User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 02/15/2002 Time: 09:29:53

Sample: AE02973
Analysis: $GCMS GCMS Scan For Unknowns
Units: ug
Started: 2/15/02 09:29 Ended: 2/15/02 09:29 Analyst: DLV
Page: 1

	Component Name	Vial	Result	Qualifier	MDL
1	Other unknown compounds		.		
2	Surr_2,4-DDD		80		5.0

Comments for Sample AE02973 - Analysis $GCMS

Westchester Dept. of Labs & Research
User: Vinci, David L
Date: 02-15-2002 Time: 09:30:01

The GCMS scan, from 35 to 550 amu, does not reveal the presence of unusual compounds.
The recoveries for 7 of the 43 compounds in the LCS Standard were below their acceptance ranges.

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\FEB0802\2973.D Vial: 37
Acq On : 9 Feb 2002 8:17 pm Operator: 209 GALOS
Sample : GC/MS SCAN 2/7FB Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Feb 13 10:51 2002 Quant Results File: GCMS0208.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Fri Feb 08 15:29:22 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	419654	20.00	ug/l	-0.01
10) Naphthalene-d8	15.99	136	1608990	20.00	ug/l	-0.01
17) Acenaphthene-d10	21.30	164	877914	20.00	ug/l	-0.01
29) Phenanthrene-d10	25.76	188	1252962	20.00	ug/l	-0.01
38) Chrysene-d12	33.83	240	836354	20.00	ug/l	-0.02
44) Perylene-d12	38.11	264	544403	20.00	ug/l	-0.01
System Monitoring Compounds						
37) 2,4-DDD	30.83	235	62366	4.00	ug/mL	0.33
Spiked Amount 5.000			Recovery = 80.00%			
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	16.05	128	319	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	20.64	163	204	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	553	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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Title : 209 625/TCLP calibration table
Last Update : Fri Feb 08 15:29:22 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	25.82	178	366	N.D.	
34)	Anthracene	25.95	178	323	N.D.	
35)	Di-n-butylphthalate	27.71	149	3458	N.D.	
36)	Fluoranthene	29.44	202	703	N.D.	
39)	Pyrene	30.13	202	861	N.D.	
40)	Butylbenzylphthalate	0.00	149	0	N.D.	
41)	Benzo(a)anthracene	33.83	228	2475	N.D.	
42)	Bis(2-ethylhexyl)phthalate	34.03	149	3352	N.D.	
43)	Chrysene	33.89	228	273	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	2499	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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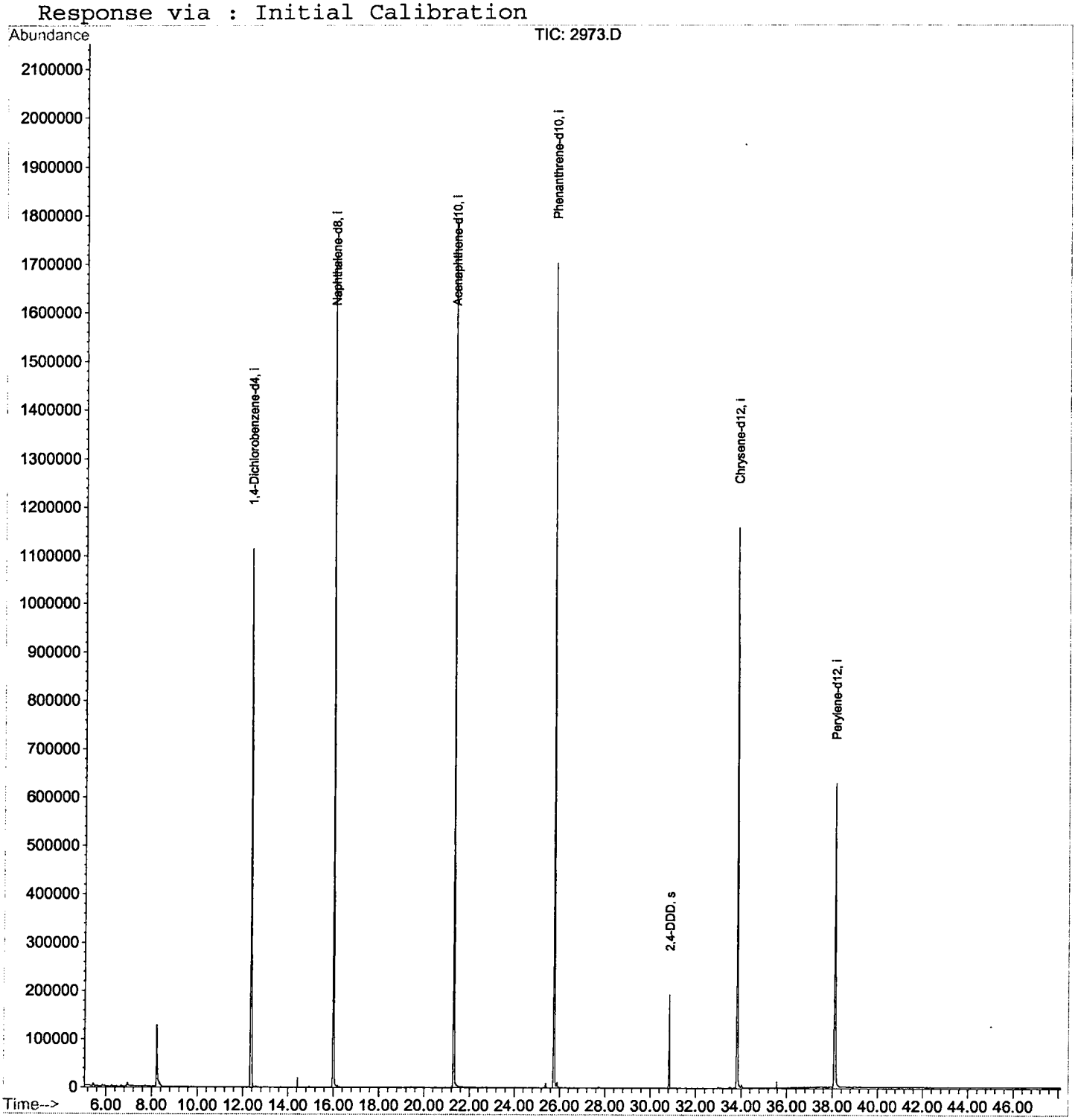
Acq On : 9 Feb 2002 8:17 pm Operator: 209 GALOS
Sample : GC/MS SCAN 2/7FB Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Feb 13 10:51 2002 Quant Results File: GCMS0208.RES

Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Last Update : Fri Feb 08 15:29:22 2002
Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\FEB0802\2973.D Vial: 37
Acq On : 9 Feb 2002 8:17 pm Operator: 209 GALOS
Sample : GC/MS SCAN 2/7FB 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

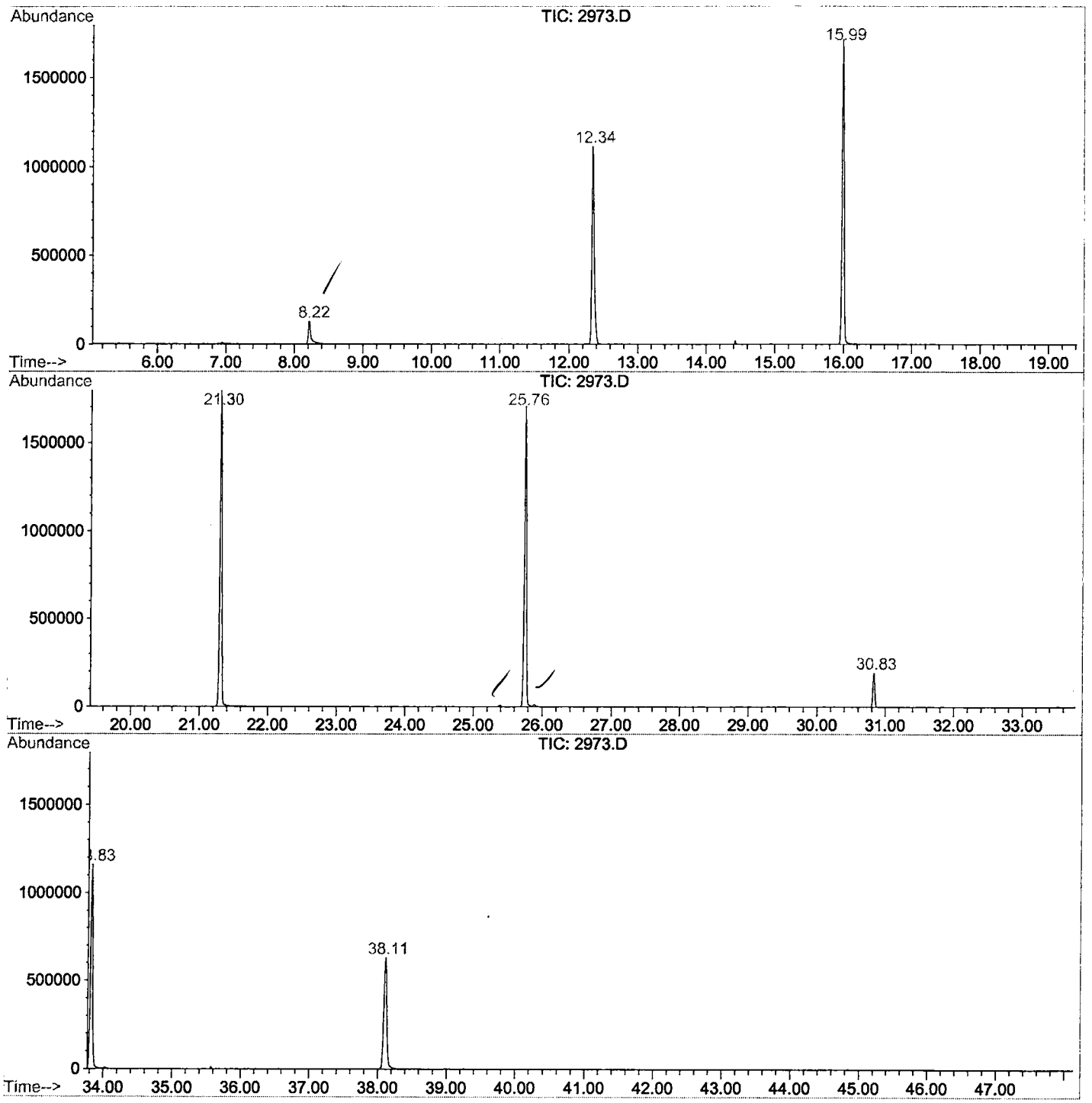
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	peak area	peak % max.	% of total
1	8.222	341	346	361	rBV	127057	351946	9.62%	1.926%
2	12.344	789	795	810	rBV	1113332	2649879	72.46%	14.503%
3	15.988	1186	1192	1208	rBV	1713730	3328642	91.02%	18.218%
4	21.304	1764	1771	1791	rBV	1792689	3656923	100.00%	20.014%
5	25.756	2249	2256	2266	rBV	1701734	3458310	94.57%	18.927%
6	30.833	2804	2809	2815	rVB	193428	383938	10.50%	2.101%
7	33.826	3128	3135	3153	rBV	1157976	2549303	69.71%	13.952%
8	38.113	3593	3602	3622	rBV	627858	1892500	51.75%	10.358%

Sum of corrected areas: 18271441

2973.D GCMS0208.M Wed Feb 13 11:12:40 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\FEB0802\2973.D
Operator : 209 GALOS
Acquired : 9 Feb 2002 8:17 pm using AcqMethod GCMS0208
Instrument : GC/MS 209
Sample Name: GC/MS SCAN 2/7FB
Misc Info :
Vial Number: 37
Quant File :GCMS0208.RES (RTE Integrator)



2973.D GCMS0208.M Wed Feb 13 11:12:46 2002

Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\FEB0802\2973.D
Acq On : 9 Feb 2002 8:17 pm
Sample : GC/MS SCAN 2/7FB
Misc :
MS Integration Params: LSCINT.P

Vial: 37
Operator: 209 GALOS
Inst : GC/MS 209
Multiplr: 1.00

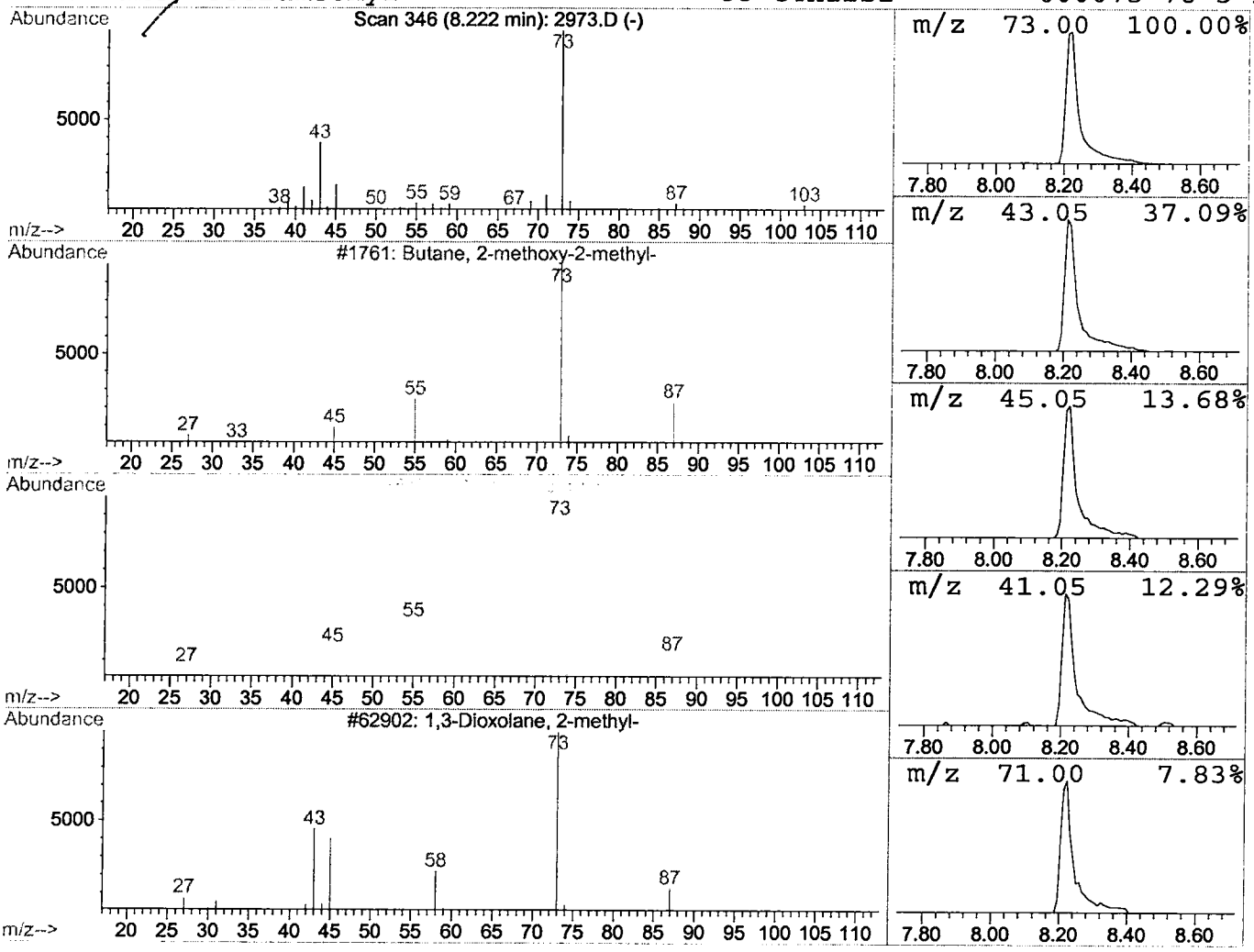
Quant Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
Title : 209 625/TCLP calibration table
Library : C:\DATABASE\NBS75K.L

****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.22	2.66 ug/l	351946	1,4-Dichlorobenzene-d4	12.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	36
2	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	28
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



Tentatively Identified Compound (LSC) summary

Operator ID: 209 GALOS Date Acquired: 9 Feb 2002 8:17 pm
Data File: C:\HPCHEM\1\DATA\FEB0802\2973.D
Name: GC/MS SCAN 2/7FB
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
Butane, 2 methoxy-2-	8.22	2.7	ug/l	351946	ISTD01	12.34	2649880	20.0

2973.D GCMS0208.M Wed Feb 13 11:12:55 2002