User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 02/25/2002 Time: 15:47:47

Sample: AD31253
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 2/25/02 15:46 Ended: 2/25/02 15:46 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		85		5.0

Comments for Sample AD31253 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 02-25-2002 Time: 15:48:02

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 78% and 85%. 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\31253.D Vial: 10
Acq On : 9 Jan 2002 10:34 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 11:09 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	11.52	152	866504	20.00	ug/l	0.00
10) Naphthalene-d8	15.12	136	3391749	20.00	ug/l	0.00
17) Acenaphthene-d10	20.39	164	1705627	20.00	ug/l	0.00
29) Phenanthrene-d10	24.81	188	2393839	20.00	ug/l	0.00
38) Chrysene-d12	32.84	240	1430776	20.00	ug/l	0.00
44) Perylene-d12	36.91	264	852745	20.00	ug/l	0.01
System Monitoring Compounds						
37) 2,4-DDD	29.88	235	106895	3.89	ug/mL	-0.19
Spiked Amount 5.000			Recovery = 77.80%			
Target Compounds						Qvalue
2) N-nitroso-dimethyl amine	5.15	74	401	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	1212	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	19.80	163	252	N.D.		
21) 2,6-Dinitrotoluene	0.00	165	0	N.D.	d	
22) Acenaphthylene	0.00	152	0	N.D.		
23) Acenaphthene	0.00	153	0	N.D.		
24) 2,4-Dinitrotoluene	20.78	165	203	N.D.		
25) Diethylphthalate	21.92	149	1753	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\31253.D Vial: 10
Acq On : 9 Jan 2002 10:34 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 11:09 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	24.87	178	650	N.D.	
34) Anthracene	24.87	178	650	N.D.	
35) Di-n-butylphthalate	26.83	149	12373	N.D.	
36) Fluoranthene	28.50	202	266	N.D.	
39) Pyrene	29.17	202	105	N.D.	
40) Butylbenzylphthalate	0.00	149	0	N.D.	
41) Benzo(a)anthracene	32.84	228	3982	N.D.	
42) Bis(2-ethylhexyl)phthalate	33.13	149	6907	N.D.	
43) Chrysene	32.84	228	3982	N.D.	
45) Di-n-Octylphthalate	34.89	149	223	N.D.	
46) Benzo(b)fluoranthene	35.94	252	190	N.D.	
47) Benzo(k)fluoranthene	35.94	252	189	N.D.	
48) Benzo(a)pyrene	36.91	252	3606	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.	

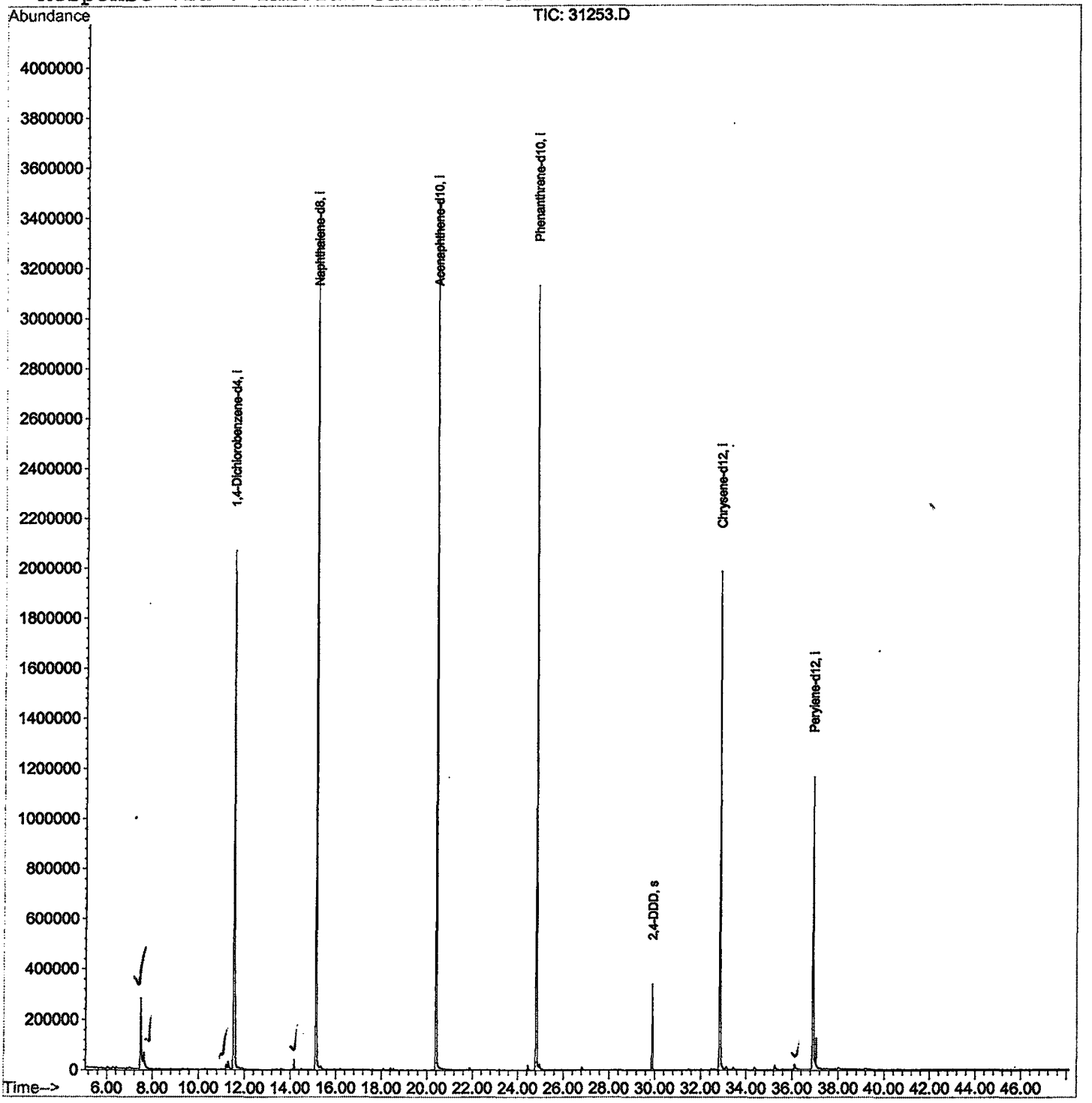
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\JAN0802\31253.D
Acq On : 9 Jan 2002 10:34 pm
Sample : gcms scan 12/24
Misc :
MS Integration Params: GOOFY.P
Quant Time: Feb 6 11:09 2002

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00
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\31253.D Vial: 10
Acq On : 9 Jan 2002 10:34 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

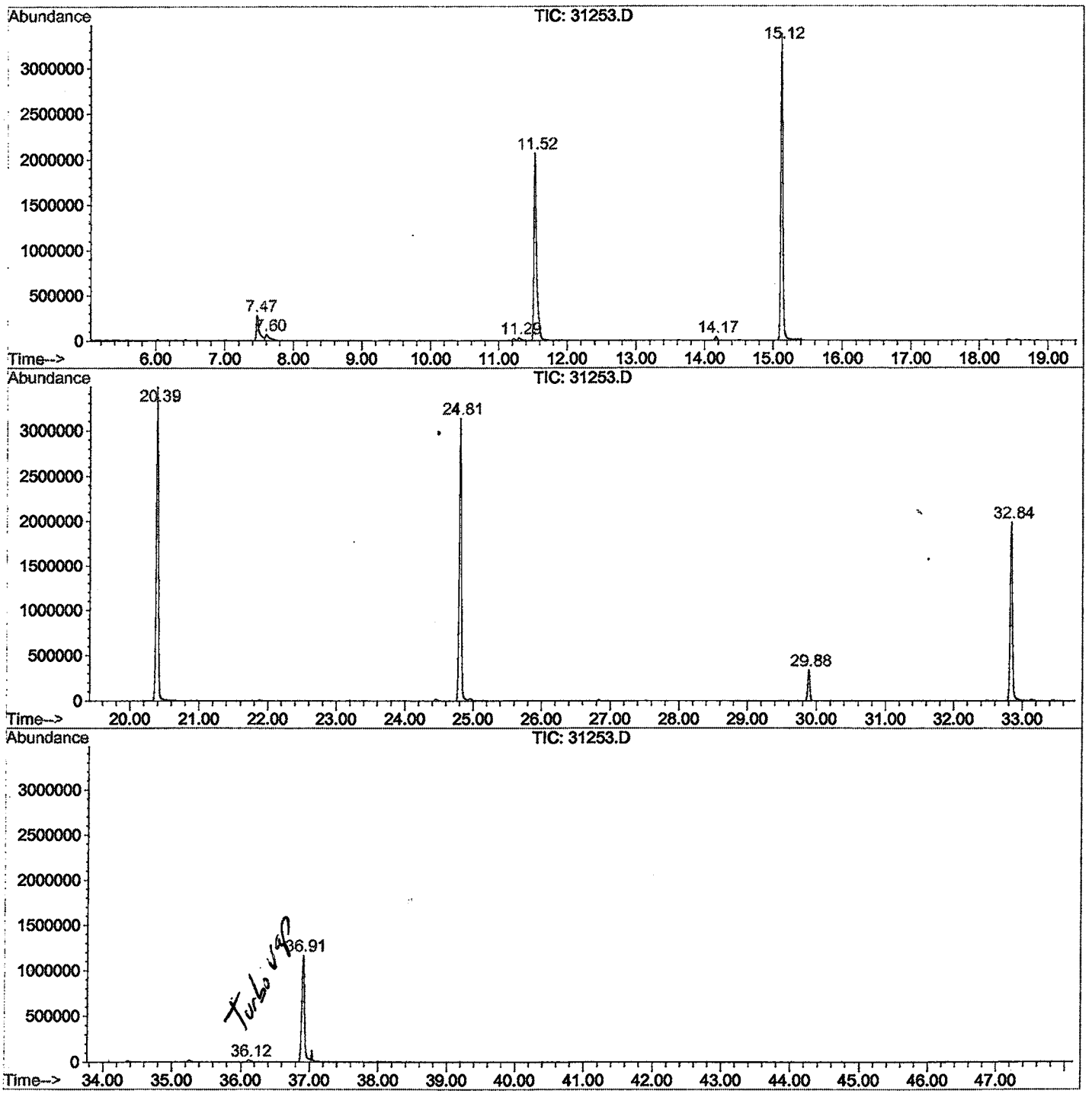
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	7.466	260	264	276	rBV	278465	783451	11.00%	2.173%
2	7.604	276	279	305	rVB	64803	279836	3.93%	0.776%
3	11.294	678	681	699	rVB2	30244	86325	1.21%	0.239%
4	11.524	701	706	725	rBV	2067889	5557491	78.06%	15.415%
5	14.168	988	994	1001	rVB	42043	79651	1.12%	0.221%
6	15.123	1093	1098	1116	rBV	3370249	6919563	97.19%	19.194%
7	20.392	1666	1672	1689	rBV	3475232	7119739	100.00%	19.749%
8	24.808	2147	2153	2165	rBV2	3130596	6723020	94.43%	18.648%
9	29.884	2701	2706	2711	rBV	338475	656802	9.23%	1.822%
10	32.841	3021	3028	3047	rBV2	1983754	4595623	64.55%	12.747%
11	36.118	3378	3385	3393	rBV2	21281	72312	1.02%	0.201%
12	36.907	3463	3471	3484	rBV	1164515	3177510	44.63%	8.814%

Sum of corrected areas: 36051323

31253.D GCMS0103.M Wed Feb 06 11:21:05 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\JAN0802\31253.D
Operator : 209 GALOS
Acquired : 9 Jan 2002 10:34 pm using AcqMethod GCMS0103
Instrument : GC/MS 209
Sample Name: gcms scan 12/24
Misc Info :
Vial Number: 10
Quant File :GCMS0103.RES (RTE Integrator)



31253.D GCMS0103.M Wed Feb 06 11:21:11 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\JAN0802\31253.D
Acq On : 9 Jan 2002 10:34 pm
Sample : gcms scan 12/24
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

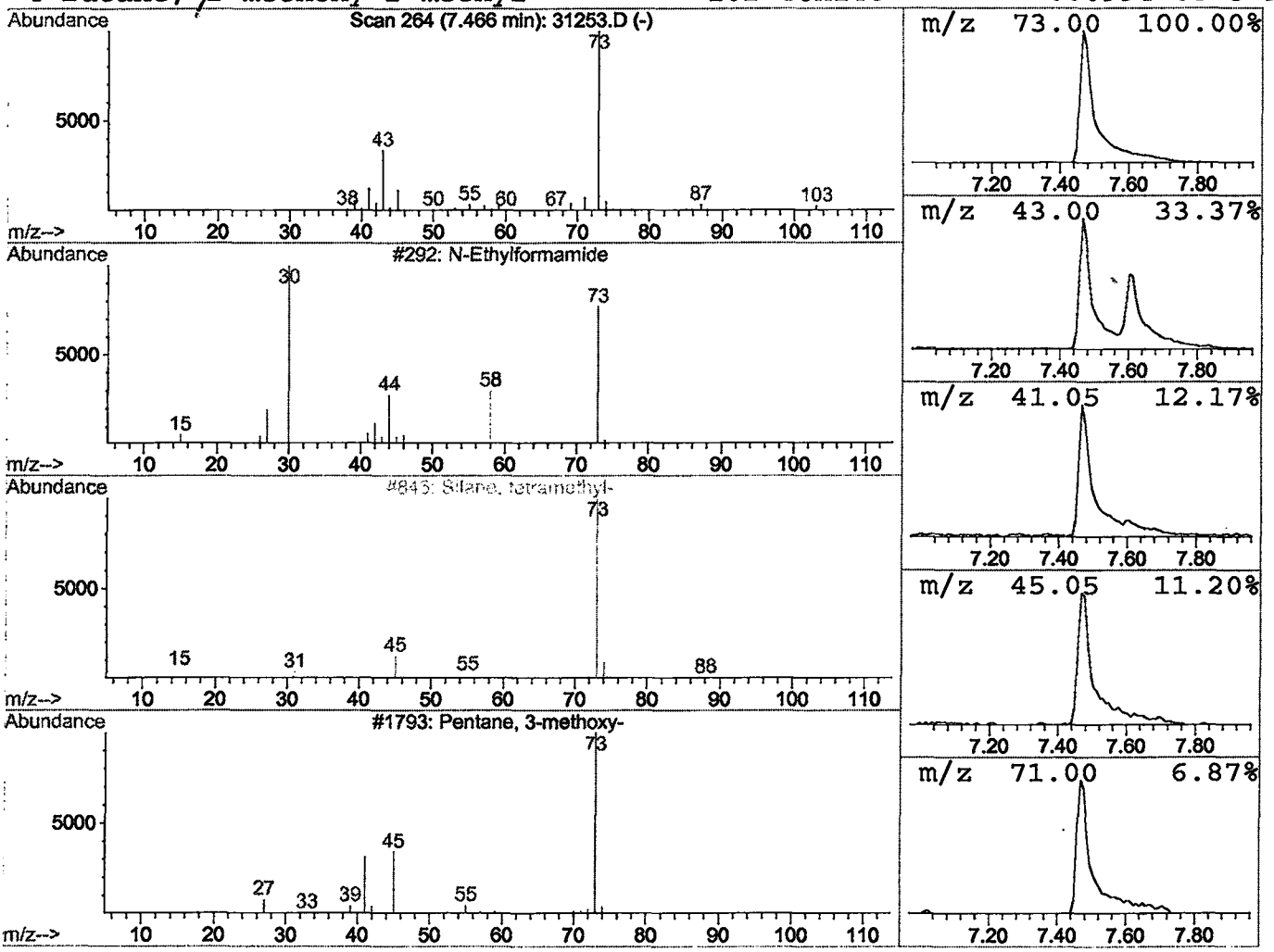
Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 N-Ethylformamide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.47	2.82 ug/l	783451	1,4-Dichlorobenzene-d4	11.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N-Ethylformamide	73	C3H7NO	000627-45-2	9
2	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
4	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\JAN0802\31253.D
Acq On : 9 Jan 2002 10:34 pm
Sample : gcms scan 12/24
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

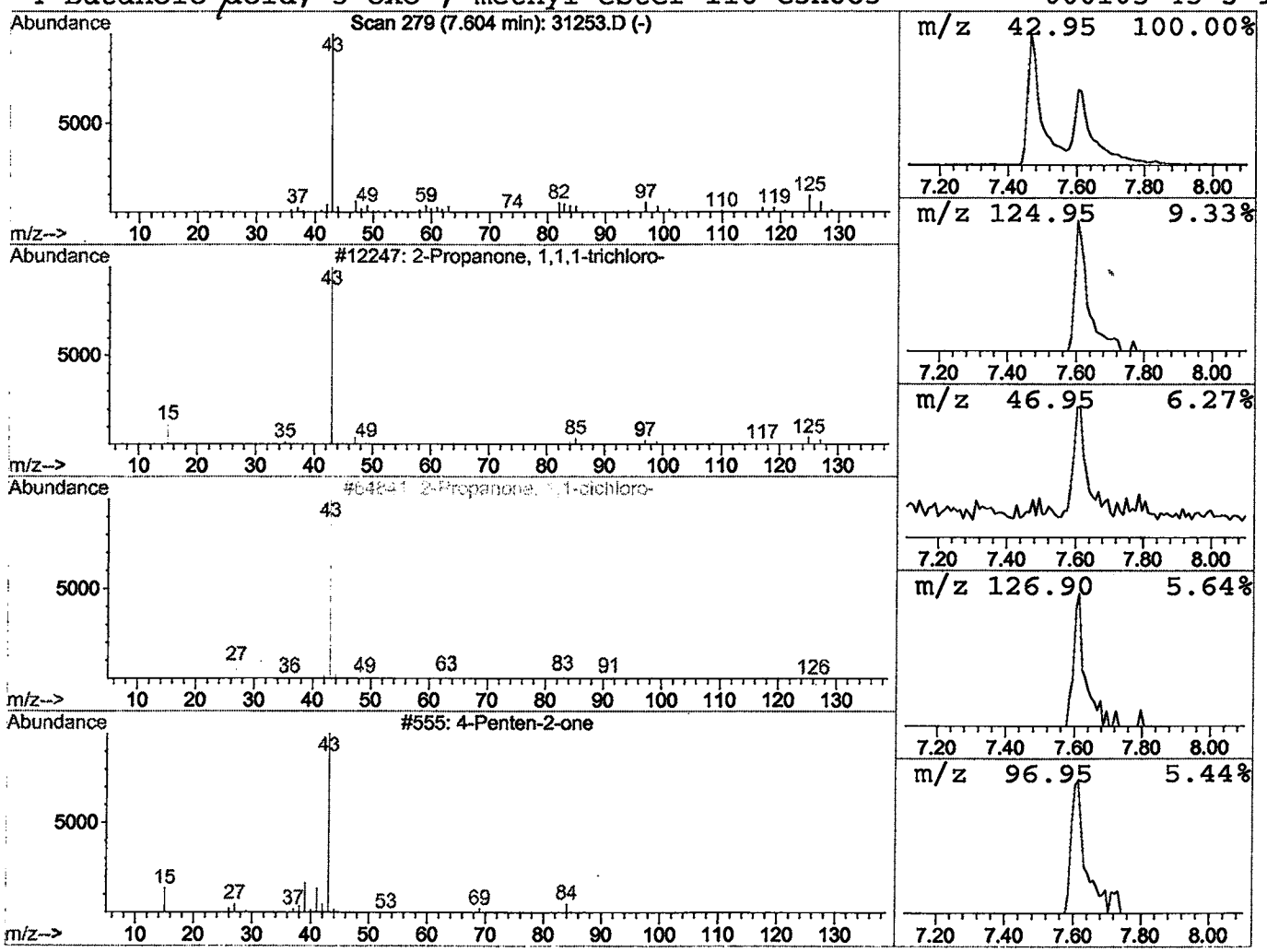
Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	1.01 ug/l	279836	1,4-Dichlorobenzene-d4	11.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanone, 1,1,1-trichloro-	160	C3H3Cl3O	000918-00-3	32
2	2-Propanone, 1,1-dichloro-	126	C3H4Cl2O	000513-88-2	12
3	4-Penten-2-one	84	C5H8O	013891-87-7	9
4	Butanoic acid, 3-oxo-, methyl ester	116	C5H8O3	000105-45-3	9



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\JAN0802\31253.D
Acq On : 9 Jan 2002 10:34 pm
Sample : gcms scan 12/24
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

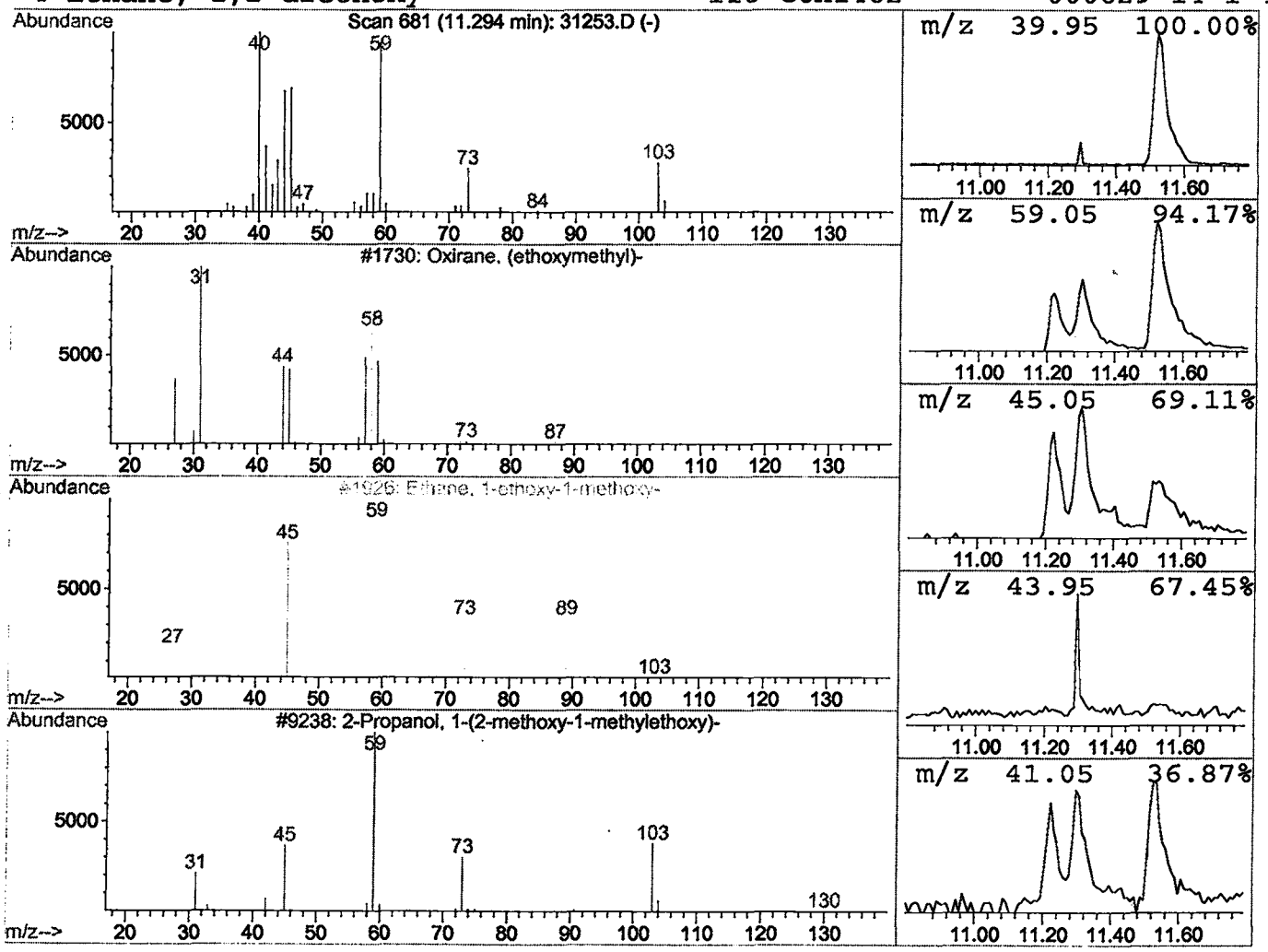
Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 3 Oxirane, (ethoxymethyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.29	0.31 ug/l	86325	1,4-Dichlorobenzene-d4	11.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, (ethoxymethyl)-	102	C5H10O2	004016-11-9	45
2	Ethane, 1-ethoxy-1-methoxy-	104	C5H12O2	010471-14-4	40
3	2-Propanol, 1-(2-methoxy-1-methylet	148	C7H16O3	020324-32-7	40
4	Ethane, 1,2-diethoxy-	118	C6H14O2	000629-14-1	40



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\JAN0802\31253.D
Acq On : 9 Jan 2002 10:34 pm
Sample : gcms scan 12/24
Misc :
MS Integration Params: LSCINT.P

Vial: 10
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

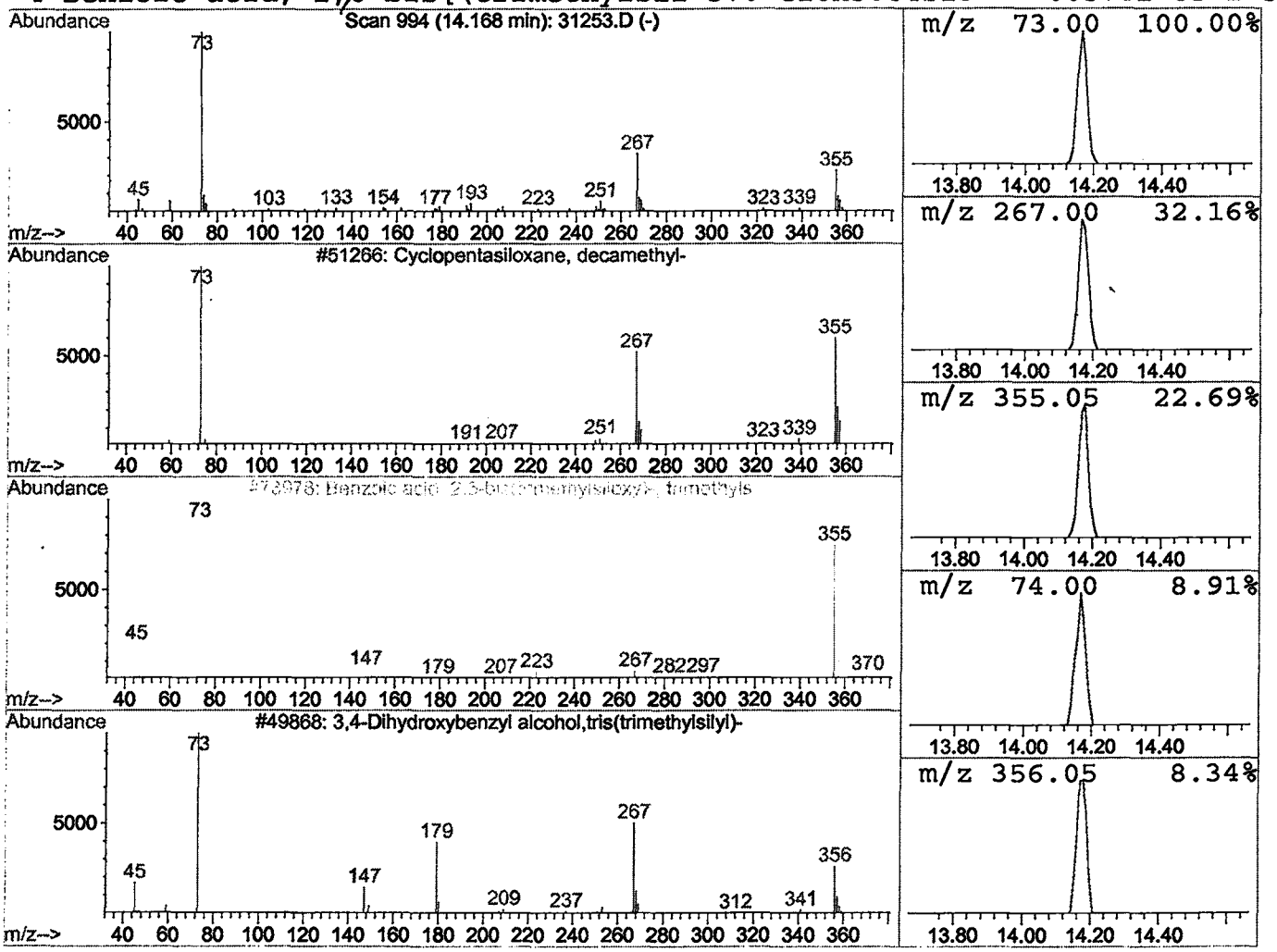
Quant Method : C:\HPCHEM\1\METHODS\GCMS0103.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 4 Cyclopentasiloxane, decamethyl Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.17	0.23 ug/l	79651	Naphthalene-d8	15.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	74
2	Benzoic acid, 2,5-bis(trimethylsilo	370	C16H30O4Si3	003618-20-0	47
3	3,4-Dihydroxybenzyl alcohol,tris(tr	356	C16H32O3Si3	068595-79-9	43
4	Benzoic acid, 2,6-bis[(trimethylsil	370	C16H30O4Si3	003782-85-2	33



Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\JAN0802\31253.D
Acq On : 9 Jan 2002 10:34 pm
Sample : gcms scan 12/24
Misc :
MS Integration Params: LSCINT.P

Vial: 10
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

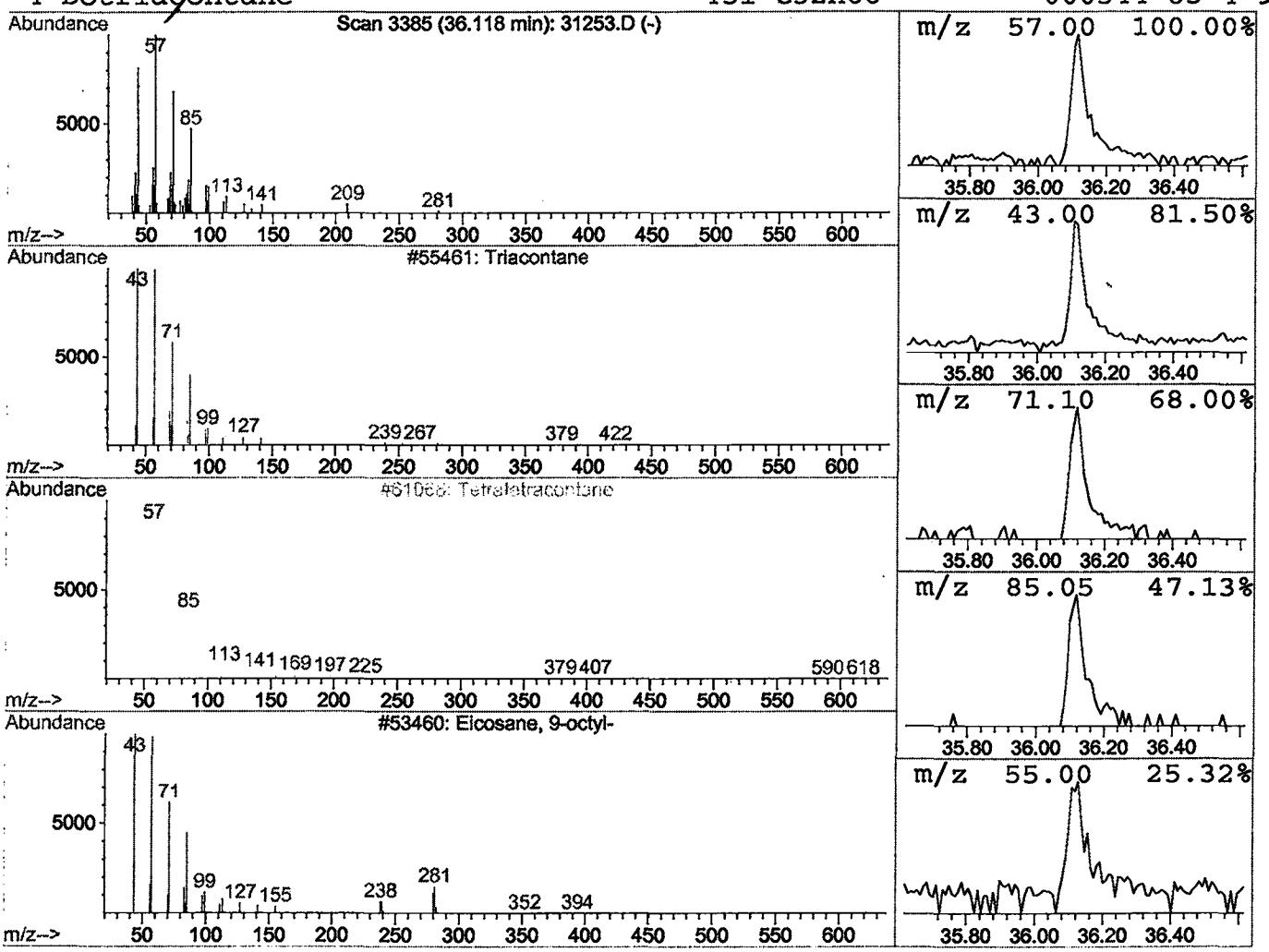
****

Peak Number 5 Triacontane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
36.12	0.46 ug/l	72312	Perylene-d12	36.91

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triacontane	422	C30H62	000638-68-6	91
2	Tetratetracontane	619	C44H90	007098-22-8	91
3	Eicosane, 9-octyl-	394	C28H58	013475-77-9	90
4	Dotriacontane	451	C32H66	000544-85-4	90

TV grad



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 9 Jan 2002 10:34 pm
Data File: C:\HPCHEM\1\DATA\JAN0802\31253.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
N-Ethylformamide	7.47	2.8	ug/l	783451	ISTD01	11.52	5557490	20.0
2-Propanone, 1,1,1-t	7.60	1.0	ug/l	279836	ISTD01	11.52	5557490	20.0
Oxirane, (ethoxymeth	11.29	0.3	ug/l	86325	ISTD01	11.52	5557490	20.0
Cyclopentasiloxane,	14.17	0.2	ug/l	79651	ISTD02	15.12	6919560	20.0
~~Triacontane~~	36.12	0.5	ug/l	72312	ISTD06	36.91	3177510	20.0

31253.D GCMS0103.M Wed Feb 06 11:21:41 2002

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\FEB1402\31253.D Vial: 29
Acq On : 15 Feb 2002 6:21 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 12:03 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	348149	20.00	ug/l	-0.01
10) Naphthalene-d8	15.98	136	1351619	20.00	ug/l	-0.01
17) Acenaphthene-d10	21.30	164	730488	20.00	ug/l	-0.01
29) Phenanthrene-d10	25.75	188	1063653	20.00	ug/l	-0.01
38) Chrysene-d12	33.82	240	736469	20.00	ug/l	-0.02
44) Perylene-d12	38.11	264	455034	20.00	ug/l	-0.01
System Monitoring Compounds						
37) 2,4-DDD	30.83	235	45424	4.25	ug/mL	0.33
Spiked Amount 5.000			Recovery = 85.00%			
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	0.00	128	0	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	217	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\31253.D Vial: 29
Acq On : 15 Feb 2002 6:21 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 12:03 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	25.74	178	324	N.D.	
34) Anthracene	25.74	178	324	N.D.	
35) Di-n-butylphthalate	27.72	149	6470	N.D.	
36) Fluoranthene	0.00	202	0	N.D.	
39) Pyrene	0.00	202	0	N.D.	
40) Butylbenzylphthalate	0.00	149	0	N.D.	
41) Benzo(a)anthracene	33.82	228	1873	N.D.	
42) Bis(2-ethylhexyl)phthalate	34.02	149	960	N.D.	
43) Chrysene	33.82	228	1873	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	1884	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.	

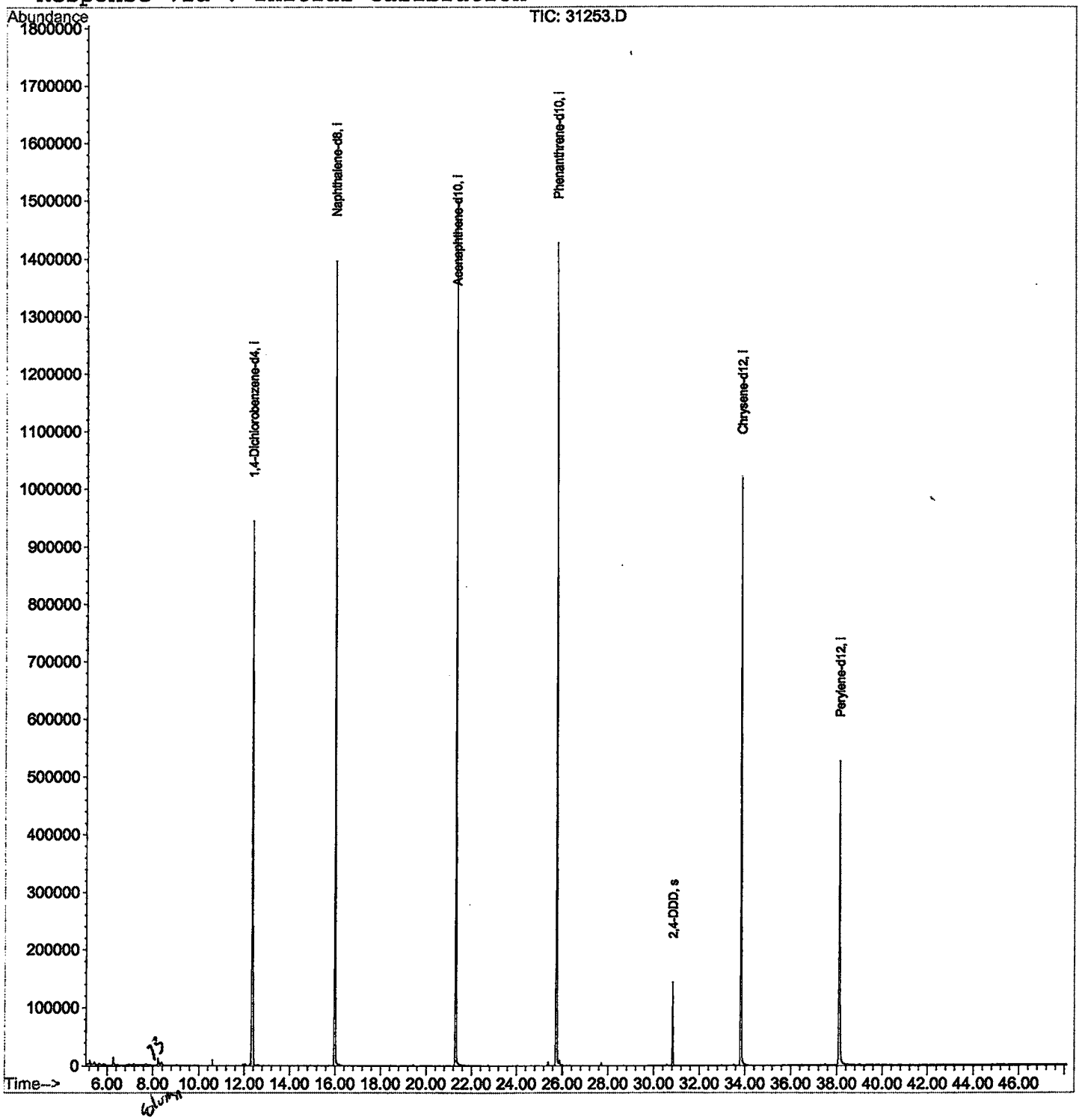
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\FEB1402\31253.D
Acq On : 15 Feb 2002 6:21 am
Sample : GC/MS SCAN 2/11
Misc :
MS Integration Params: GOOFY.P
Quant Time: Feb 15 12:03 2002

Vial: 29
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\31253.D Vial: 29
Acq On : 15 Feb 2002 6:21 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

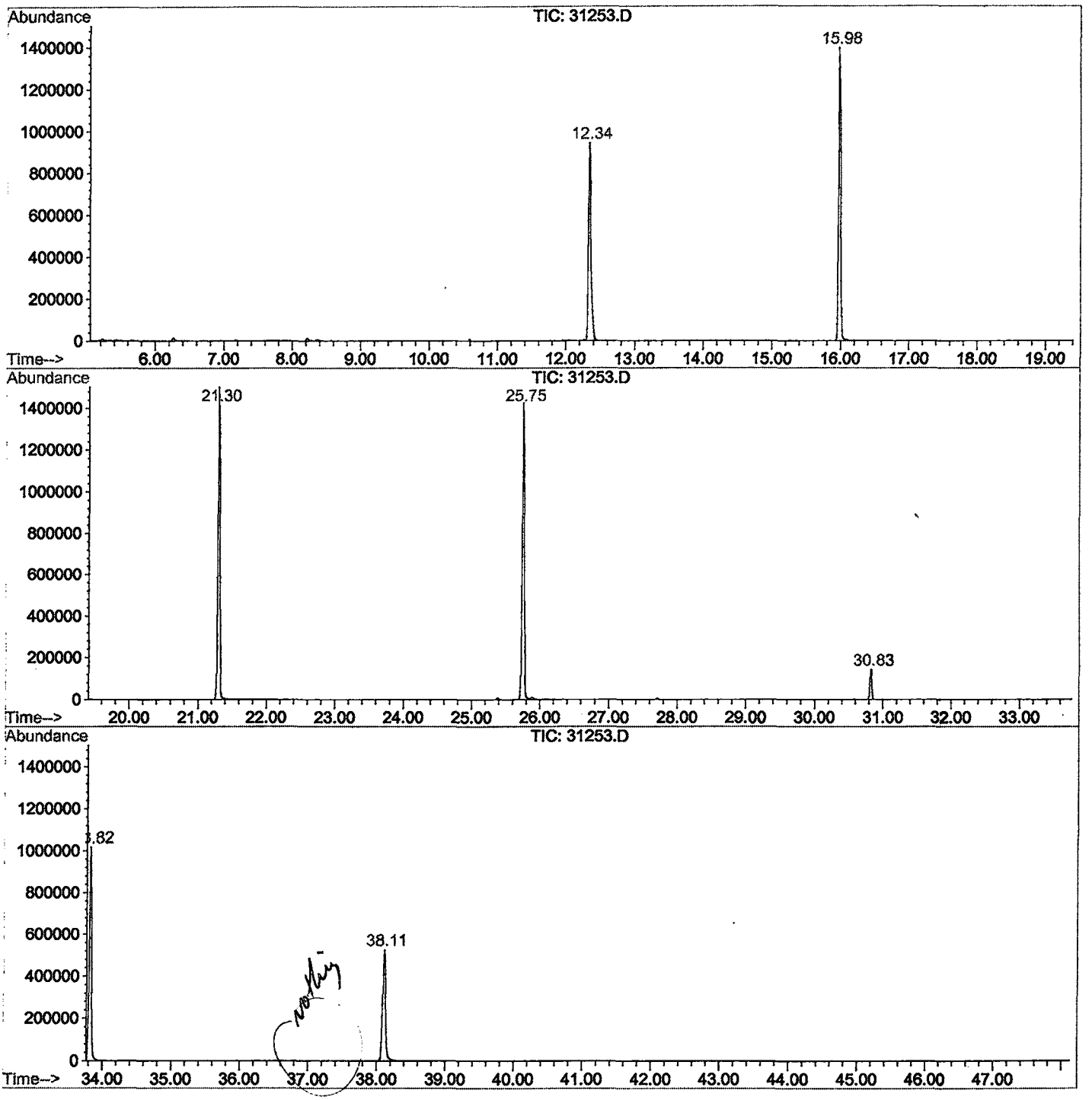
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	12.339	790	795	809	rVB	943311	2221354	73.49%	14.674%
2	15.984	1186	1192	1210	rBV	1396441	2813149	93.06%	18.583%
3	21.299	1765	1771	1780	rBV	1503593	3022848	100.00%	19.968%
4	25.752	2249	2256	2267	rBV2	1426760	2967151	98.16%	19.600%
5	30.829	2804	2809	2816	rVB	143108	280864	9.29%	1.855%
6	33.821	3128	3135	3151	rBV2	1017898	2264560	74.91%	14.959%
7	38.109	3593	3602	3624	rBV	524665	1568437	51.89%	10.361%

Sum of corrected areas: 15138363

31253.D GCMS0208.M Fri Feb 15 14:02:58 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\FEB1402\31253.D
Operator : 209 GALOS
Acquired : 15 Feb 2002 6:21 am using AcqMethod GCMS0208
Instrument : GC/MS 209
Sample Name: GC/MS SCAN 2/11
Misc Info :
Vial Number: 29
Quant File :GCMS0208.RES (RTE Integrator)



31253.D GCMS0208.M Fri Feb 15 14:03:03 2002

Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 15 Feb 2002 6:21 am
Data File: C:\HPCHEM\1\DATA\FEB1402\31253.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

31253.D GCMS0208.M Fri Feb 15 14:03:08 2002