

Data File : C:\HPCHEM\1\DATA\MAY2302\12093.D

Vial: 8

Acq On : 23 May 2002 7:01 pm

Operator: Morgan

Sample : 5/21 bn

Inst : 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: May 31 15:47 2002

Quant Results File: GCMS0520.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0520.M (RTE Integrator)

Title : 209 625/TCLP calibration table

Last Update : Fri May 31 15:47:14 2002

Response via : Initial Calibration

DataAcq Meth : GCMS0520

original

to be re-ent

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	12.14	152	461300	40.00	ug/l	0.00
10) Naphthalene-d8	15.77	136	1678368	40.00	ug/l	-0.02
17) Acenaphthene-d10	21.09	164	914468	40.00	ug/l	0.00
30) Phenanthrene-d10	25.53	188	1198920	40.00	ug/l	0.00
38) Chrysene-d12	33.59	240	666071	40.00	ug/l	-0.01
44) Perylene-d12	37.81	264	457862	40.00	ug/l	0.00

System Monitoring Compounds

28) TCMX	23.23	209	127109	27.66	ug/L	0.00
Spiked Amount	40.000		Recovery		69.15%	

Target Compounds

					Qvalue
2) N-nitrosodimethylamine	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-Nitrosodi-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	22.58	149	239	N.D.	
26) 4-Chlorophenylphenylether	23.23	204	454	N.D.	
27) Fluorene	0.00	166	0	N.D.	
29) Azobenzene	0.00	77	0	N.D.	
31) N-Nitrosodiphenylamine	23.24	168	345	N.D.	
32) 4-Bromophenylphenylether	0.00	248	0	N.D.	
33) Hexachlorobenzene	0.00	284	0	N.D.	
34) Phenanthrene	25.53	178	114	N.D.	

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

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Quantitation Report

(QT Reviewed)

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Misc :

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MS Integration Params: GOOFY.P

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DataAcq Meth : GCMS0520

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
35) Anthracene	25.60	178	108		N.D.	
36) Di-n-butylphthalate	27.52	149	1821		N.D.	
37) 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	33.60	228	2050		N.D.	
42) Bis(2-ethylhexyl)phthalate	33.82	149	1461		N.D.	
43) Chrysene	33.66	228	818		N.D.	
45) Di-n-Octylphthalate	35.56	149	315		N.D.	
46) Benzo(b)fluoranthene	36.65	252	1147		N.D.	
47) Benzo(k)fluoranthene	36.71	252	1289		N.D.	
48) Benzo(a)pyrene	37.61	252	423		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

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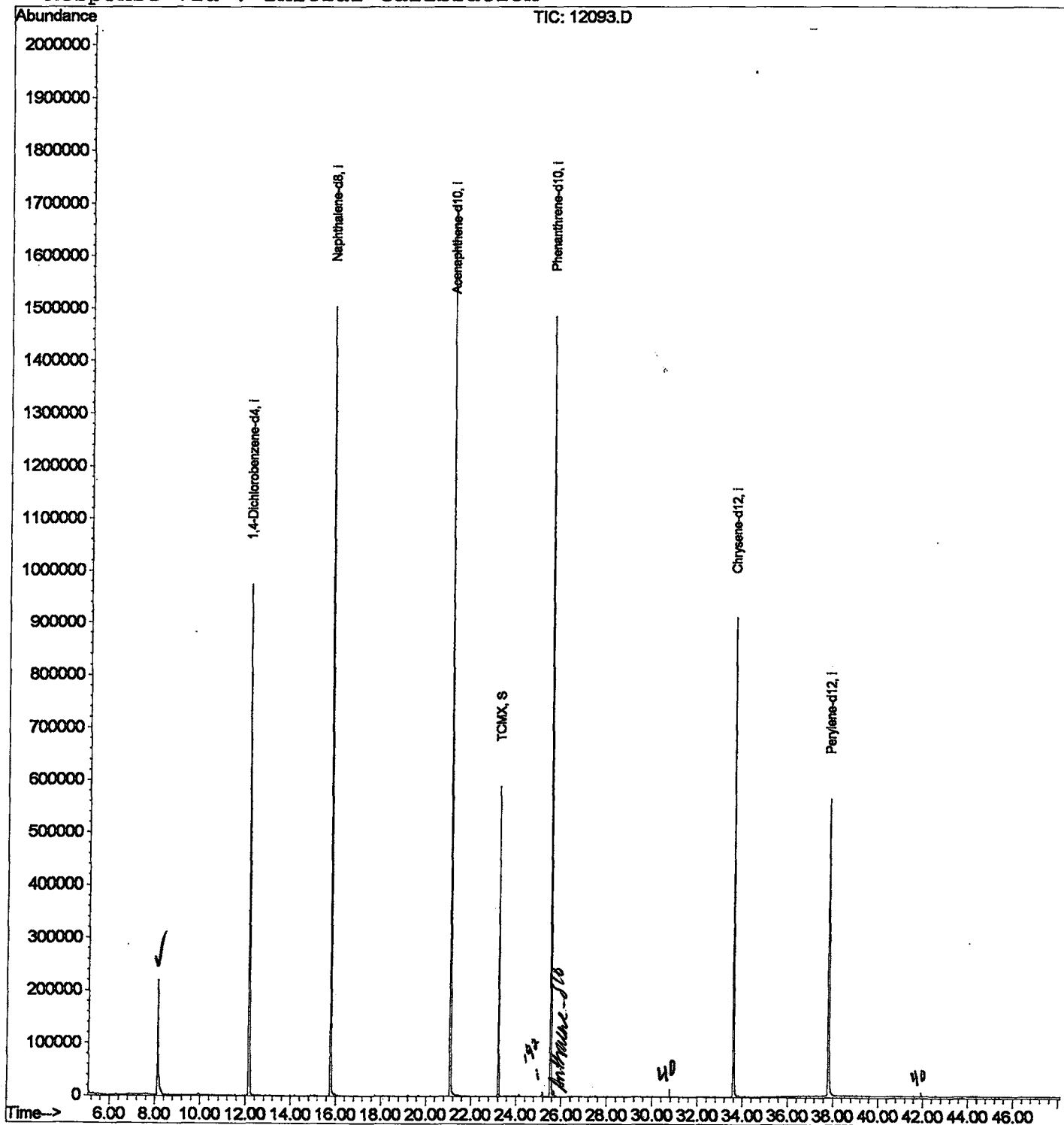
Quantitation Report

Data File : C:\HPCHEM\1\DATA\MAY2302\12093.D
 Acq On : 23 May 2002 7:01 pm
 Sample : 5/21 bn
 Misc :
 MS Integration Params: GOOFY.P
 Quant Time: May 31 15:47 2002

Vial: 8
 Operator: Morgan
 Inst : 209
 Multiplr: 1.00

Quant Results File: GCMS0520.RES

Method : C:\HPCHEM\1\METHODS\GCMS0520.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Last Update : Fri May 31 15:47:14 2002
 Response via : Initial Calibration



LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\MAY2302\12093.D

Vial: 8

Acq On : 23 May 2002 7:01 pm

Operator: Morgan

Sample : 5/21 bn

Inst : 209

Misc :

Multiplr: 1.00

MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\GCMS0520.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	8.118	331	335	366	rBV	217395	579925	16.74%	3.338%
2	12.139	768	773	788	rBV	972269	2326668	67.14%	13.392%
3	15.774	1163	1169	1187	rBV	1502885	3061593	88.35%	17.623%
4	21.090	1741	1748	1768	rBV	1696316	3465142	100.00%	19.946%
5	23.229	1973	1981	1991	rBV	588378	1190411	34.35%	6.852%
6	25.533	2225	2232	2243	rBV2	1485277	3106091	89.64%	17.879%
7	33.593	3103	3110	3130	rBV2	910547	2023828	58.41%	11.649%
8	37.807	3560	3569	3591	rBV2	563183	1619326	46.73%	9.321%

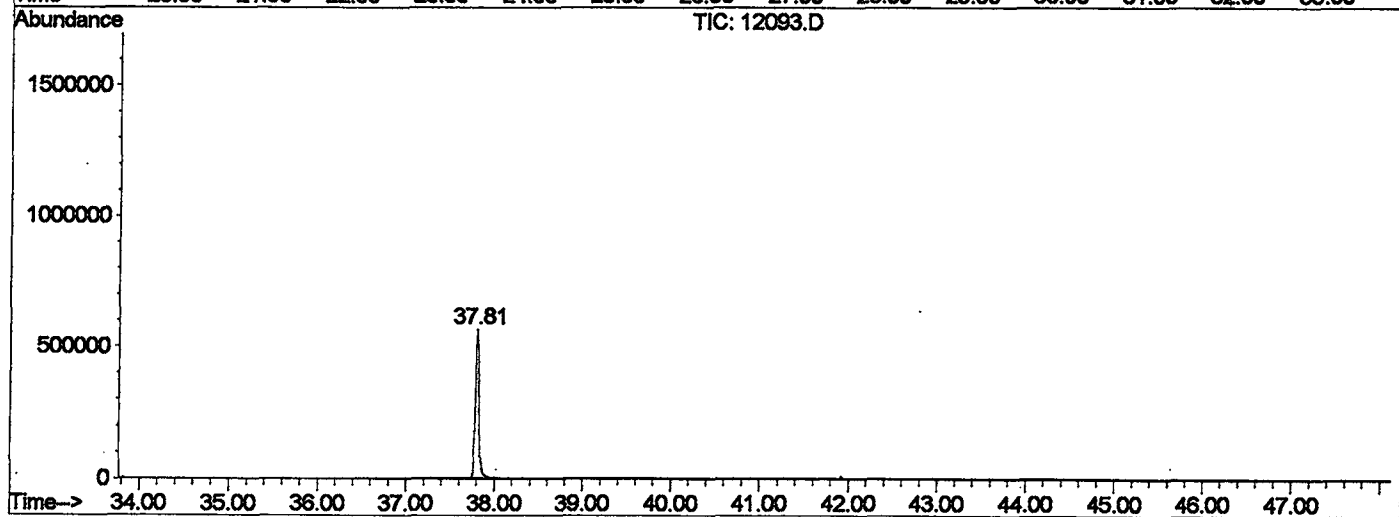
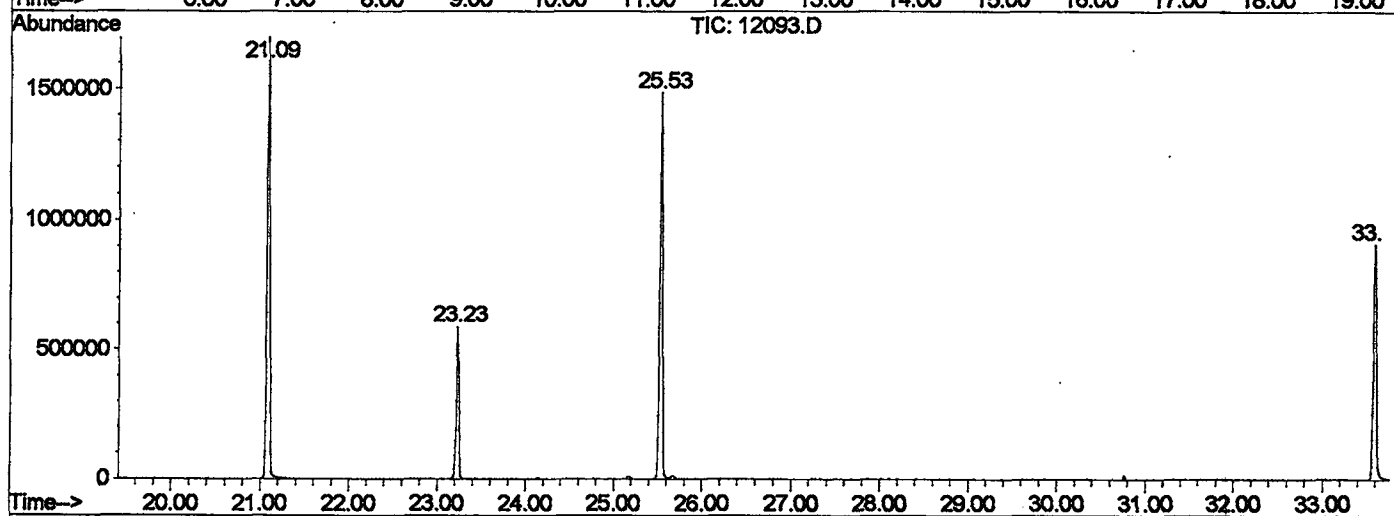
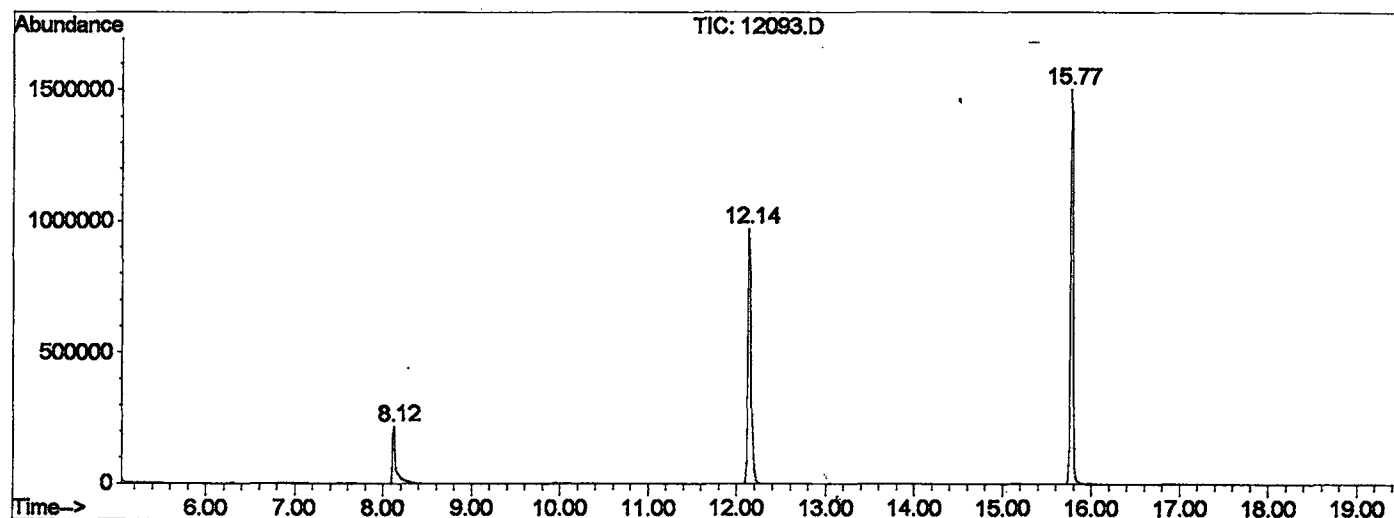
Sum of corrected areas: 17372984

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Fri May 31 15:46:15 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\MAY2302\12093.D
 Operator : Morgan
 Acquired : 23 May 2002 7:01 pm using AcqMethod GCMS0520
 Instrument : 209
 Sample Name: 5/21 bn
 Misc Info :
 Vial Number: 8
 Quant File :GCMS0520.RES (RTE Integrator)



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Library Search Compound Report

Data File : C:\HPCHEM\1\DATA\MAY2302\12093.D

Acq On : 23 May 2002 7:01 pm

Sample : 5/21 bn

Misc :

MS Integration Params: LSCINT.P

Vial: 8

Operator: Morgan

Inst : 209

Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\GCMS0520.M (RTE Integrator)

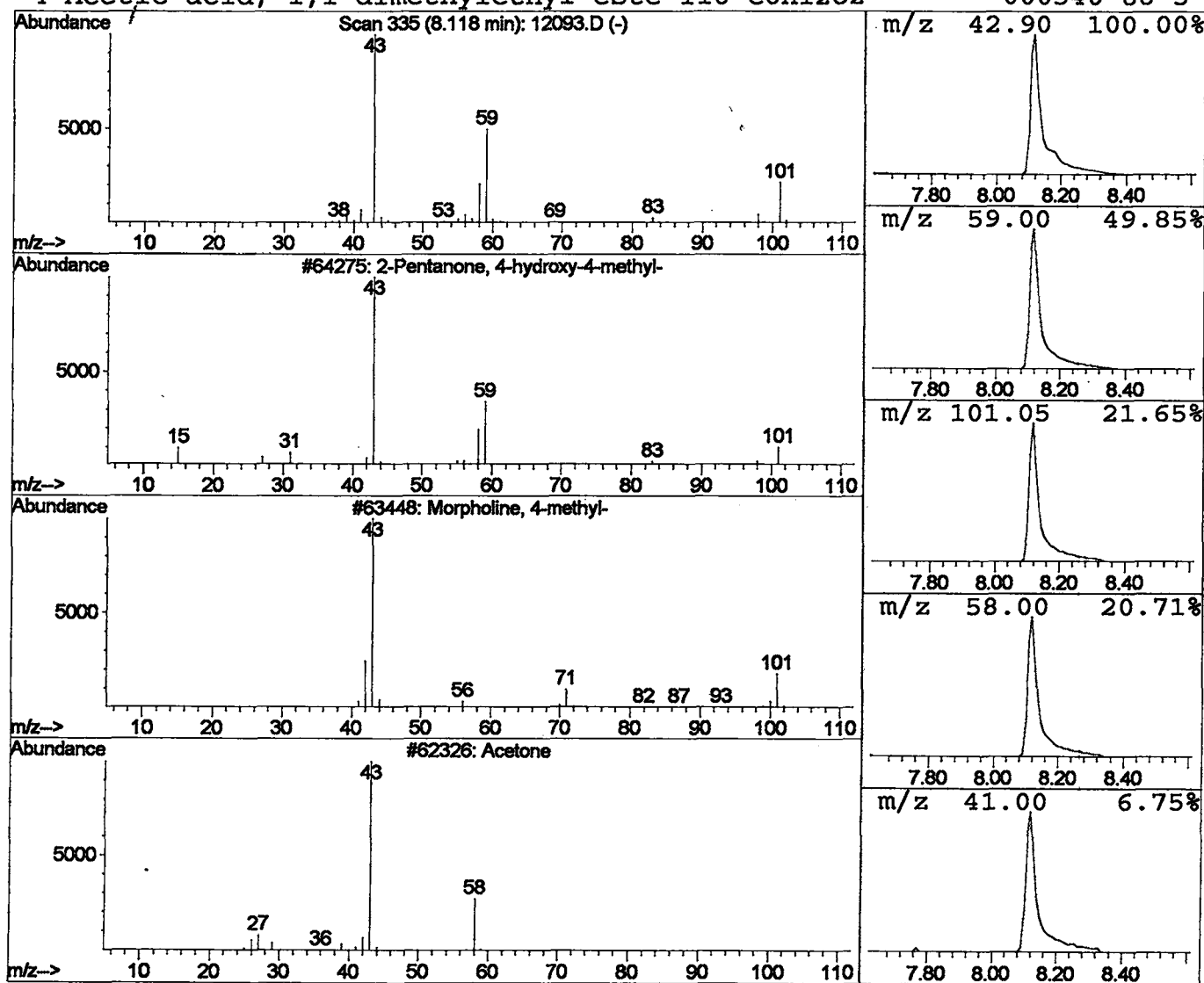
Title : 209 625/TCLP calibration table

Library : C:\DATABASE\NBS75K.L

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.12	9.97 ug/l	579925	1,4-Dichlorobenzene-d4	12.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
3			Acetone	58	C3H6O	000067-64-1	9
4			Acetic acid, 1,1-dimethylethyl este	116	C6H12O2	000540-88-5	9



Tentatively Identified Compound (LSC) summary

Operator ID: Morgan Date Acquired: 23 May 2002 7:01 pm
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Misc:
Method: C:\HPCHEM\1\METHODS\GCMS0520.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-Pentanone, 4-hydro	8.12	10.0	ug/l	579925	ISTD01	12.14	2326670	40.0

12093.D GCMS0520.M Fri May 31 15:46:17 2002