User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 03/11/2002 Time: 16:08:51

Sample: AE02788
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 3/11/02 16:08 Ended: 3/11/02 16:08 Analyst: DLV
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		112		5.0

Comments for Sample AE02788 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 03-11-2002 Time: 16:09:01

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 below its acceptance range.

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\MAR0102\2788.D Vial: 22
Acq On : 2 Mar 2002 5:00 am Operator:
Sample : gc/ms scan 2/6 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Mar 5 10:41 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 27 11:30:15 2002
Response via : Initial Calibration
DataAcq Meth : GCMS0208

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	12.31	152	302025	20.00	ug/l	-0.04
10) Naphthalene-d8	15.96	136	1166931	20.00	ug/l	-0.04
17) Acenaphthene-d10	21.27	164	618040	20.00	ug/l	-0.04
29) Phenanthrene-d10	25.71	188	887788	20.00	ug/l	-0.05
38) Chrysene-d12	33.78	240	623430	20.00	ug/l	-0.06
44) Perylene-d12	38.05	264	423949	20.00	ug/l	-0.07
System Monitoring Compounds						
37) 2,4-DDD	30.79	235	51802	5.59	ug/mL	0.29
Spiked Amount 5.000			Recovery = 111.80%			
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	16.01	128	414	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	0.00	149	0	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

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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	0.00	178	0	N.D.	
34)	Anthracene	0.00	178	0	N.D.	
35)	Di-n-butylphthalate	27.68	149	13194	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	33.78	228	1486	N.D.	
42)	Bis(2-ethylhexyl)phthalate	33.99	149	2002	N.D.	
43)	Chrysene	33.78	228	1486	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	38.05	252	1436	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

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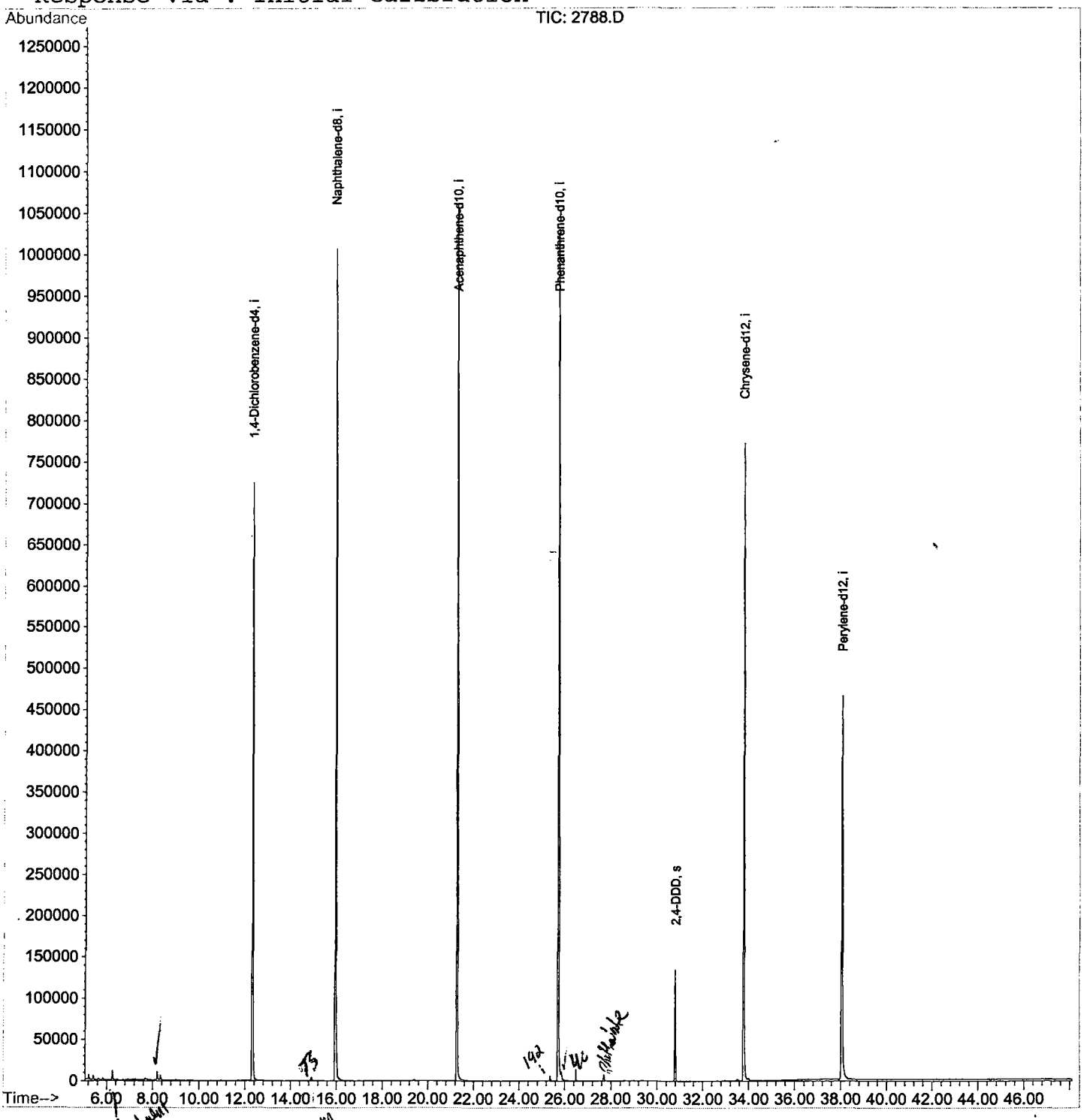
Acq On : 2 Mar 2002 5:00 am Operator:
Sample : gc/ms scan 2/6 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Mar 5 10:41 2002 Quant Results File: GCMS0208.RES

Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Wed Feb 27 11:30:15 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\MAR0102\2788.D Vial: 22
Acq On : 2 Mar 2002 5:00 am Operator:
Sample : gc/ms scan 2/6 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

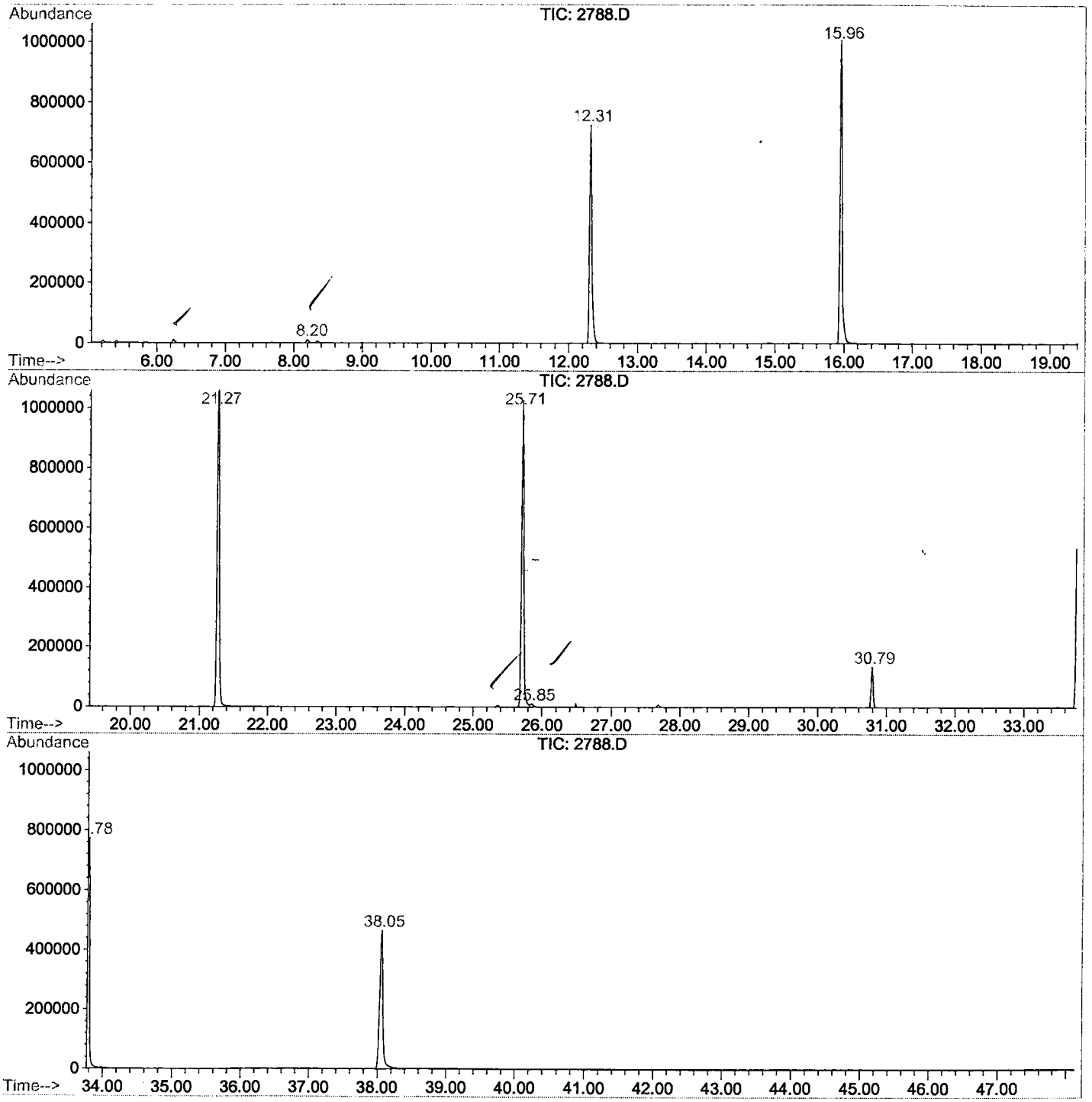
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	8.199	340	344	354	rBV	10155	25103	1.03%	0.202%
2	12.312	787	792	809	rBV	724252	1689196	69.17%	13.581%
3	15.956	1183	1189	1206	rBV	1006164	2261849	92.62%	18.185%
4	21.272	1761	1768	1781	rBV	1060672	2442147	100.00%	19.634%
5	25.715	2245	2252	2264	rBV2	1025336	2327044	95.29%	18.709%
6	25.853	2264	2267	2276	rVB3	9755	29557	1.21%	0.238%
7	30.792	2800	2805	2813	rBV	135182	271166	11.10%	2.180%
8	33.784	3123	3131	3150	rBV2	771779	1893885	77.55%	15.227%
9	38.053	3586	3596	3620	rBV2	464145	1498133	61.34%	12.045%

Sum of corrected areas: 12438080

2788.D GCMS0208.M Tue Mar 05 10:53:21 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\MAR0102\2788.D
Operator :
Acquired : 2 Mar 2002 5:00 am using AcqMethod GCMS0208
Instrument : GC/MS 209
Sample Name: gc/ms scan 2/6
Misc Info :
Vial Number: 22
Quant File :GCMS0208.RES (RTE Integrator)



2788.D GCMS0208.M Tue Mar 05 10:53:27 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\MAR0102\2788.D Vial: 22
Acq On : 2 Mar 2002 5:00 am Operator:
Sample : gc/ms scan 2/6 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

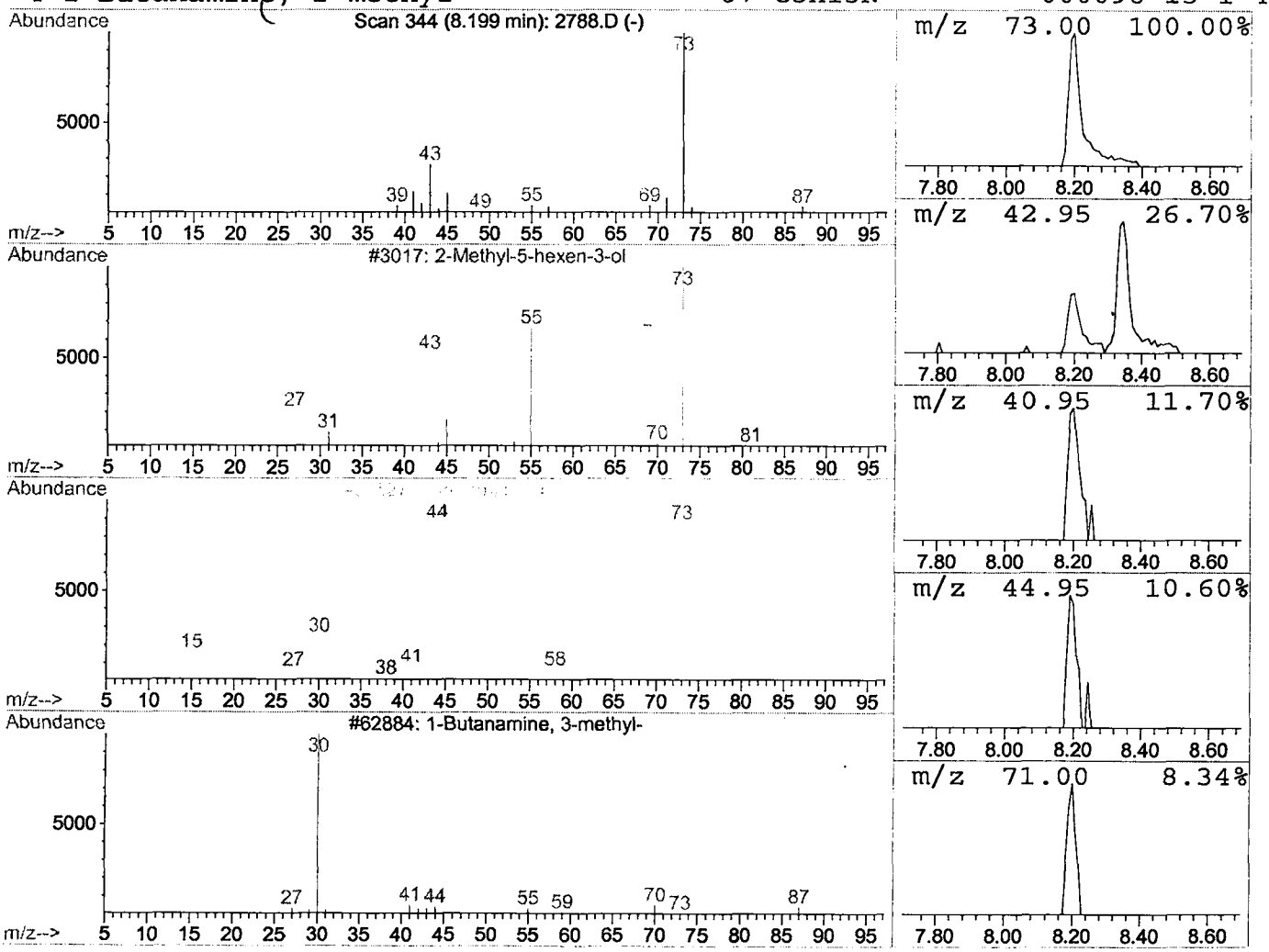
Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 2-Methyl-5-hexen-3-ol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.20	0.30 ug/l	25103	1,4-Dichlorobenzene-d4	12.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	9
2	Formamide, N,N-dimethyl-	73	C3H7NO	000068-12-2	5
3	1-Butanamine, 3-methyl-	87	C5H13N	000107-85-7	4
4	1-Butanamine, 2-methyl-	87	C5H13N	000096-15-1	4



Data File : C:\HPCHEM\1\DATA\MAR0102\2788.D Vial: 22
Acq On : 2 Mar 2002 5:00 am Operator:
Sample : gc/ms scan 2/6 Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: LSCINT.P

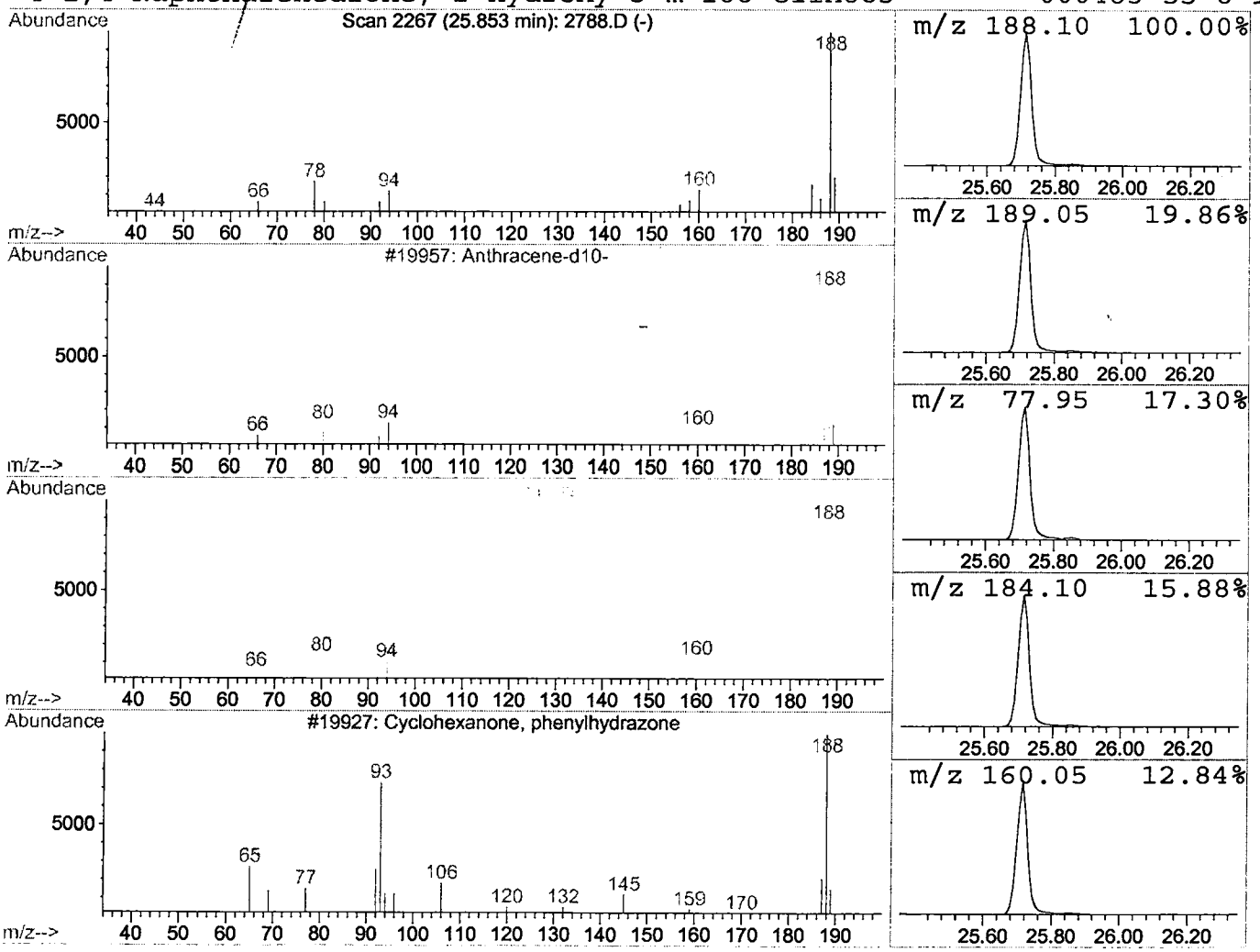
Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 2 Anthracene-d10- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.85	0.25 ug/l	29557	Phenanthrene-d10	25.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene-d10-	188	C14D10	001719-06-8	86
2	Phenanthrene-d10	188	C14D10	001517-22-2	86
3	Cyclohexanone, phenylhydrazone	188	C12H16N2	000946-82-7	9
4	1,4-Naphthalenedione, 2-hydroxy-3-m	188	C11H8O3	000483-55-6	9



Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 2 Mar 2002 5:00 am
Data File: C:\HPCHEM\1\DATA\MAR0102\2788.D
Name: gc/ms scan 2/6
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
2-Methyl-5-hexen-3-o	8.20	0.3	ug/l	25103	ISTD01	12.31	1689200	20.0
Anthracene-d10-	25.85	0.3	ug/l	29557	ISTD04	25.72	2327040	20.0

2788.D GCMS0208.M Tue Mar 05 10:53:41 2002

Library Searched : C:\DATABASE\nbs75k.l
Quality : 42
ID : Methanethioamide, N,N-dimethyl-