User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 02/25/2002 Time: 15:45:53

Sample: AD31255
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 2/25/02 15:42 Ended: 2/25/02 15:42 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		92		5.0

Comments for Sample AD31255 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 02-25-2002 Time: 15:46:02

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds. This sample was extracted twice with Surrogate Standard recoveries of 74% and 92%. The LCS Standards for both extractions had an unusually high number of compounds with results below their acceptance ranges. This sample will receive an analysis cost discount.

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\JAN0802\31255.D Vial: 9
Acq On : 9 Jan 2002 9:35 pm Operator: 209 GALOS
Sample : gcms scan 12/24 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Feb 6 10:03 2002 Quant Results File: GCMS0103.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Thu Dec 13 14:40:46 2001
Response via : Initial Calibration
DataAcq Meth : GCMS0103

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	11.53	152	858927	20.00	ug/l	0.00
10) Naphthalene-d8	15.13	136	3466709	20.00	ug/l	0.00
17) Acenaphthene-d10	20.40	164	1765499	20.00	ug/l	0.00
29) Phenanthrene-d10	24.81	188	2499005	20.00	ug/l	0.00
38) Chrysene-d12	32.84	240	1469242	20.00	ug/l	0.00
44) Perylene-d12	36.90	264	851932	20.00	ug/l	0.00
System Monitoring Compounds						
37) 2,4-DDD	29.89	235	106030	3.70	ug/mL	-0.18
Spiked Amount 5.000			Recovery = 74.00%			
Target Compounds						Qvalue
2) N-nitroso-dimethyl amine	5.16	74	302	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-Chloropropan	0.00	45	0	N.D.		
8) N-Nitroso-di-n-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)methane	0.00	93	0	N.D.		
14) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.		
15) Naphthalene	15.18	128	1786	N.D.		
16) Hexachlorobutadiene	0.00	225	0	N.D.		
18) Hexachlorocyclopentadiene	0.00	237	0	N.D.		
19) 2-Chloronaphthalene	0.00	162	0	N.D.		
20) Dimethylphthalate	0.00	163	0	N.D.		
21) 2,6-Dinitrotoluene	19.96	165	240	N.D.		
22) Acenaphthylene	0.00	152	0	N.D.		
23) Acenaphthene	0.00	153	0	N.D.		
24) 2,4-Dinitrotoluene	0.00	165	0	N.D.		
25) Diethylphthalate	21.92	149	1620	N.D.		
26) 4-Chlorophenyl-phenylether	0.00	204	0	N.D.		
27) Fluorene	0.00	166	0	N.D.		
28) Azobenzen/ 1,2-Diphenylhyd	0.00	77	0	N.D.		

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

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\JAN0802\31255.D Vial: 9
Acq On : 9 Jan 2002 9:35 pm Operator: 209 GALOS
Sample : gcms scan 12/24 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Feb 6 10:03 2002 Quant Results File: GCMS0103.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Thu Dec 13 14:40:46 2001
Response via : Initial Calibration
DataAcq Meth : GCMS0103

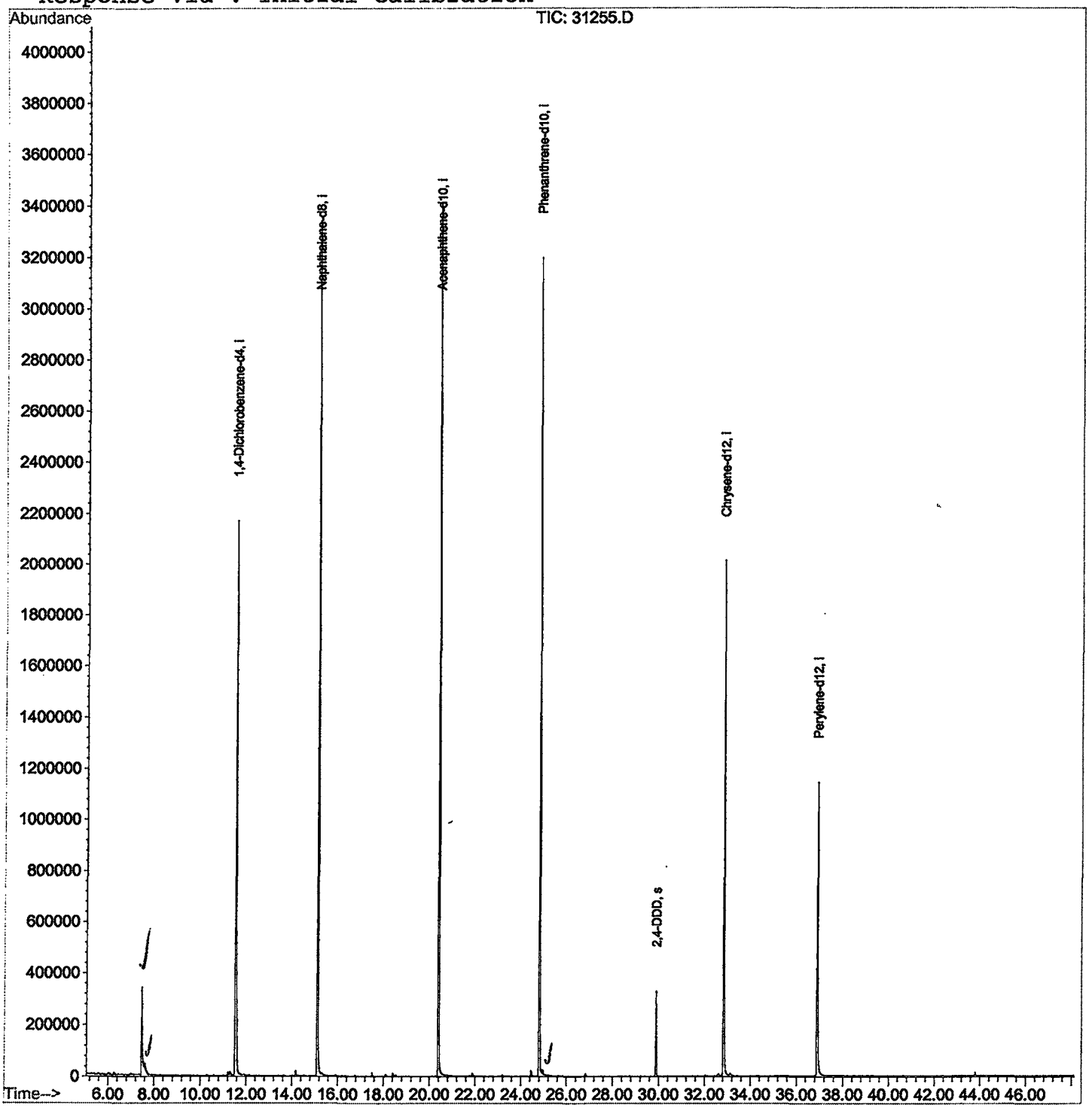
	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
30)	N-Nitrosodiphenylamine	0.00	168	0	N.D.	
31)	4-Bromophenyl-phenylether	0.00	248	0	N.D.	
32)	Hexachlorobenzene	0.00	284	0	N.D.	
33)	Phenanthrene	24.88	178	995	N.D.	
34)	Anthracene	24.88	178	995	N.D.	
35)	Di-n-butylphthalate	26.83	149	11615	N.D.	
36)	Fluoranthene	28.50	202	417	N.D.	
39)	Pyrene	29.18	202	191	N.D.	
40)	Butylbenzylphthalate	0.00	149	0	N.D.	
41)	Benzo(a)anthracene	32.84	228	4188	N.D.	
42)	Bis(2-ethylhexyl)phthalate	33.14	149	4907	N.D.	
43)	Chrysene	32.84	228	4188	N.D.	
45)	Di-n-Octylphthalate	0.00	149	0	N.D.	
46)	Benzo(b)fluoranthene	35.98	252	504	N.D.	
47)	Benzo(k)fluoranthene	35.98	252	504	N.D.	
48)	Benzo(a)pyrene	36.91	252	3677	N.D.	
49)	Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.	
50)	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

Quantitation Report

Data File : C:\HPCHEM\1\DATA\JAN0802\31255.D Vial: 9
Acq On : 9 Jan 2002 9:35 pm Operator: 209 GALOS
Sample : gcms scan 12/24 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Feb 6 10:03 2002 Quant Results File: GCMS0103.RES

Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Mon Jan 07 10:31:02 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\JAN0802\31255.D Vial: 9
Acq On : 9 Jan 2002 9:35 pm Operator: 209 GALOS
Sample : gcms scan 12/24 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

Signal : TIC

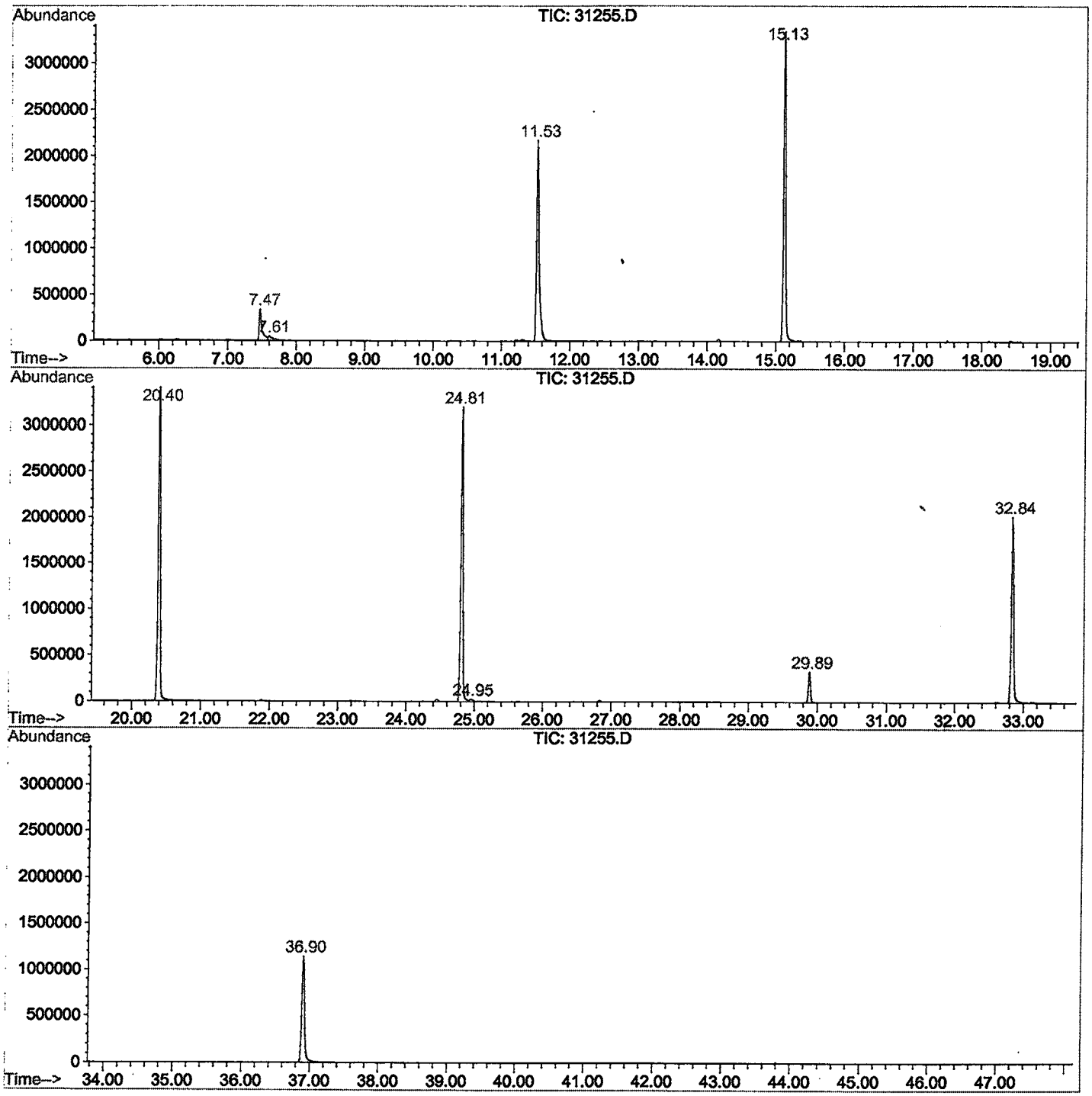
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	7.470	260	264	276	rBV	337054	927709	12.52%	2.524%
2	7.608	276	279	296	rVB2	44278	182020	2.46%	0.495%
3	11.528	701	706	727	rBV	2167898	5464380	73.74%	14.866%
4	15.126	1092	1098	1116	rBV	3339816	7083199	95.58%	19.270%
5	20.396	1665	1672	1691	rBV	3413071	7410509	100.00%	20.161%
6	24.812	2146	2153	2165	rBV2	3197692	7015255	94.67%	19.085%
7	24.949	2165	2168	2184	rVB	22540	81081	1.09%	0.221%
8	29.888	2700	2706	2711	rBV	328301	664933	8.97%	1.809%
9	32.835	3021	3027	3048	rBV2	2012584	4739053	63.95%	12.893%
10	36.902	3462	3470	3497	rBV2	1143432	3189153	43.04%	8.676%

Sum of corrected areas: 36757292

31255.D GCMS0103.M Wed Feb 06 11:23:21 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\JAN0802\31255.D
Operator : 209 GALOS
Acquired : 9 Jan 2002 9:35 pm using AcqMethod GCMS0103
Instrument : GC/MS 209
Sample Name: gcms scan 12/24
Misc Info :
Vial Number: 9
Quant File :GCMS0103.RES (RTE Integrator)



31255.D GCMS0103.M Wed Feb 06 11:23:27 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\JAN0802\31255.D
Acq On : 9 Jan 2002 9:35 pm
Sample : gcms scan 12/24
Misc :
MS Integration Params: LSCINT.P

Vial: 9
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

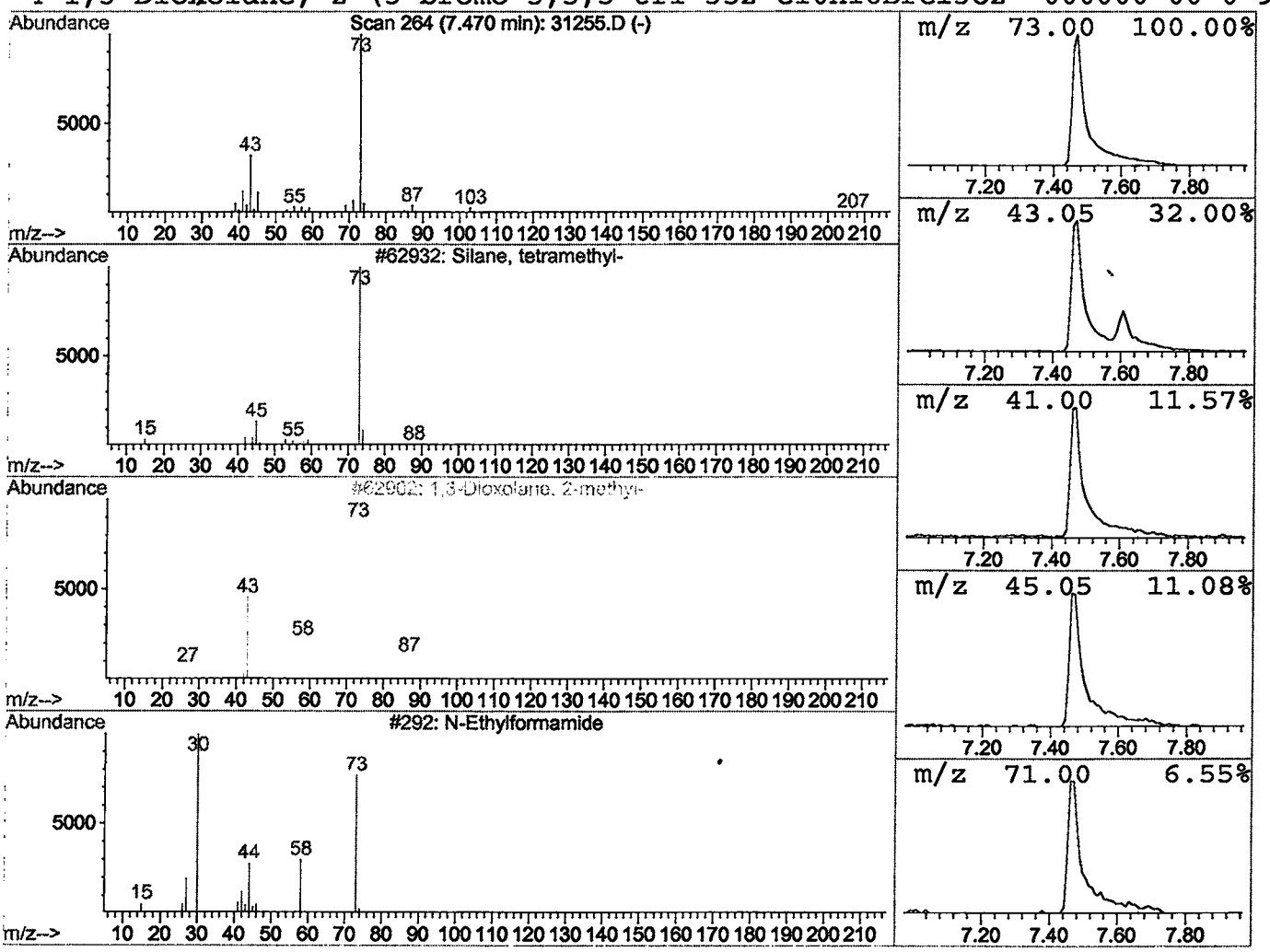
Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Silane, tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.47	3.40 ug/l	927709	1,4-Dichlorobenzene-d4	11.53

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
2	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
3	N-Ethylformamide	73	C3H7NO	000627-45-2	9
4	1,3-Dioxolane, 2-(3-bromo-5,5,5-tri	352	C10H16BrCl3O2	000000-00-0	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\JAN0802\31255.D Vial: 9
Acq On : 9 Jan 2002 9:35 pm Operator: 209 GALOS
Sample : gcms scan 12/24 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

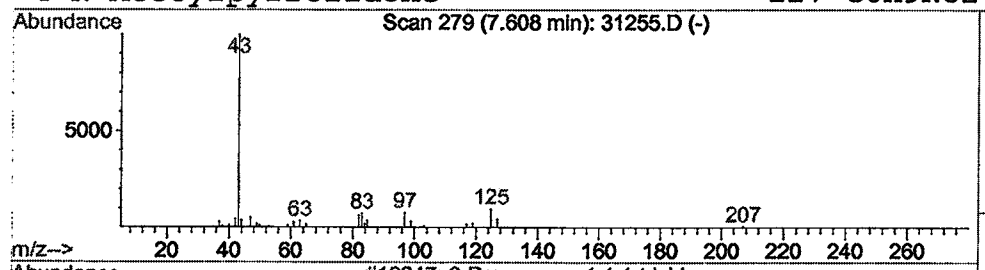
Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

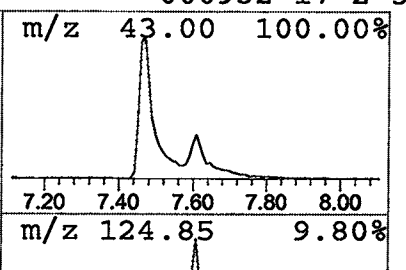
****

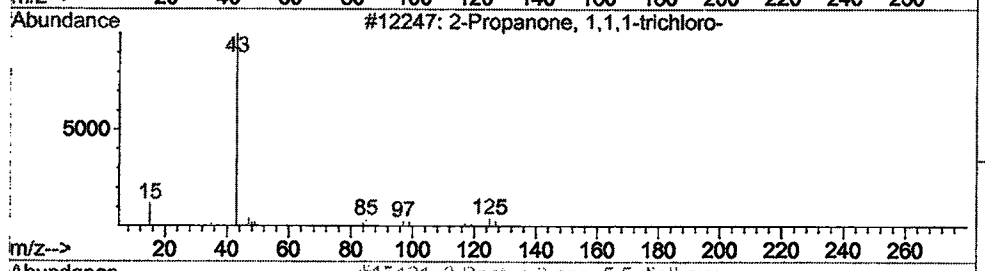
Peak Number 2 2-Propanone, 1,1,1-trichloro- Concentration Rank 2

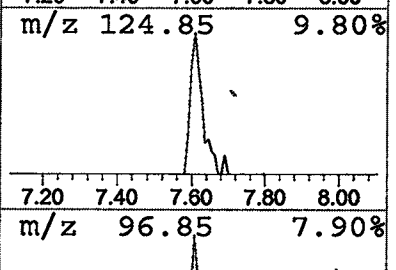
R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	0.67 ug/l	182020	1,4-Dichlorobenzene-d4	11.53

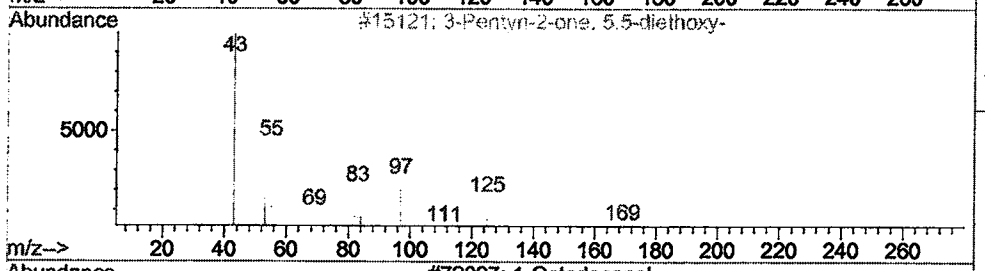
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanone, 1,1,1-trichloro-	160	C3H3Cl3O	000918-00-3	56
2	3-Pentyn-2-one, 5,5-diethoxy-	170	C9H14O3	055402-04-5	9
3	1-Octadecanol	270	C18H38O	000112-92-5	7
4	N-Acetylpyrrolidone	127	C6H9NO2	000932-17-2	5

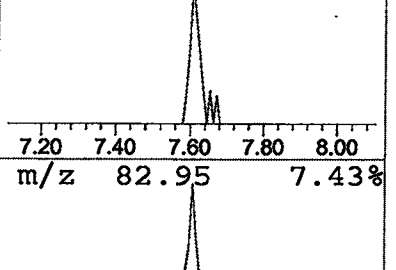


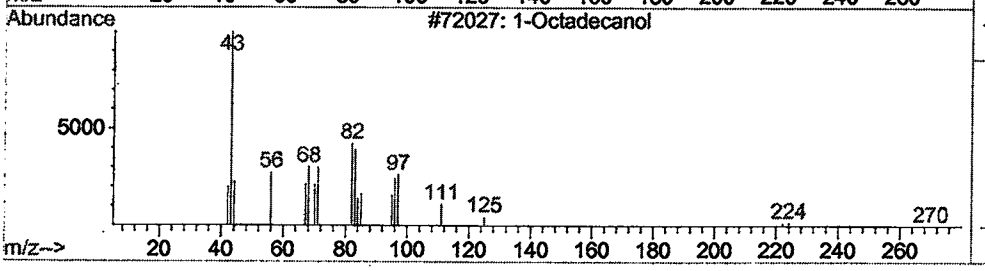


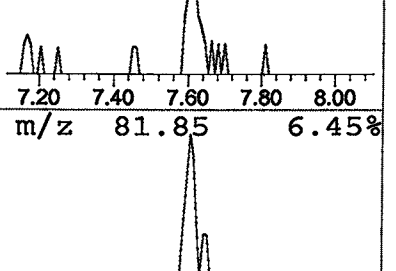












Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\JAN0802\31255.D Vial: 9
Acq On : 9 Jan 2002 9:35 pm Operator: 209 GALOS
Sample : gcms scan 12/24 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

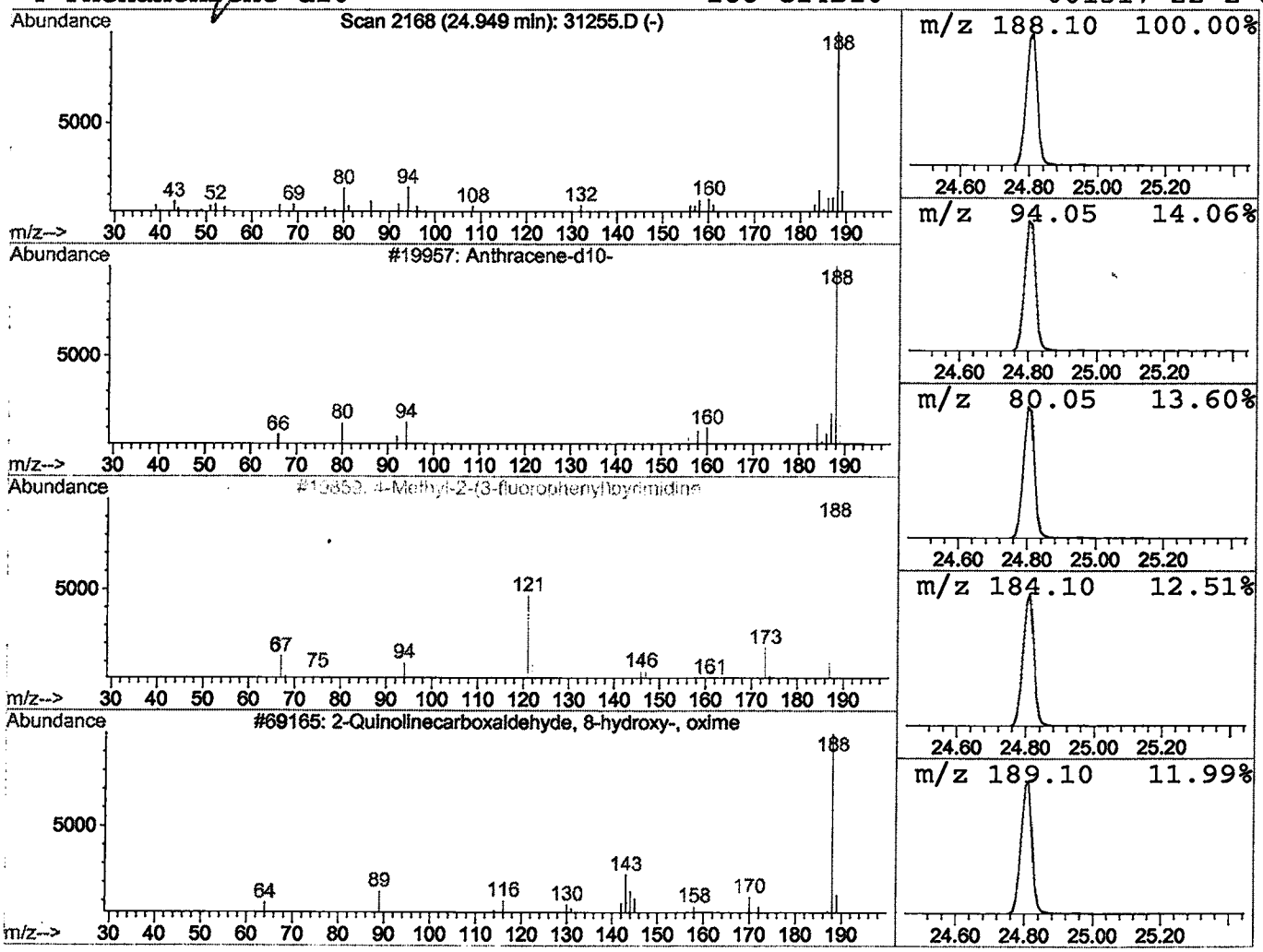
Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 Anthracene-d10- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.95	0.23 ug/l	81081	Phenanthrene-d10	24.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene-d10	188	C14D10	001719-06-8	90
2	4-Methyl-2-(3-fluorophenyl)pyrimidi	188	C11H9FN2	076128-66-0	45
3	2-Quinolinecarboxaldehyde, 8-hydrox	188	C10H8N2O2	005603-22-5	40
4	Phenanthrene-d10	188	C14D10	001517-22-2	38



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 9 Jan 2002 9:35 pm
Data File: C:\HPCHEM\1\DATA\JAN0802\31255.D
Name: gcms scan 12/24
Misc:
Method: C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title: 209 625/TCLP calibration table
Library Searched: C:\DATABASE\NBS75K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
Silane, tetramethyl-	7.47	3.4	ug/l	927709	ISTD01	11.53	5464380	20.0
2-Propanone, 1,1,1-t	7.61	0.7	ug/l	182020	ISTD01	11.53	5464380	20.0
Anthracene-d10	24.95	0.2	ug/l	81081	ISTD04	24.81	7015260	20.0

31255.D GCMS0103.M Wed Feb 06 11:23:46 2002

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\FEB1402\31255.D Vial: 28
Acq On : 15 Feb 2002 5:23 am Operator: 209 GALOS
Sample : GC/MS SCAN 2/11 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Feb 15 12:01 2002 Quant Results File: GCMS0208.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Wed Feb 13 11:43:23 2002
Response via : Initial Calibration
DataAcq Meth : GCMS0208

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	12.34	152	378051	20.00	ug/l	-0.01
10) Naphthalene-d8	15.99	136	1489548	20.00	ug/l	-0.01
17) Acenaphthene-d10	21.30	164	798204	20.00	ug/l	-0.01
29) Phenanthrene-d10	25.76	188	1199732	20.00	ug/l	-0.01
38) Chrysene-d12	33.83	240	862376	20.00	ug/l	-0.02
44) Perylene-d12	38.10	264	543414	20.00	ug/l	-0.02
System Monitoring Compounds						
37) 2,4-DDD	30.83	235	55435	4.60	ug/mL	0.33
Spiked Amount 5.000			Recovery = 92.00%			
Target Compounds						Qvalue
2) N-nitroso-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-Chloropropan	0.00	45	0	N.D.		
8) N-Nitroso-di-n-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)methane	0.00	93	0	N.D.		
14) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.		
15) Naphthalene	16.03	128	297	N.D.		
16) Hexachlorobutadiene	0.00	225	0	N.D.		
18) Hexachlorocyclopentadiene	0.00	237	0	N.D.		
19) 2-Chloronaphthalene	0.00	162	0	N.D.		
20) Dimethylphthalate	0.00	163	0	N.D.		
21) 2,6-Dinitrotoluene	0.00	165	0	N.D.		
22) Acenaphthylene	0.00	152	0	N.D.		
23) Acenaphthene	0.00	153	0	N.D.		
24) 2,4-Dinitrotoluene	0.00	165	0	N.D.		
25) Diethylphthalate	22.78	149	335	N.D.		
26) 4-Chlorophenyl-phenylether	0.00	204	0	N.D.		
27) Fluorene	0.00	166	0	N.D.		
28) Azobenzen/ 1,2-Diphenylhyd	0.00	77	0	N.D.		

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

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\FEB1402\31255.D Vial: 28
Acq On : 15 Feb 2002 5:23 am Operator: 209 GALOS
Sample : GC/MS SCAN 2/11 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Feb 15 12:01 2002 Quant Results File: GCMS0208.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Wed Feb 13 11:43:23 2002
Response via : Initial Calibration
DataAcq Meth : GCMS0208

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
30)	N-Nitrosodiphenylamine	0.00	168	0	N.D.	
31)	4-Bromophenyl-phenylether	0.00	248	0	N.D.	
32)	Hexachlorobenzene	0.00	284	0	N.D.	
33)	Phenanthrene	25.81	178	1085	N.D.	
34)	Anthracene	25.81	178	1563	N.D.	
35)	Di-n-butylphthalate	27.71	149	6942	N.D.	
36)	Fluoranthene	29.45	202	248	N.D.	
39)	Pyrene	30.12	202	115	N.D.	
40)	Butylbenzylphthalate	0.00	149	0	N.D.	
41)	Benzo(a)anthracene	33.83	228	2139	N.D.	
42)	Bis(2-ethylhexyl)phthalate	34.02	149	2534	N.D.	
43)	Chrysene	33.83	228	2139	N.D.	
45)	Di-n-Octylphthalate	0.00	149	0	N.D.	
46)	Benzo(b)fluoranthene	0.00	252	0	N.D.	
47)	Benzo(k)fluoranthene	0.00	252	0	N.D.	
48)	Benzo(a)pyrene	38.11	252	2230	N.D.	
49)	Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.	
50)	Dibenz(a,h)anthracene	0.00	278	0	N.D.	
51)	Benzo(g,h,i)perylene	0.00	276	0	N.D.	

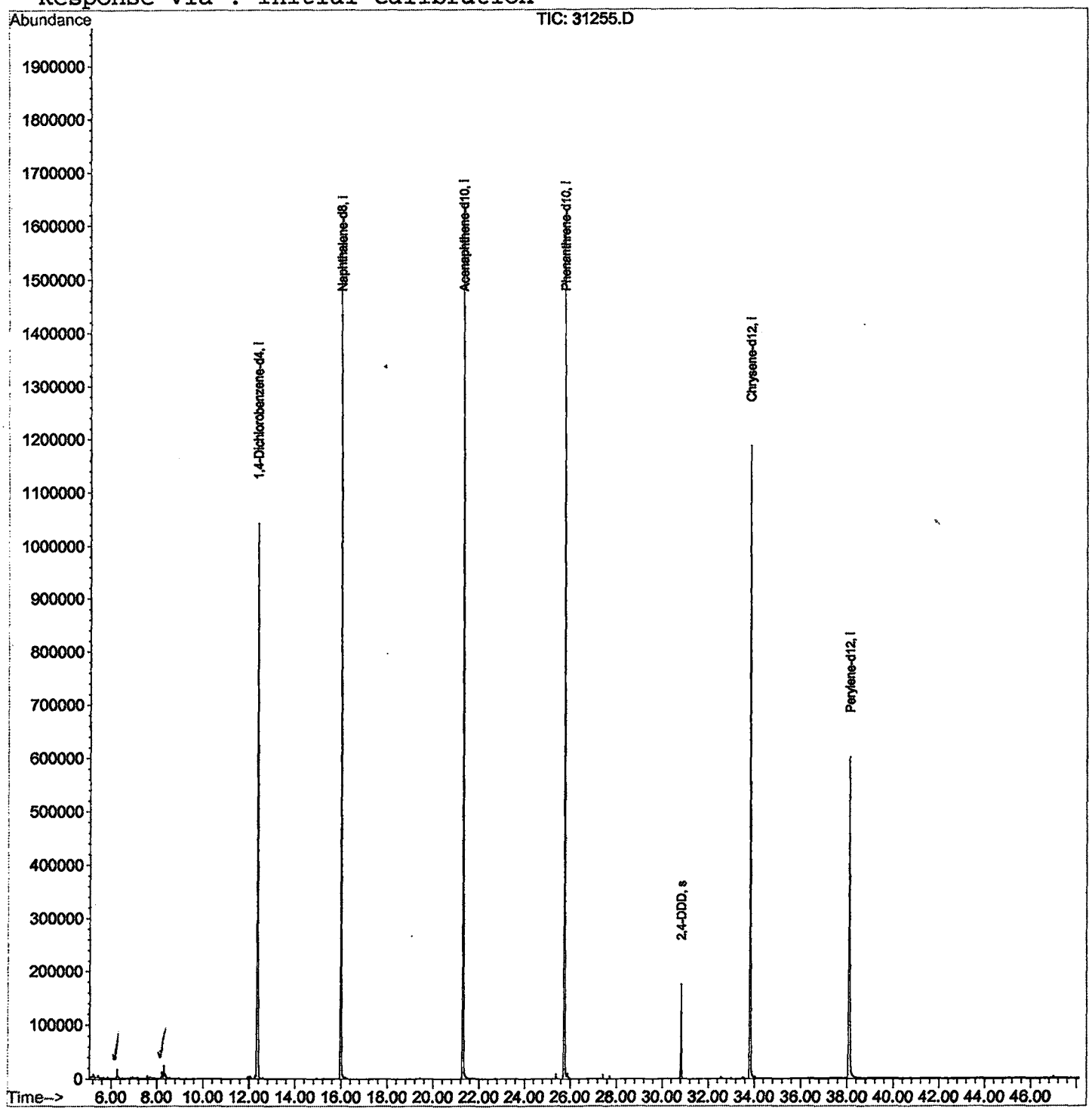
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\FEB1402\31255.D
Acq On : 15 Feb 2002 5:23 am
Sample : GC/MS SCAN 2/11
Misc :
MS Integration Params: GOOFY.P
Quant Time: Feb 15 12:01 2002

Vial: 28
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00
Quant Results File: GCMS0208.RES

Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Wed Feb 13 11:43:23 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\FEB1402\31255.D Vial: 28
Acq On : 15 Feb 2002 5:23 am Operator: 209 GALOS
Sample : GC/MS SCAN 2/11 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

Signal : TIC

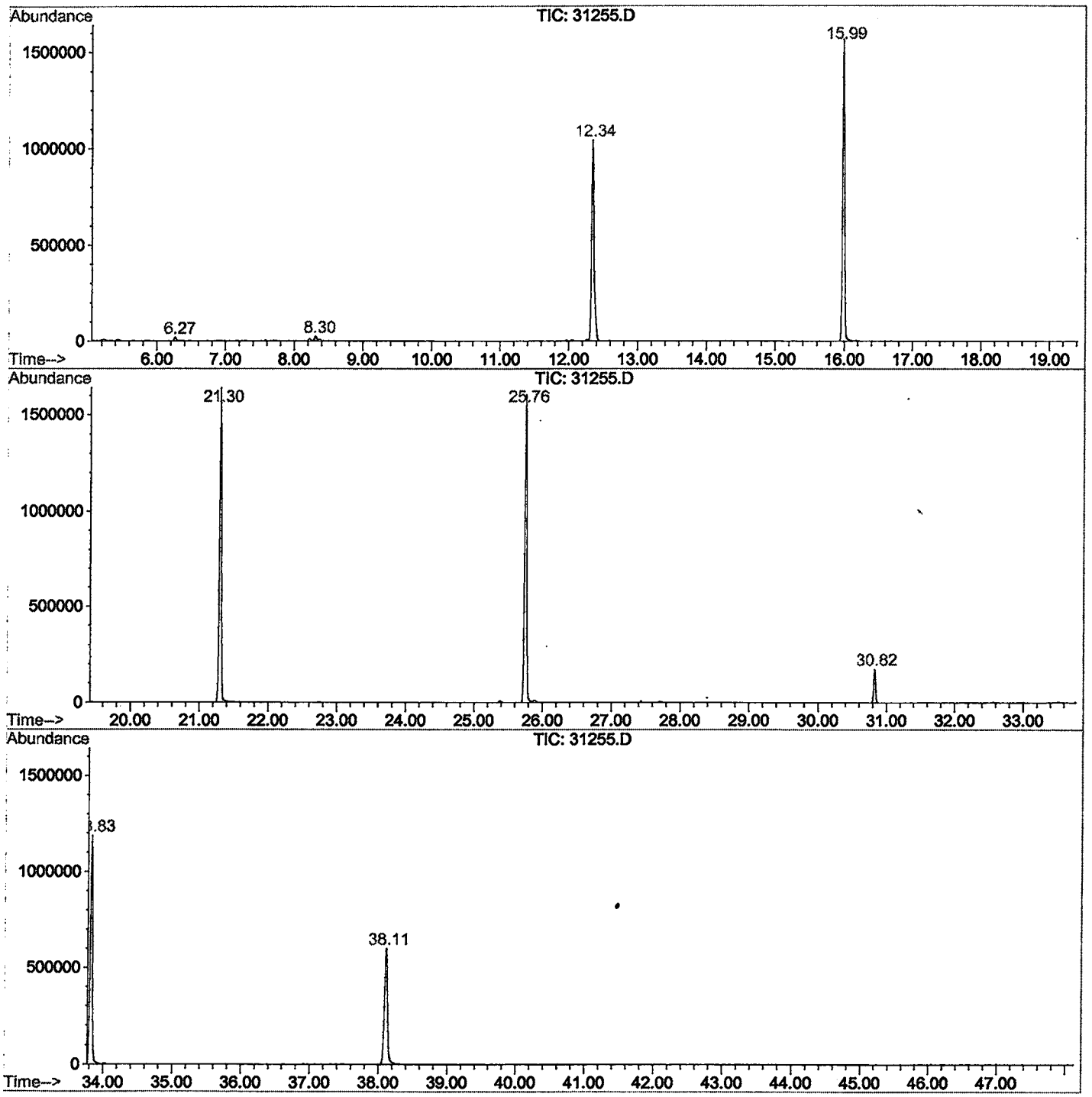
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	6.266	128	133	137	rBV	17353	35177	1.05%	0.205%
2	8.304	352	355	360	rBV	22260	48349	1.45%	0.282%
3	12.344	789	795	809	rVB	1041993	2465438	73.75%	14.365%
4	15.988	1186	1192	1208	rBV	1563970	3099645	92.72%	18.060%
5	21.304	1764	1771	1794	rBV	1642264	3343180	100.00%	19.479%
6	25.756	2249	2256	2266	rBV	1604029	3328289	99.55%	19.392%
7	30.824	2803	2808	2815	rVB	174952	339664	10.16%	1.979%
8	33.826	3128	3135	3153	rBV	1185669	2625715	78.54%	15.299%
9	38.113	3592	3602	3624	rBV2	598912	1877572	56.16%	10.940%

Sum of corrected areas: 17163029

31255.D GCMS0208.M Fri Feb 15 14:04:05 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\FEB1402\31255.D
Operator : 209 GALOS
Acquired : 15 Feb 2002 5:23 am using AcqMethod GCMS0208
Instrument : GC/MS 209
Sample Name: GC/MS SCAN 2/11
Misc Info :
Vial Number: 28
Quant File :GCMS0208.RES (RTE Integrator)



31255.D GCMS0208.M Fri Feb 15 14:04:10 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\FEB1402\31255.D
Acq On : 15 Feb 2002 5:23 am
Sample : GC/MS SCAN 2/11
Misc :
MS Integration Params: LSCINT.P

Vial: 28
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

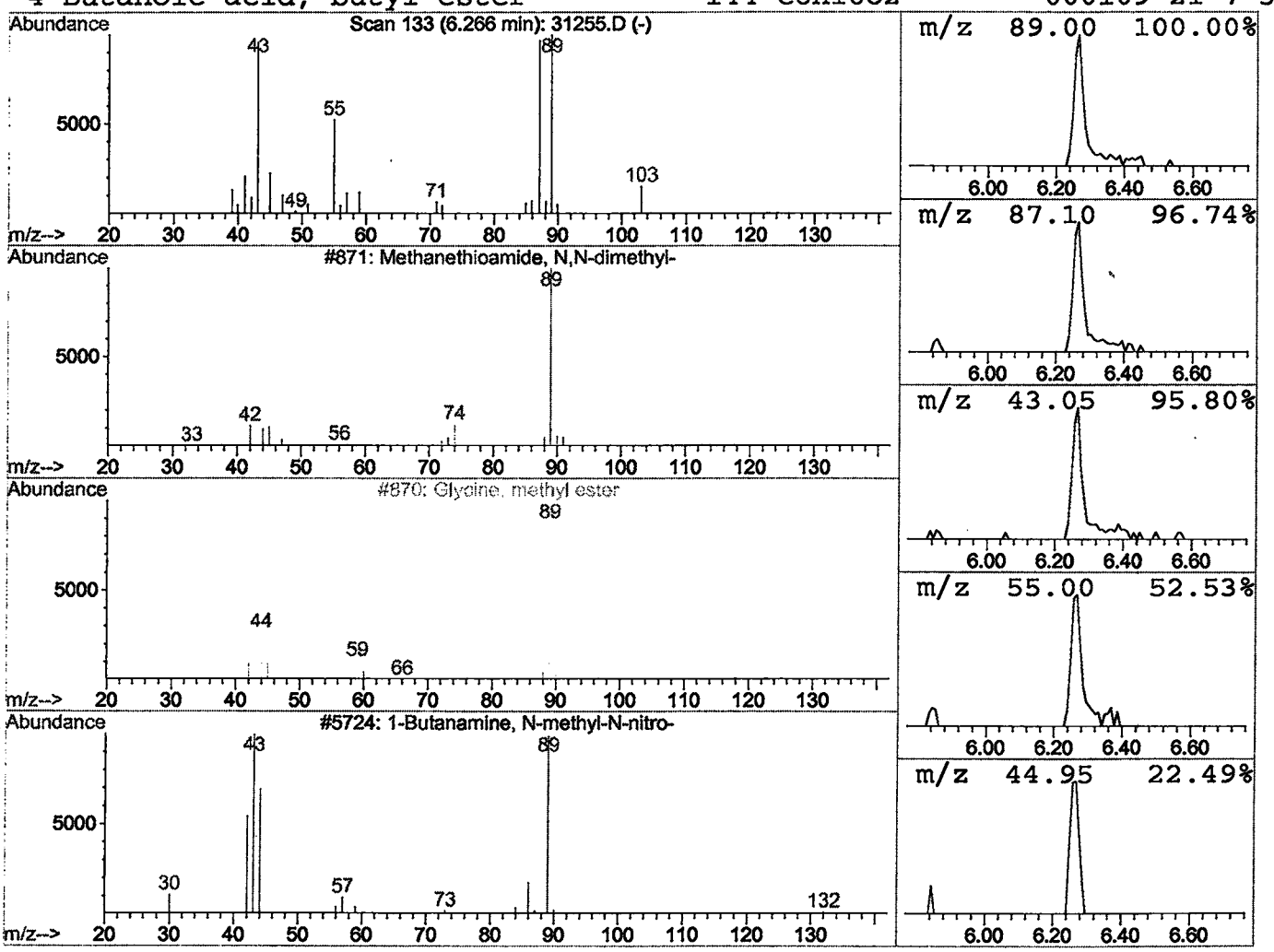
Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Methanethioamide, N,N-dimethyl Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.27	0.29 ug/l	35177	1,4-Dichlorobenzene-d4	12.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methanethioamide, N,N-dimethyl-	89	C3H7NS	000758-16-7	53
2	Glycine, methyl ester	89	C3H7NO2	000616-34-2	38
3	1-Butanamine, N-methyl-N-nitro-	132	C5H12N2O2	052330-07-1	36
4	Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	36



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\FEB1402\31255.D
Acq On : 15 Feb 2002 5:23 am
Sample : GC/MS SCAN 2/11
Misc :
MS Integration Params: LSCINT.P

Vial: 28
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

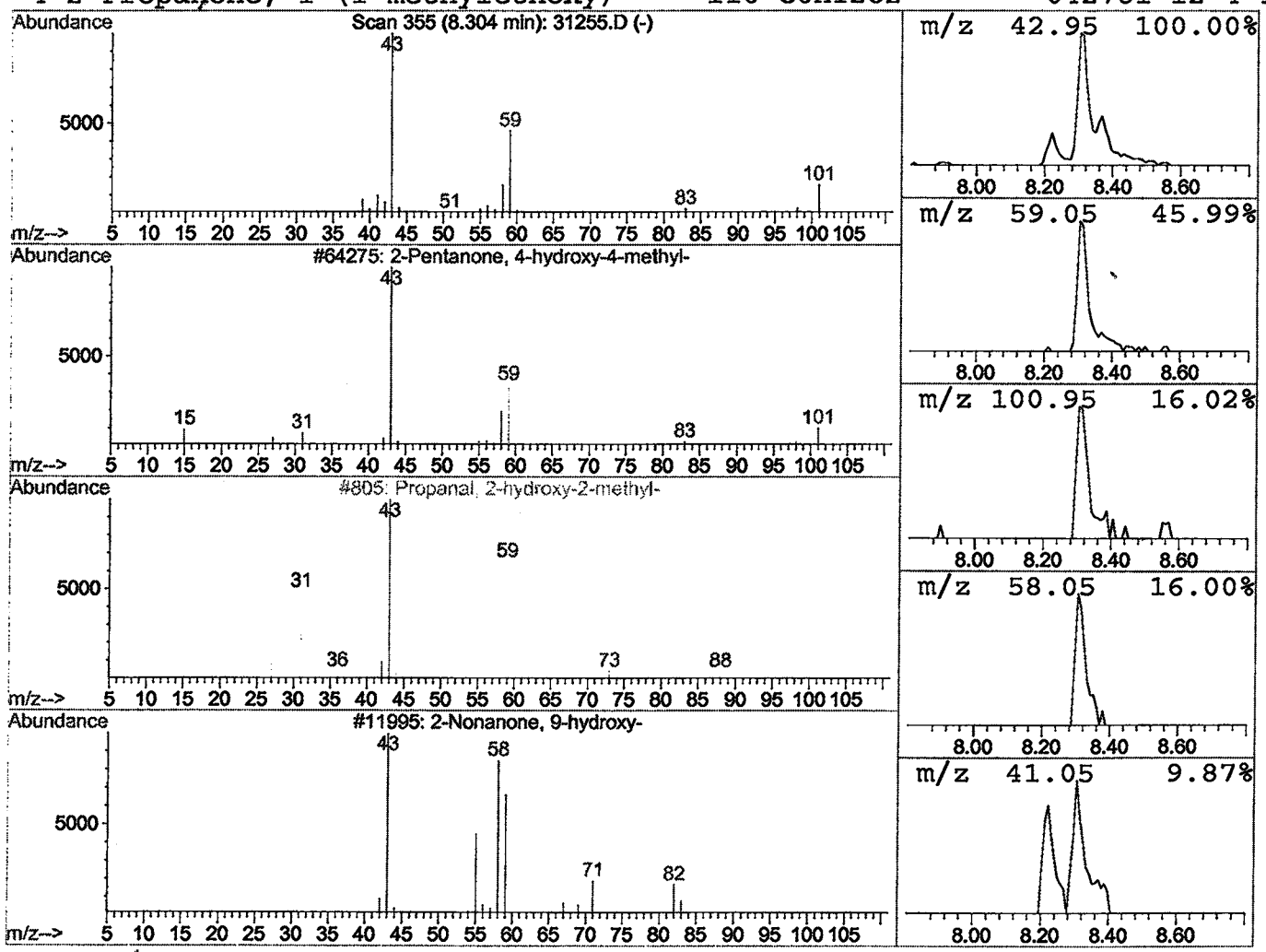
Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 2-Pentanone, 4-hydroxy-4-methy Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.30	0.39 ug/l	48349	1,4-Dichlorobenzene-d4	12.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	Propanal, 2-hydroxy-2-methyl-	88	C4H8O2	020818-81-9	33
3	2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	9
4	2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 15 Feb 2002 5:23 am
Data File: C:\HPCHEM\1\DATA\FEB1402\31255.D
Name: GC/MS SCAN 2/11
Misc:
Method: C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title: 209 625/TCLP calibration table
Library Searched: C:\DATABASE\NBS75K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
Methanethioamide, N,	6.27	0.3	ug/l	35177	ISTD01	12.34	2465440	20.0
2-Pentanone, 4-hydro	8.30	0.4	ug/l	48349	ISTD01	12.34	2465440	20.0

31255.D GCMS0208.M Fri Feb 15 14:04:24 2002