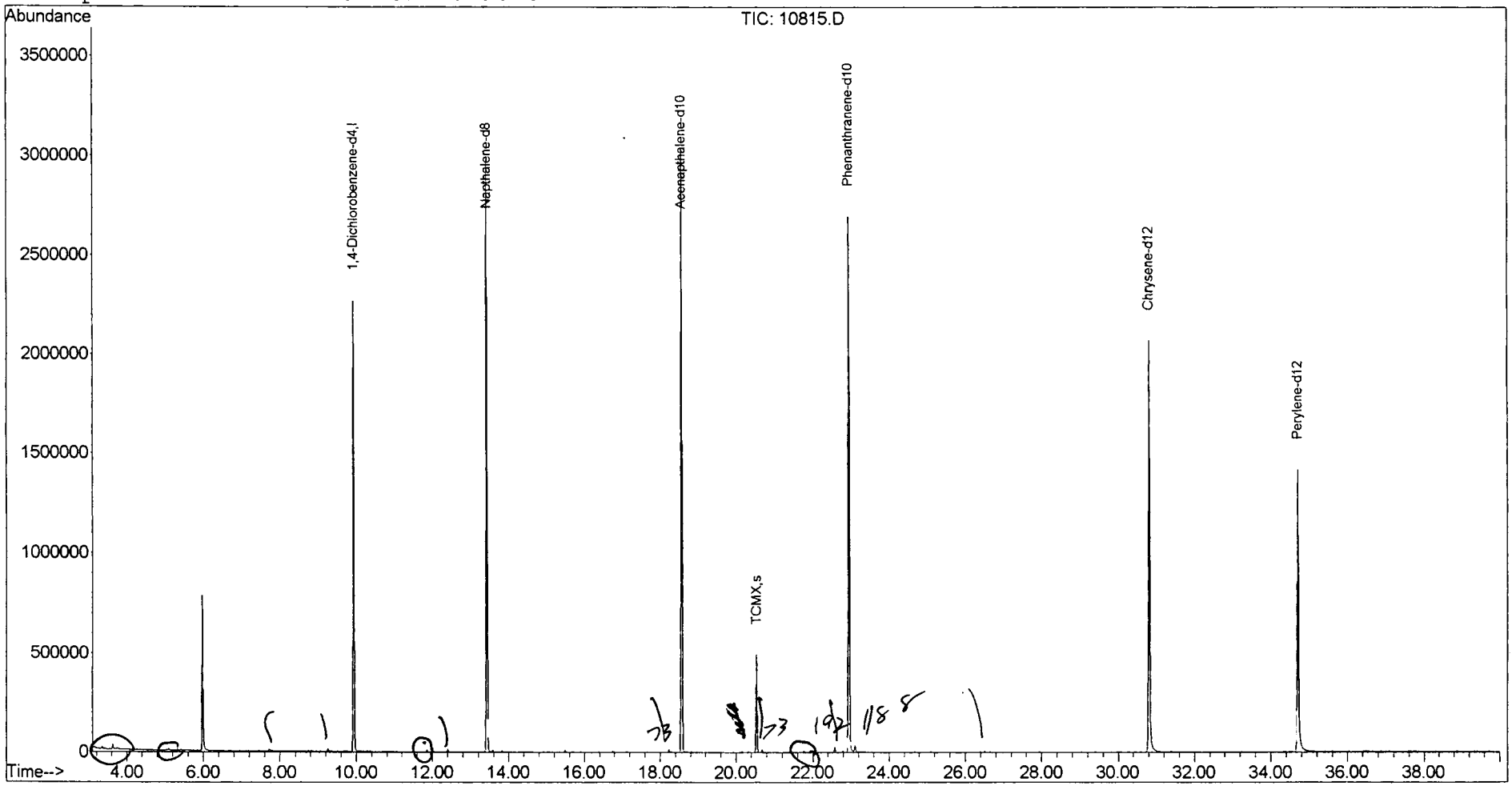
	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
36)	Fluoroanthene	0.00	202	0	N.D.	
38)	Pyrene	0.00	202	0	N.D.	
39)	Butylbenzylphthalate	0.00	149	0	N.D.	
40)	Benzo(a)anthracene	30.81	228	40042	N.D.	
41)	Bis(2-ethylhexyl)phthalate	0.00	149	0	N.D.	
42)	Chrysene	0.00	228	0	N.D.	
44)	Di-n-octylphthalate	0.00	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\051402\10815.D Vial: 25
Acq On : 15 May 2002 9:55 am Operator: Mclean
Sample : 5/7 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: EVENTS.E
Quant Time: May 16 14:16 2002 Quant Results File: 050102.RES

Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Last Update : Thu May 16 12:34:41 2002
Response via : Initial Calibration



Data File : C:\MSDCHEM\1\DATA\051402\10815.D Vial: 25
Acq On : 15 May 2002 9:55 am Operator: Mclean
Sample : 5/7 kb 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

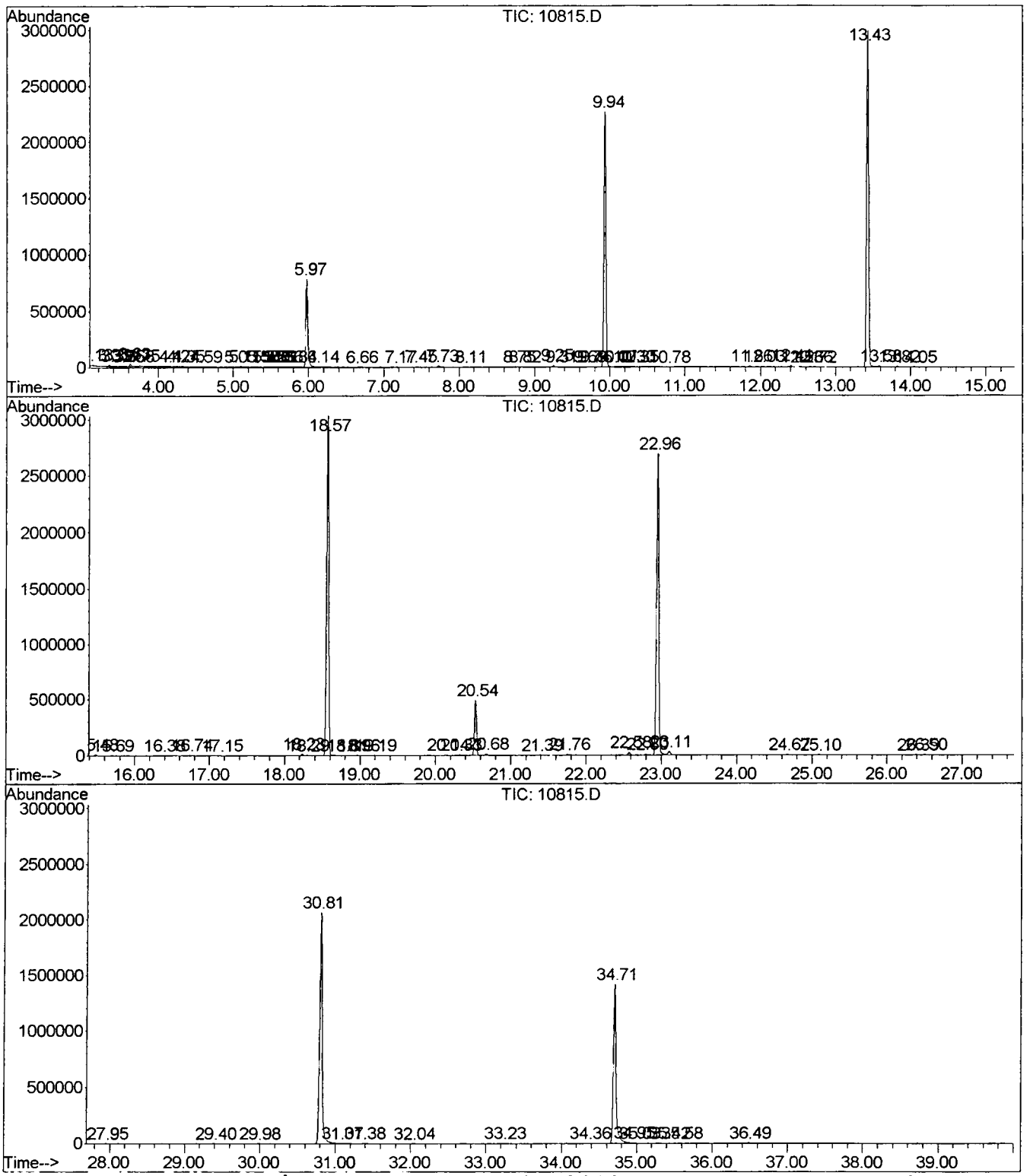
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.132	3	9	20	BV 3	6151	123546	0.20%	0.037%
2	3.349	40	46	51	PV	11347	193326	0.32%	0.058%
3	3.396	51	54	66	VV 7	4159	106306	0.17%	0.032%
4	3.543	75	79	88	PV 6	3414	84188	0.14%	0.025%
5	3.626	88	93	101	VV	23635	401875	0.66%	0.121%
6	3.684	101	103	105	VV 3	3557	49914	0.08%	0.015%
7	3.755	105	115	119	VV 2	9081	260091	0.43%	0.078%
8	4.172	183	186	196	VV 7	3558	105043	0.17%	0.032%
9	4.272	196	203	207	VV 3	3457	76283	0.13%	0.023%
10	4.354	212	217	223	VV 3	5192	91880	0.15%	0.028%
11	4.589	252	257	266	PV 10	2369	47089	0.08%	0.014%
12	5.047	330	335	340	VV 6	3200	63683	0.10%	0.019%
13	5.106	340	345	356	VV 10	7817	174620	0.29%	0.052%
14	5.347	381	386	395	VV 6	2014	58742	0.10%	0.018%
15	5.423	395	399	403	VV 3	4998	82963	0.14%	0.025%
16	5.529	414	417	418	PV 2	2205	16996	0.03%	0.005%
17	5.553	418	421	423	VV 4	2510	42825	0.07%	0.013%
18	5.576	423	425	428	VV 4	2807	37773	0.06%	0.011%
19	5.612	428	431	433	VV 3	3482	43202	0.07%	0.013%
20	5.664	433	440	448	VV 7	4398	111098	0.18%	0.033%
21	5.847	467	471	480	VV 8	2301	50924	0.08%	0.015%
22	5.976	484	493	520	VV	785255	14066516	23.10%	4.221%
23	6.140	520	521	525	VV 4	2649	33802	0.06%	0.010%
24	6.663	605	610	615	PV 4	1976	42506	0.07%	0.013%
25	7.174	692	697	702	VV 5	2060	39677	0.07%	0.012%
26	7.445	739	743	753	VV 4	3974	78207	0.13%	0.023%
27	7.727	786	791	810	VV	10893	319813	0.53%	0.096%
28	8.109	852	856	861	VV 2	3151	58816	0.10%	0.018%
29	8.749	958	965	971	PV 4	3843	77746	0.13%	0.023%
30	8.820	971	977	987	VV 7	3889	100044	0.16%	0.030%
31	9.254	1045	1051	1056	VV 2	14511	241472	0.40%	0.072%
32	9.307	1056	1060	1065	VV 4	2817	62068	0.10%	0.019%
33	9.689	1122	1125	1130	VV 3	2557	52210	0.09%	0.016%
34	9.736	1130	1133	1135	VV 3	2349	30170	0.05%	0.009%
35	9.766	1135	1138	1141	VV 3	3148	45869	0.08%	0.014%
36	9.936	1156	1167	1186	VV	2235873	41898697	68.80%	12.574%
37	10.065	1186	1189	1197	VV 7	2349	51622	0.08%	0.015%

39	10.353	1235	1238	1253	VV	10	3396	94417	0.16%	0.028%
40	10.782	1303	1311	1322	VV	6	1640	53901	0.09%	0.016%
41	11.863	1488	1495	1501	BV	4	9145	142528	0.23%	0.043%
42	12.034	1520	1524	1531	VV	3	3248	67450	0.11%	0.020%
43	12.404	1580	1587	1593	VV		13441	223990	0.37%	0.067%
44	12.527	1602	1608	1619	VV	5	1848	41373	0.07%	0.012%
45	12.662	1625	1631	1636	VV	3	2199	36588	0.06%	0.011%
46	12.721	1636	1641	1648	VV		2013	39656	0.07%	0.012%
47	13.432	1748	1762	1782	PV		2963846	56241101	92.35%	16.878%
48	13.585	1782	1788	1793	VV	2	9844	197408	0.32%	0.059%
49	13.826	1819	1829	1844	VV	5	6700	146911	0.24%	0.044%
50	14.043	1852	1866	1873	VV	5	2796	62358	0.10%	0.019%
51	15.482	2100	2111	2117	PV	2	10816	175053	0.29%	0.053%
52	15.694	2143	2147	2156	VV	5	2857	56922	0.09%	0.017%
53	16.376	2244	2263	2270	BV	4	1983	42690	0.07%	0.013%
54	16.740	2320	2325	2339	VV	4	5286	157958	0.26%	0.047%
55	17.151	2384	2395	2401	VV	5	2005	37464	0.06%	0.011%
56	18.232	2573	2579	2585	PV		15856	255485	0.42%	0.077%
57	18.291	2585	2589	2598	VV	2	3068	54146	0.09%	0.016%
58	18.573	2626	2637	2660	VV		3012650	60898547	100.00%	18.276%
59	18.814	2668	2678	2697	VV	5	2069	72105	0.12%	0.022%
60	18.961	2697	2703	2714	VV	2	2250	46271	0.08%	0.014%
61	19.190	2737	2742	2746	PV	4	1871	30380	0.05%	0.009%
62	20.142	2894	2904	2913	PV	3	4230	89056	0.15%	0.027%
63	20.324	2930	2935	2941	PV	2	4386	70517	0.12%	0.021%
64	20.535	2961	2971	2982	PV		482437	8853840	14.54%	2.657%
65	20.682	2982	2996	3002	VV		14660	239242	0.39%	0.072%
66	21.393	3111	3117	3125	BV	4	2407	43777	0.07%	0.013%
67	21.758	3170	3179	3184	BV	3	7364	107467	0.18%	0.032%
68	22.580	3312	3319	3333	VV	2	26481	477806	0.78%	0.143%
69	22.798	3349	3356	3362	PV	2	8658	135010	0.22%	0.041%
70	22.956	3369	3383	3401	BV	2	2707772	58609478	96.24%	17.589%
71	23.109	3401	3409	3421	VV	2	32430	734987	1.21%	0.221%
72	24.672	3670	3675	3683	VV	2	4347	79338	0.13%	0.024%
73	25.095	3738	3747	3756	PV	3	3409	65741	0.11%	0.020%
74	26.393	3960	3968	3973	PV	3	3248	57817	0.09%	0.017%
75	26.505	3980	3987	3997	VV	4	3796	77625	0.13%	0.023%
76	27.950	4222	4233	4244	VV	4	2703	49392	0.08%	0.015%
77	29.396	4473	4479	4490	VV	4	2412	47529	0.08%	0.014%
78	29.983	4569	4579	4592	BV	4	2098	54120	0.09%	0.016%
79	30.806	4703	4719	4762	PV	2	2040195	48004677	78.83%	14.406%
80	31.070	4762	4764	4781	VV	6	3316	148767	0.24%	0.045%
81	31.376	4810	4816	4829	PV	7	3762	104231	0.17%	0.031%
82	32.034	4924	4928	4940	VV	2	2401	50696	0.08%	0.015%
83	33.233	5120	5132	5141	VV	5	2522	55291	0.09%	0.017%
84	34.361	5313	5324	5329	PV	5	2775	70913	0.12%	0.021%
85	34.707	5369	5383	5423	VV	2	1404197	35736859	58.68%	10.725%
86	34.948	5423	5424	5436	VV	8	6066	218109	0.36%	0.065%
87	35.025	5436	5437	5447	VV	6	3648	106291	0.17%	0.032%

90	36.488	5675	5686	5693	VV	6	4392	115051	0.19%	0.035%

Sum of corrected areas: 333223945

File : C:\MSDCHEM\1\DATA\051402\10815.D
Operator : Mclean
Acquired : 15 May 2002 9:55 am using AcqMethod 508SCAN
Instrument : Instrumen
Sample Name: 5/7 kb
Misc Info :
Vial Number: 25
Quant File :050102.RES (Chemstation Integrator)



Data File : C:\MSDCHEM\1\DATA\051402\10815.D Vial: 25
Acq On : 15 May 2002 9:55 am Operator: Mclean
Sample : 5/7 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

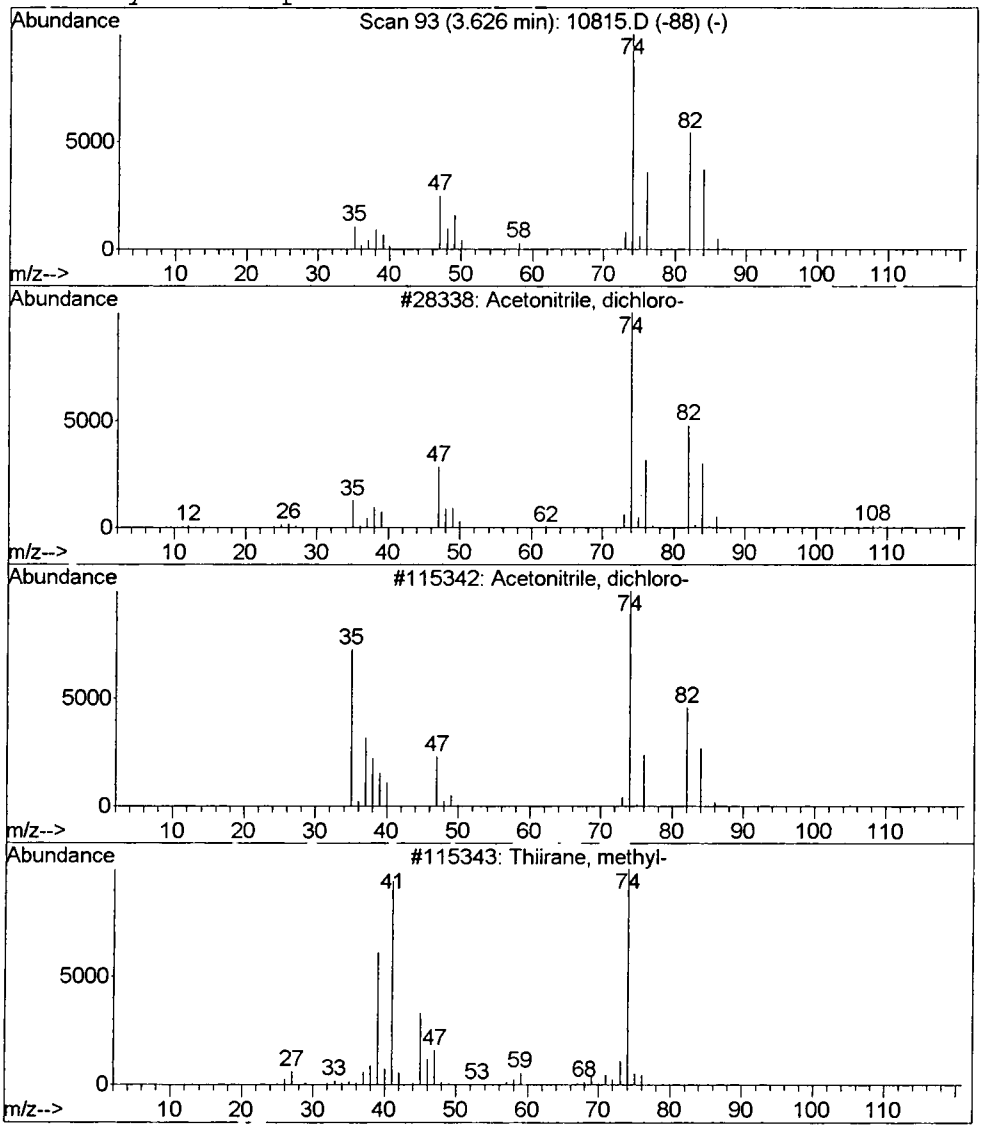
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

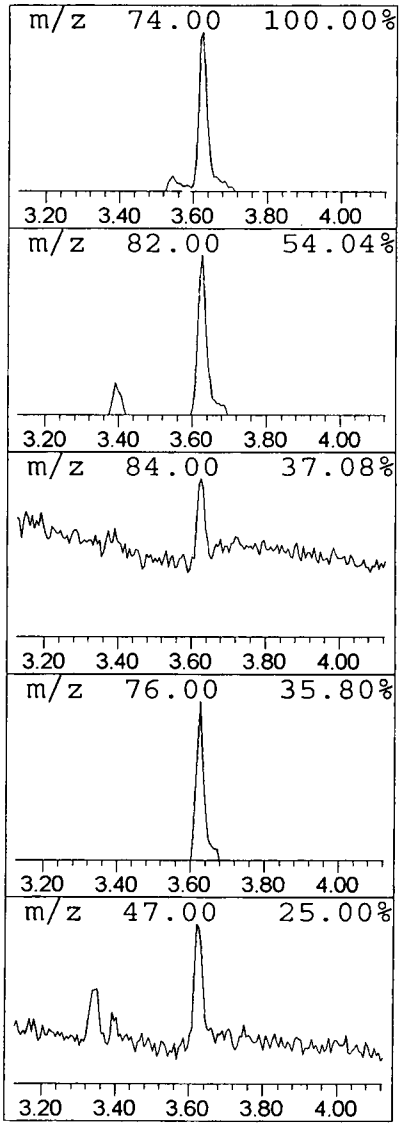
****

Peak Number 1 Acetonitrile, dichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.63	0.38 ug/L	401875	1,4-Dichlorobenzene-d4	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetonitrile, dichloro-	109	C2HCl2N	003018-12-0	91
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





Data File : C:\MSDCHEM\1\DATA\051402\10815.D Vial: 25
Acq On : 15 May 2002 9:55 am Operator: Mclean
Sample : 5/7 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

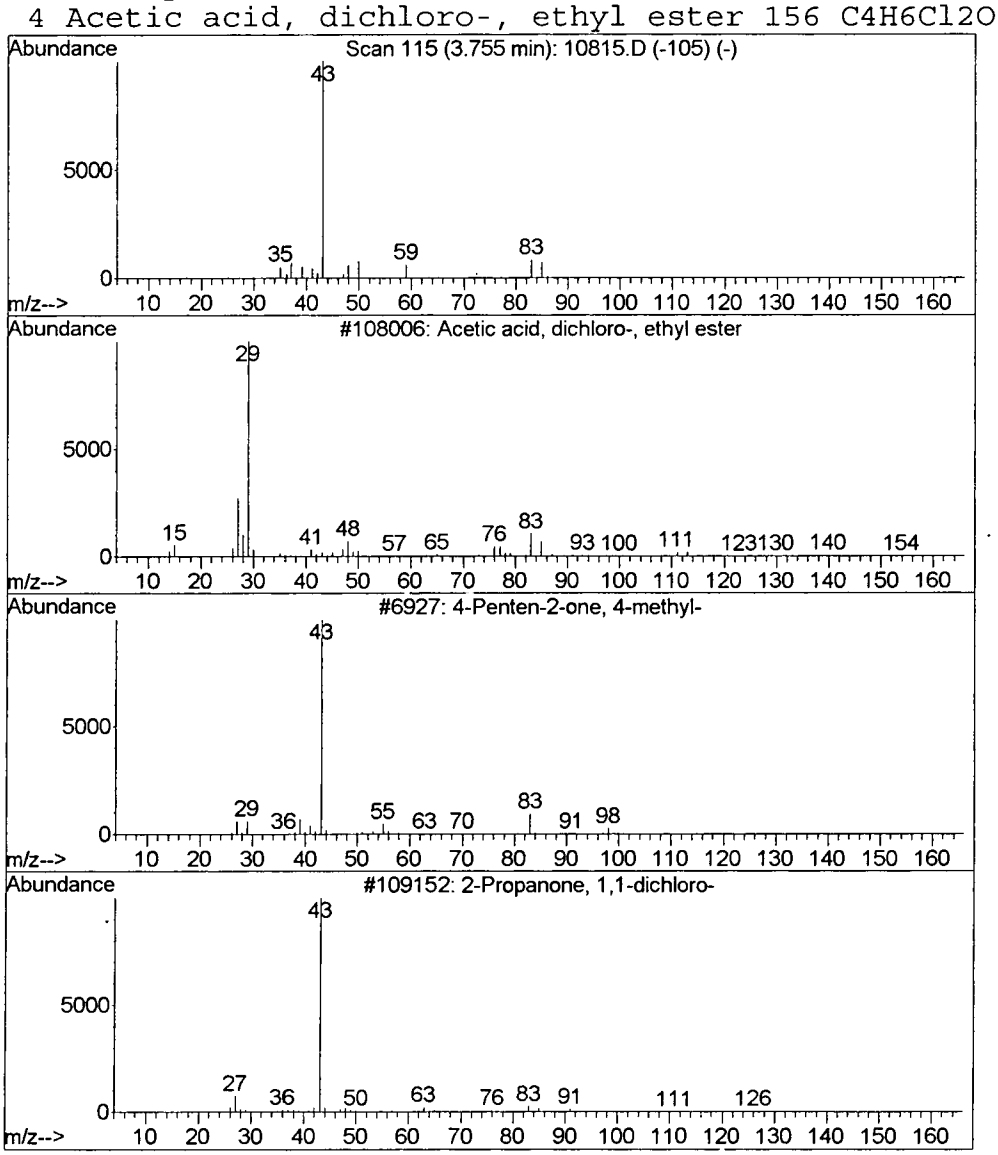
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

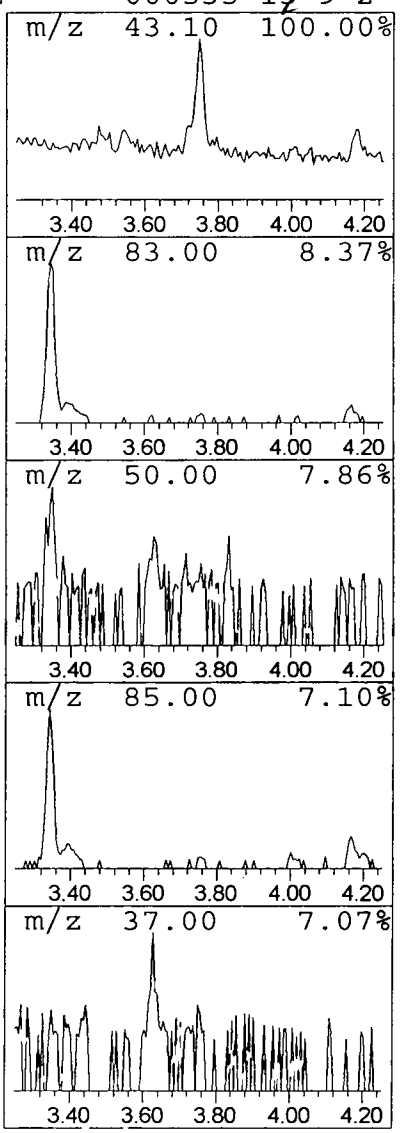
****

Peak Number 2 Acetic acid, dichloro-, eth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.75	0.25 ug/L	260091	1,4-Dichlorobenzene-d4	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, dichloro-, ethyl ester	156	C4H6Cl2O2	000535-15-9	4
2	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	2
3	2-Propanone, 1,1-dichloro-	126	C3H4Cl2O	000513-88-2	2
4	Acetic acid, dichloro-, ethyl ester	156	C4H6Cl2O2	000535-15-9	2





Data File : C:\MSDCHEM\1\DATA\051402\10815.D Vial: 25
Acq On : 15 May 2002 9:55 am Operator: Mclean
Sample : 5/7 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

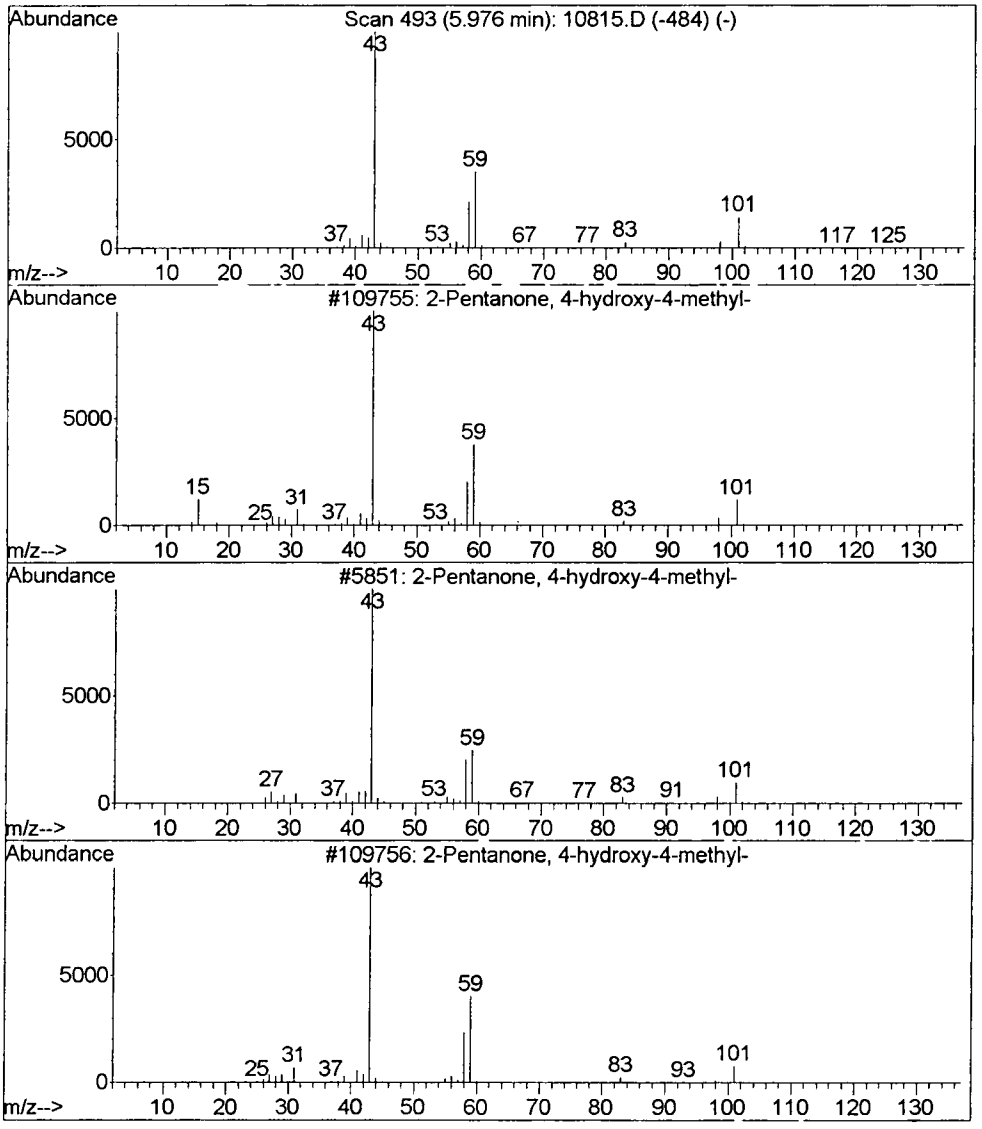
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

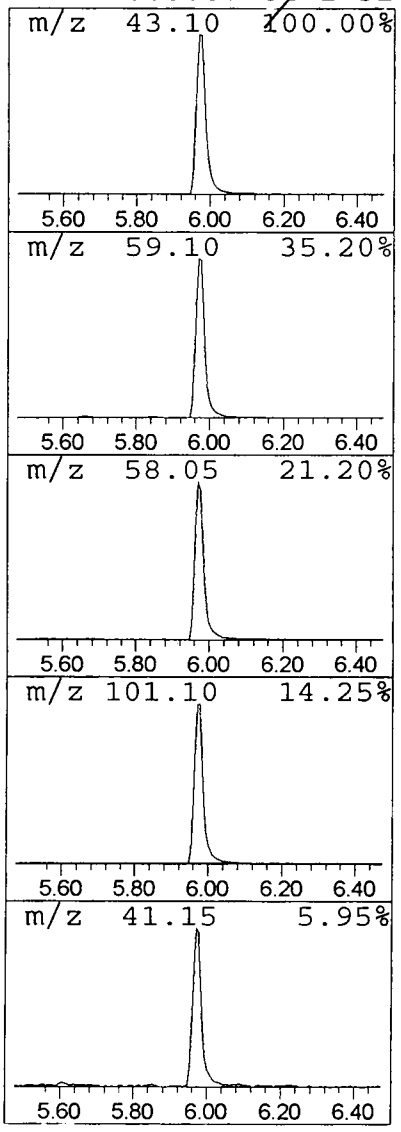
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.97	13.43 ug/L	14066500	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	78
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	64
4	Acetone	58	C3H6O	000067-64-1	32





Data File : C:\MSDCHEM\1\DATA\051402\10815.D Vial: 25
Acq On : 15 May 2002 9:55 am Operator: Mclean
Sample : 5/7 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

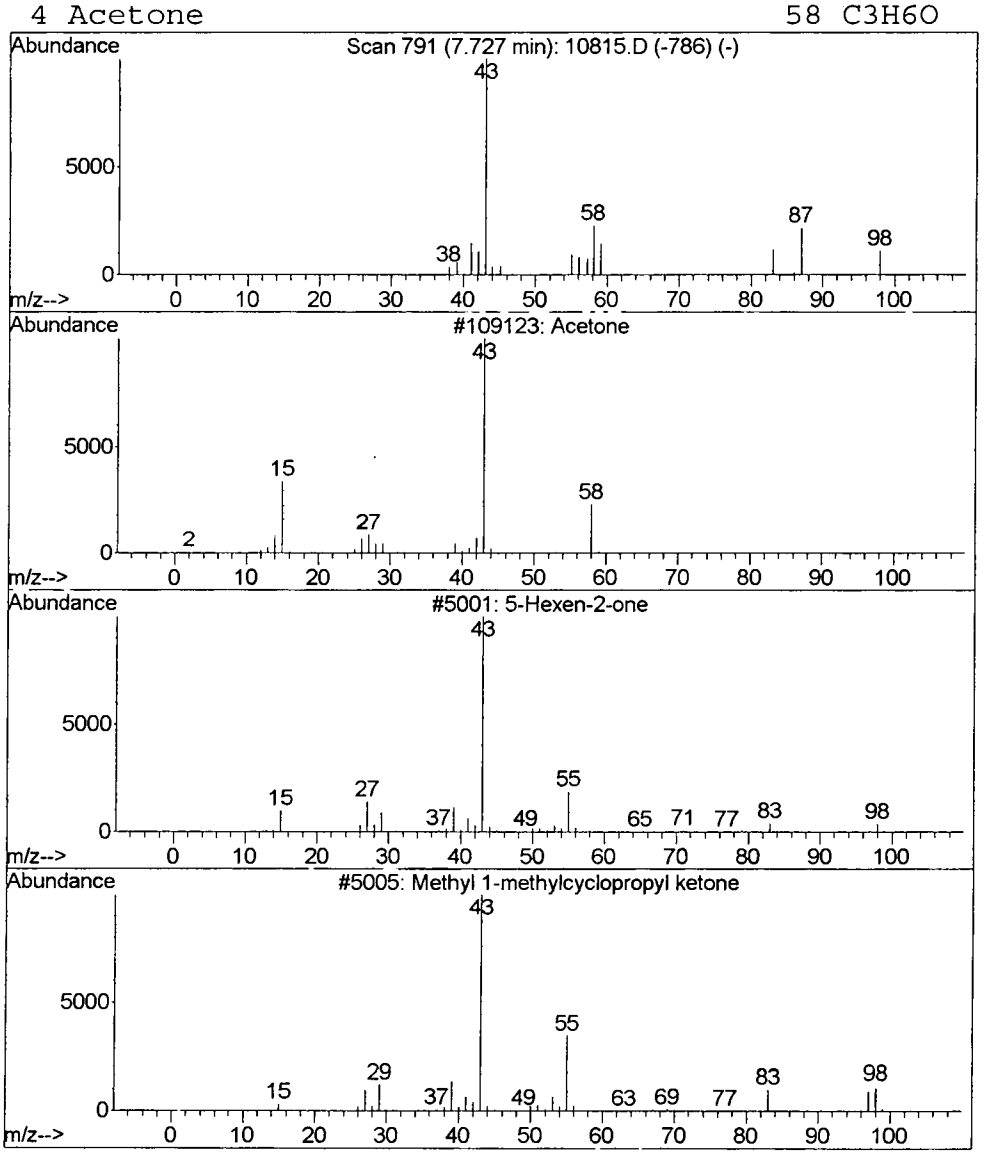
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

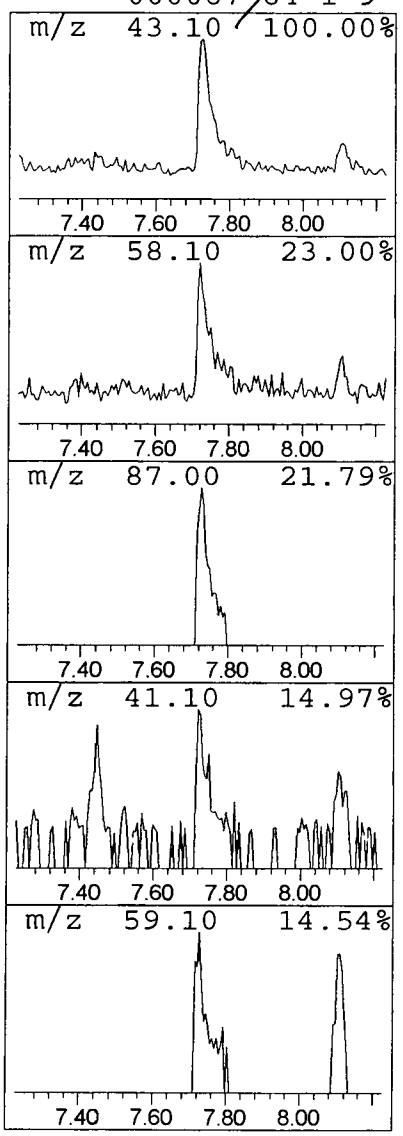
****

Peak Number 4 Acetone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.73	0.31 ug/L	319813	1,4-Dichlorobenzene-d4	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetone	58	C3H6O	000067-64-1	9
2	5-Hexen-2-one	98	C6H10O	000109-49-9	9
3	Methyl 1-methylcyclopropyl ketone	98	C6H10O	001567-75-5	9
4	Acetone	58	C3H6O	000067-64-1	9





Data File : C:\MSDCHEM\1\DATA\051402\10815.D Vial: 25
Acq On : 15 May 2002 9:55 am Operator: Mclean
Sample : 5/7 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

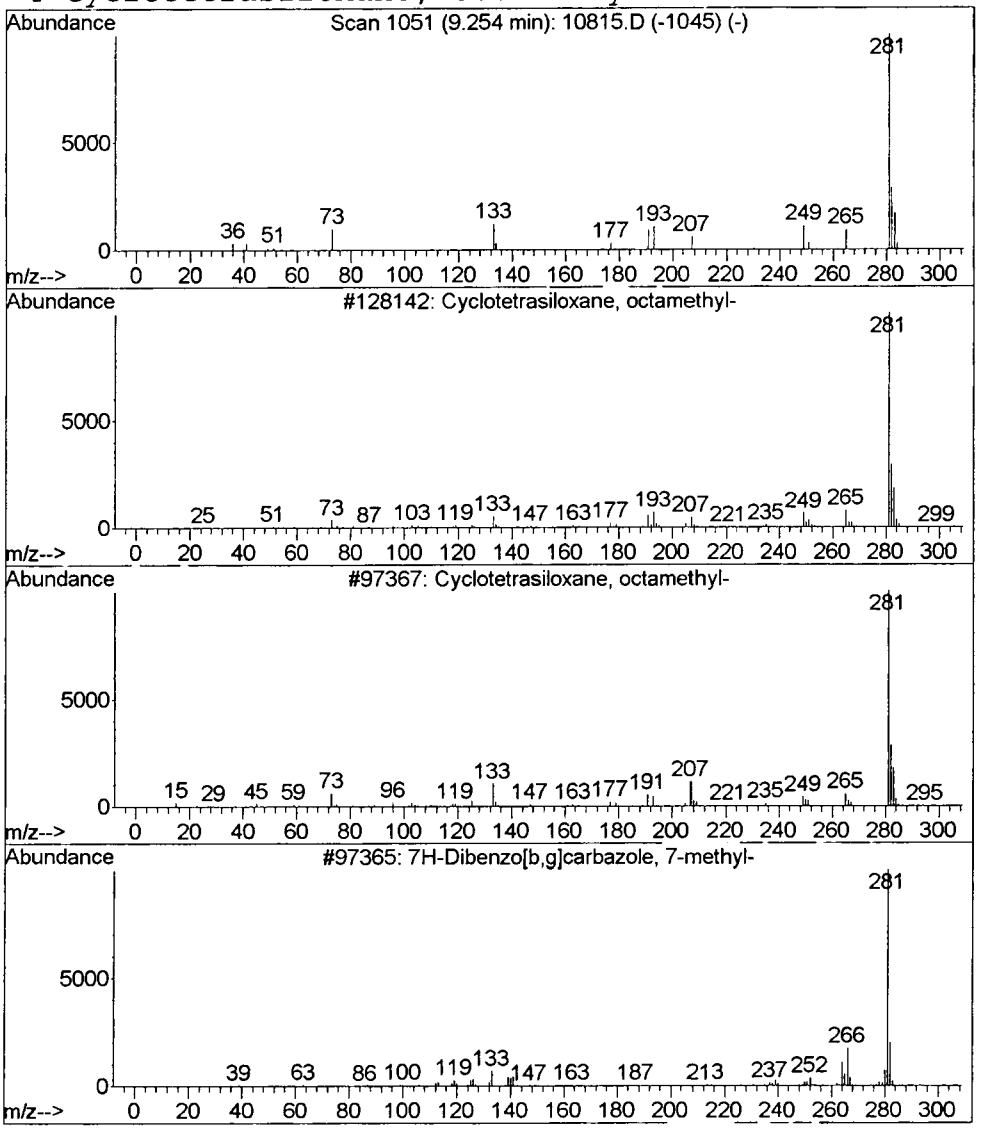
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

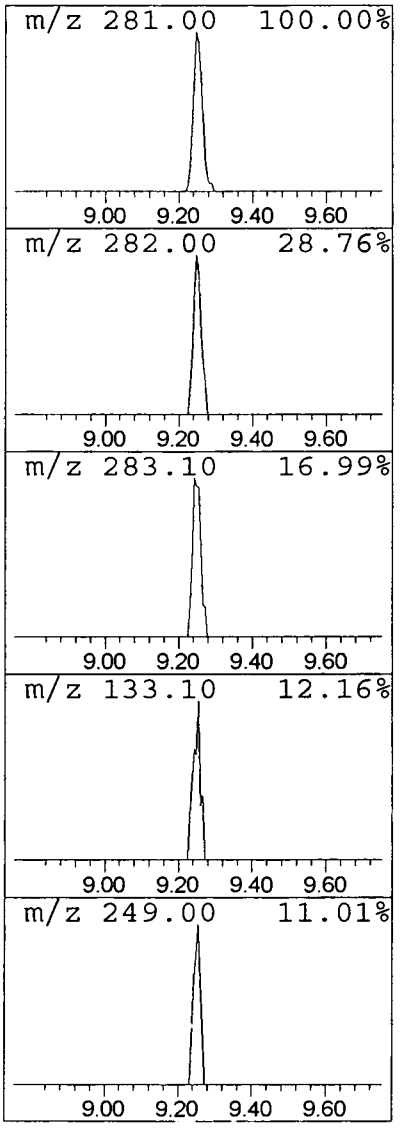
****

Peak Number 5 Cyclotetrasiloxane, octamet... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.25	0.23 ug/L	241472	1,4-Dichlorobenzene-d4	9.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	83
2	Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	64
3	7H-Dibenzo[b,g]carbazole, 7-methyl-	281	C21H15N	003557-49-1	53
4	Cyclotetrasiloxane, octamethyl-	296	C8H24O4Si4	000556-67-2	50





Data File : C:\MSDCHEM\1\DATA\051402\10815.D Vial: 25
Acq On : 15 May 2002 9:55 am Operator: Mclean
Sample : 5/7 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

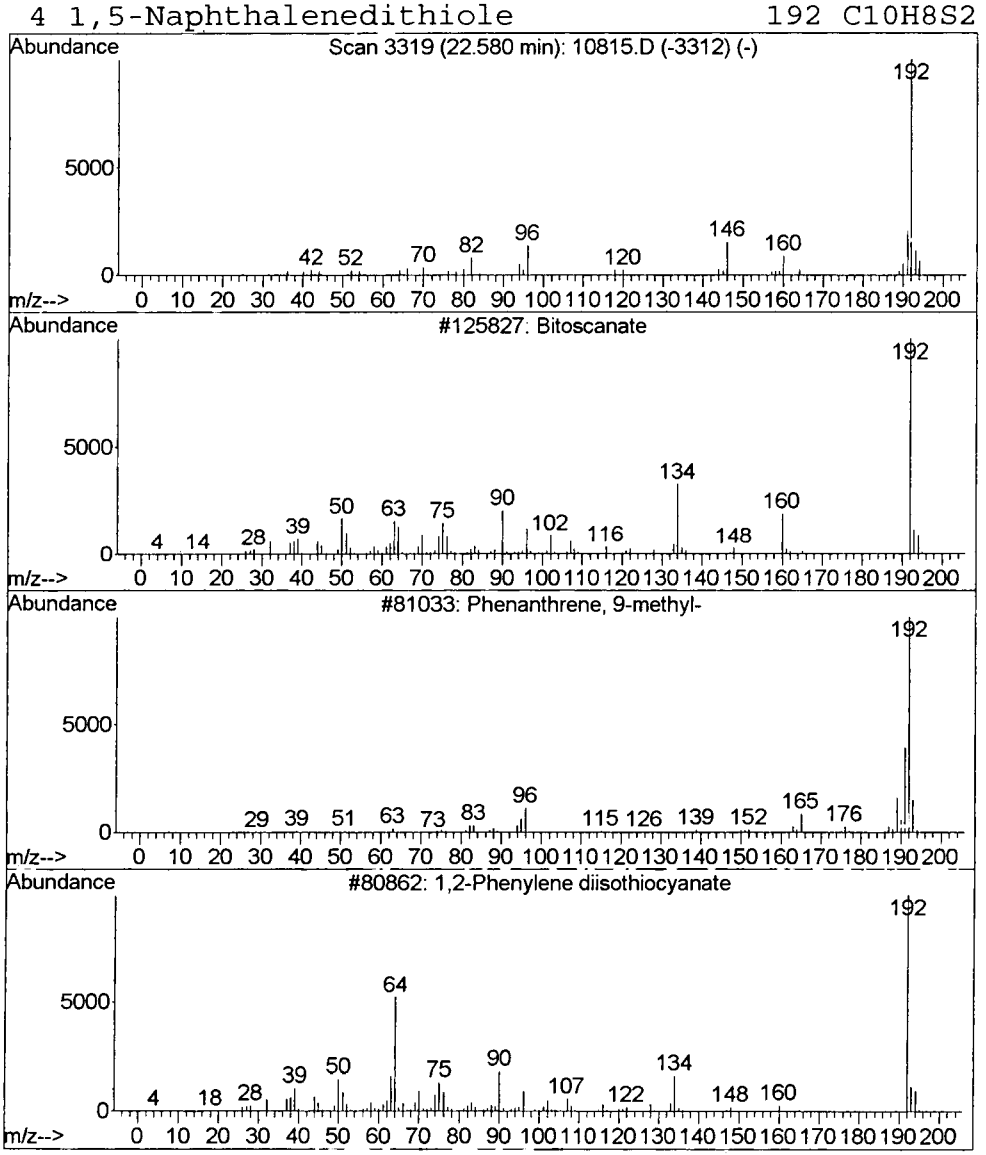
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

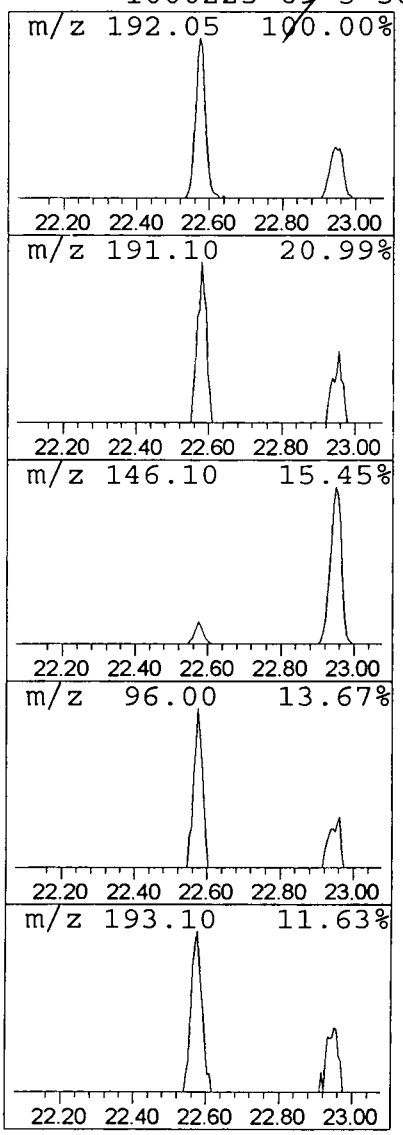
****

Peak Number 6 Bitoscanate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.58	0.33 ug/L	477806	Phenanthranene-d10	22.95

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bitoscanate	192	C8H4N2S2	004044-65-9	52
2	Phenanthrene, 9-methyl-	192	C15H12	000883-20-5	50
3	1,2-Phenylene diisothiocyanate	192	C8H4N2S2	071105-17-4	50
4	1,5-Naphthalenedithiole	192	C10H8S2	1000223-69-5	50





Data File : C:\MSDCHEM\1\DATA\051402\10815.D Vial: 25
Acq On : 15 May 2002 9:55 am Operator: Mclean
Sample : 5/7 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

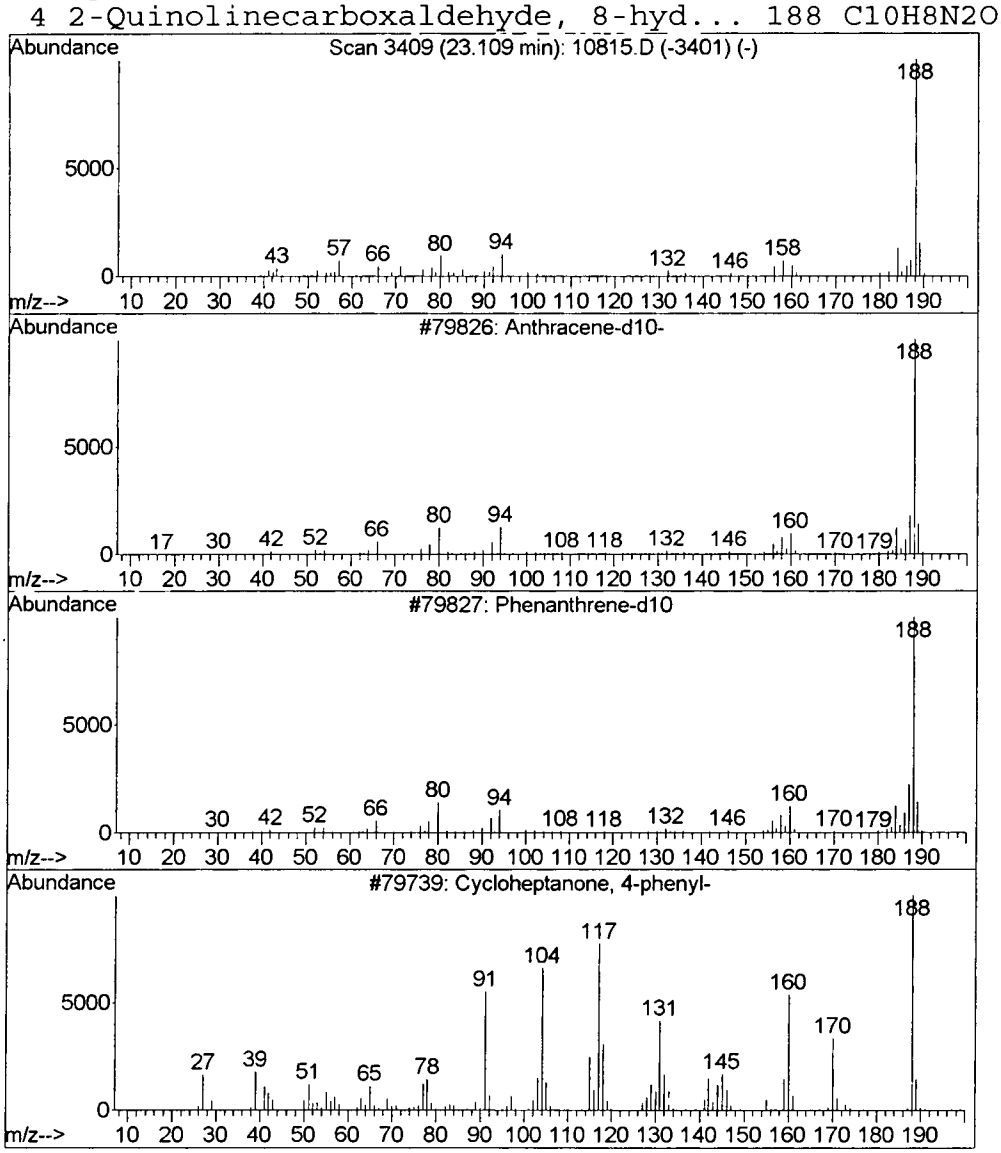
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

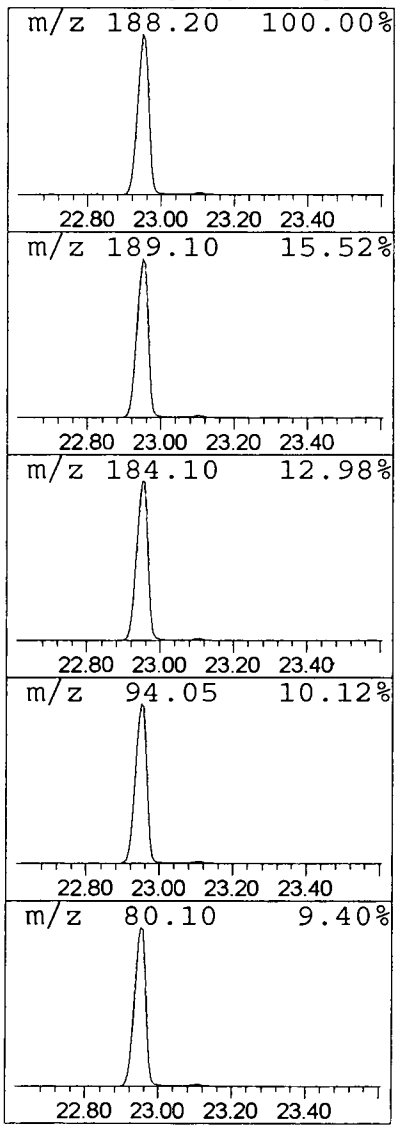
****

Peak Number 7 Anthracene-d10- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.11	0.50 ug/L	734987	Phenanthranene-d10	22.95

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	Cycloheptanone, 4-phenyl-	188	C13H16O	067688-28-2	50
4	2-Quinolinecarboxaldehyde, 8-hyd...	188	C10H8N2O2	005603-22-5	40





Data File : C:\MSDCHEM\1\DATA\051402\10815.D Vial: 25
Acq On : 15 May 2002 9:55 am Operator: Mclean
Sample : 5/7 kb Inst : Instrumen
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.e

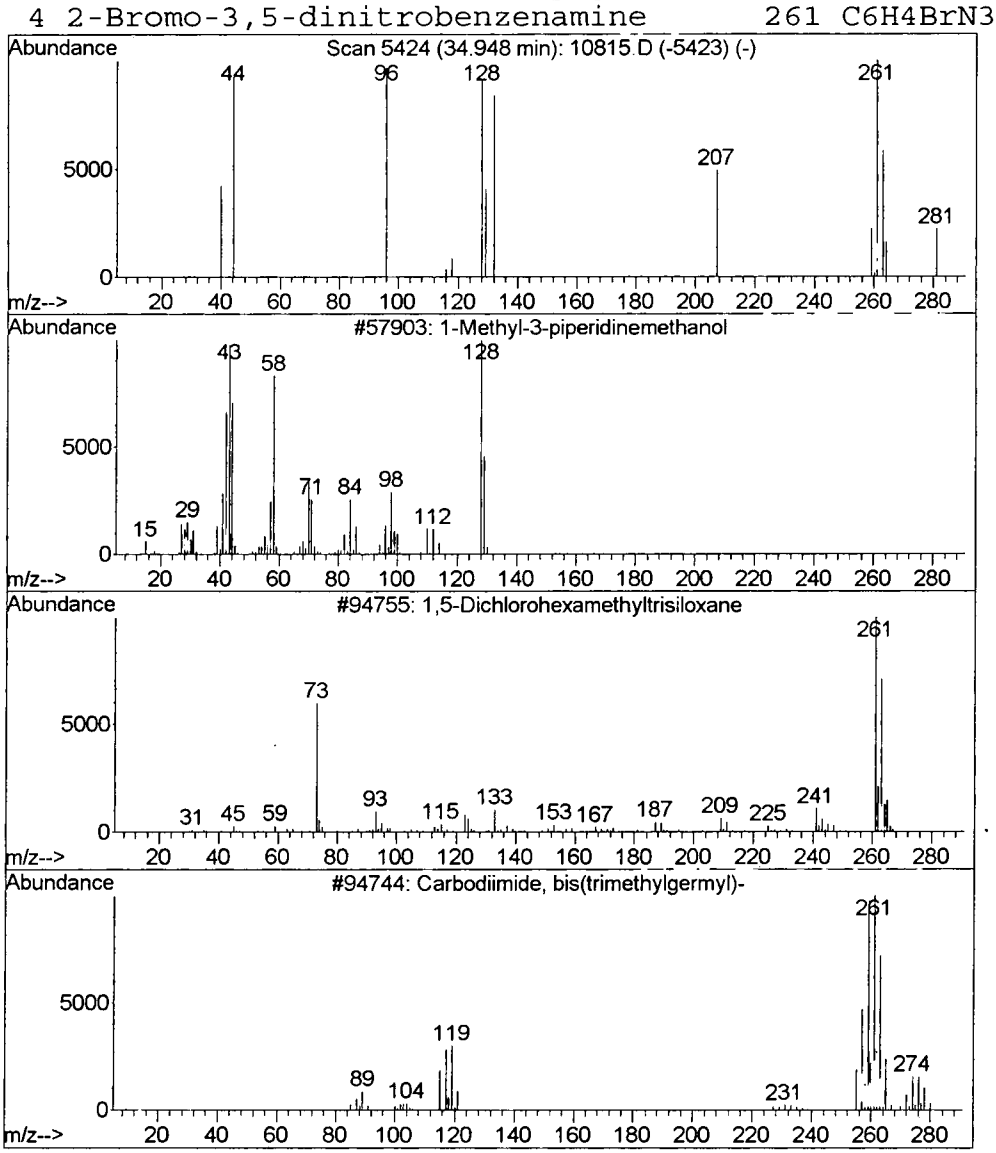
Quant Method : C:\MSDCHEM\1\METHODS\050102.M (Chemstation Integrator)
Title : 508 Modified GC/MS scan
Library : C:\DATABASE\NIST98.L

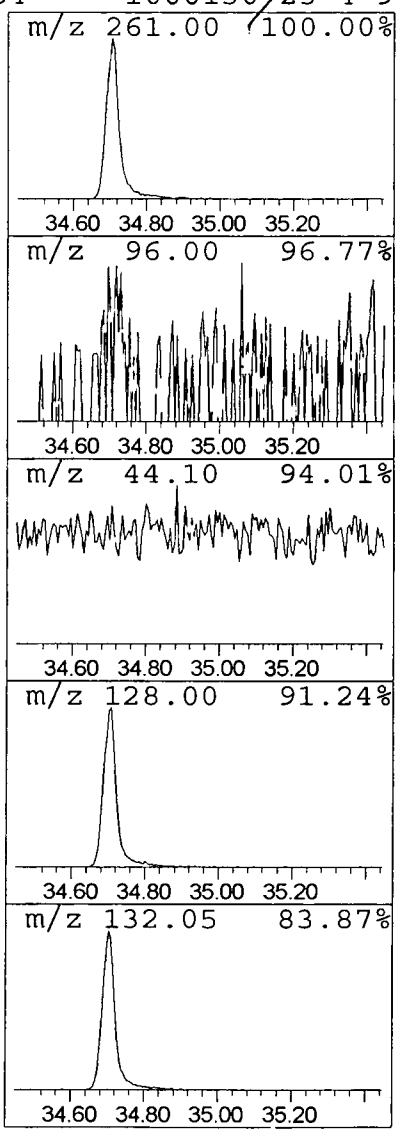
****

Peak Number 8 1-Methyl-3-piperidinemethanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
34.95	0.24 ug/L	218109	Perylene-d12	34.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Methyl-3-piperidinemethanol	129	C7H15NO	007583-53-1	9
2	1,5-Dichlorohexamethyltrisiloxane	276	C6H18Cl2O2Si3	003582-71-6	9
3	Carbodiimide, bis(trimethylgermyl)-	278	C7H18Ge2N2	059579-33-8	9
4	2-Bromo-3,5-dinitrobenzenamine	261	C6H4BrN3O4	1000130-23-4	9





Operator ID: Mclean Date Acquired: 15 May 2002 9:55 am
Data File: C:\MSDCHEM\1\DATA\051402\10815.D
Name: 5/7 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--- #	RT	Resp	Conc
Acetonitrile, dic...	3.63	0.4	ug/L	401875	1	9.94	41898700	40.0
Acetic acid, dich...	3.75	0.2	ug/L	260091	1	9.94	41898700	40.0
2-Pentanone, 4-hy...	5.97	13.4	ug/L	14066500	1	9.94	41898700	40.0
Acetone	7.73	0.3	ug/L	319813	1	9.94	41898700	40.0
Cyclotetrasiloxan...	9.25	0.2	ug/L	241472	1	9.94	41898700	40.0
Bitoscanate	22.58	0.3	ug/L	477806	4	22.95	58609500	40.0
Anthracene-d10-	23.11	0.5	ug/L	734987	4	22.95	58609500	40.0
1-Methyl-3-piperi...	34.95	0.2	ug/L	218109	6	34.71	35736900	40.0