

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\APR1802\7108.D
 Acq On : 19 Apr 2002 3:38 pm
 Sample : GC/MS SCAN 3/26 FB
 Misc :
 MS Integration Params: GOOFY.P
 Quant Time: Apr 22 8:29 2002

Vial: 27
 Operator:
 Inst : 209
 Multiplr: 1.00

Quant Results File: GCMS0419.RES

Quant Method : C:\HPCHEM\1\METHODS\GCMS0419.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Last Update : Fri Apr 19 14:51:24 2002
 Response via : Initial Calibration
 DataAcq Meth : GCMS0412

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	12.22	152	674254	20.00	ug/l	-0.02
10) Naphthalene-d8	15.84	136	2468936	20.00	ug/l	-0.03
17) Acenaphthene-d10	21.15	164	1382141	20.00	ug/l	-0.02
30) Phenanthrene-d10	25.59	188	2030352	20.00	ug/l	-0.02
39) Chrysene-d12	33.65	240	1348445	20.00	ug/l	-0.02
45) Perylene-d12	37.88	264	896171	20.00	ug/l	-0.04

System Monitoring Compounds

28) TCMX	23.29	209	30671	2.66	ug/L	-0.04
Spiked Amount	4.000		Recovery	=	66.50%	
38) 2,4-DDD	0.00	235	0	0.00	ug/mL	
Spiked Amount	5.000		Recovery	=	0.00%	

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	21.53	165	256	N.D.
25) Diethylphthalate	22.63	149	733	N.D.
26) 4-Chlorophenyl-phenylether	0.00	204	0	N.D.
27) Fluorene	0.00	166	0	N.D.
29) Azobenzene/ 1,2-Diphenylhyd	0.00	77	0	N.D.
31) N-Nitrosodiphenylamine	0.00	168	0	N.D.
32) 4-Bromophenyl-phenylether	0.00	248	0	N.D.

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

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Response via : Initial Calibration

DataAcq Meth : GCMS0412

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
33) Hexachlorobenzene	0.00	284	0	N.D.		
34) Phenanthrene	25.60	178	571	N.D.		
35) Anthracene	25.60	178	571	N.D.		
36) Di-n-butylphthalate	27.56	149	2275	N.D.		
37) Fluoranthene	0.00	202	0	N.D.		
40) Pyrene	0.00	202	0	N.D.		
41) Butylbenzylphthalate	0.00	149	0	N.D.		
42) Benzo(a)anthracene	33.65	228	4274	N.D.		
43) Bis(2-ethylhexyl)phthalate	33.86	149	40975	1.21 ug/l	#	98
44) Chrysene	33.65	228	3774	N.D.		
46) Di-n-Octylphthalate	0.00	149	0	N.D.		
47) Benzo(b)fluoranthene	36.69	252	340	N.D.		
48) Benzo(k)fluoranthene	36.78	252	205	N.D.		
49) Benzo(a)pyrene	37.88	252	3782	N.D.		
50) Indeno(1,2,3-cd)pyrene	0.00	276	0	N.D.		
51) Dibenz(a,h)anthracene	0.00	278	0	N.D.		
52) Benzo(g,h,i)perylene	0.00	276	0	N.D.		

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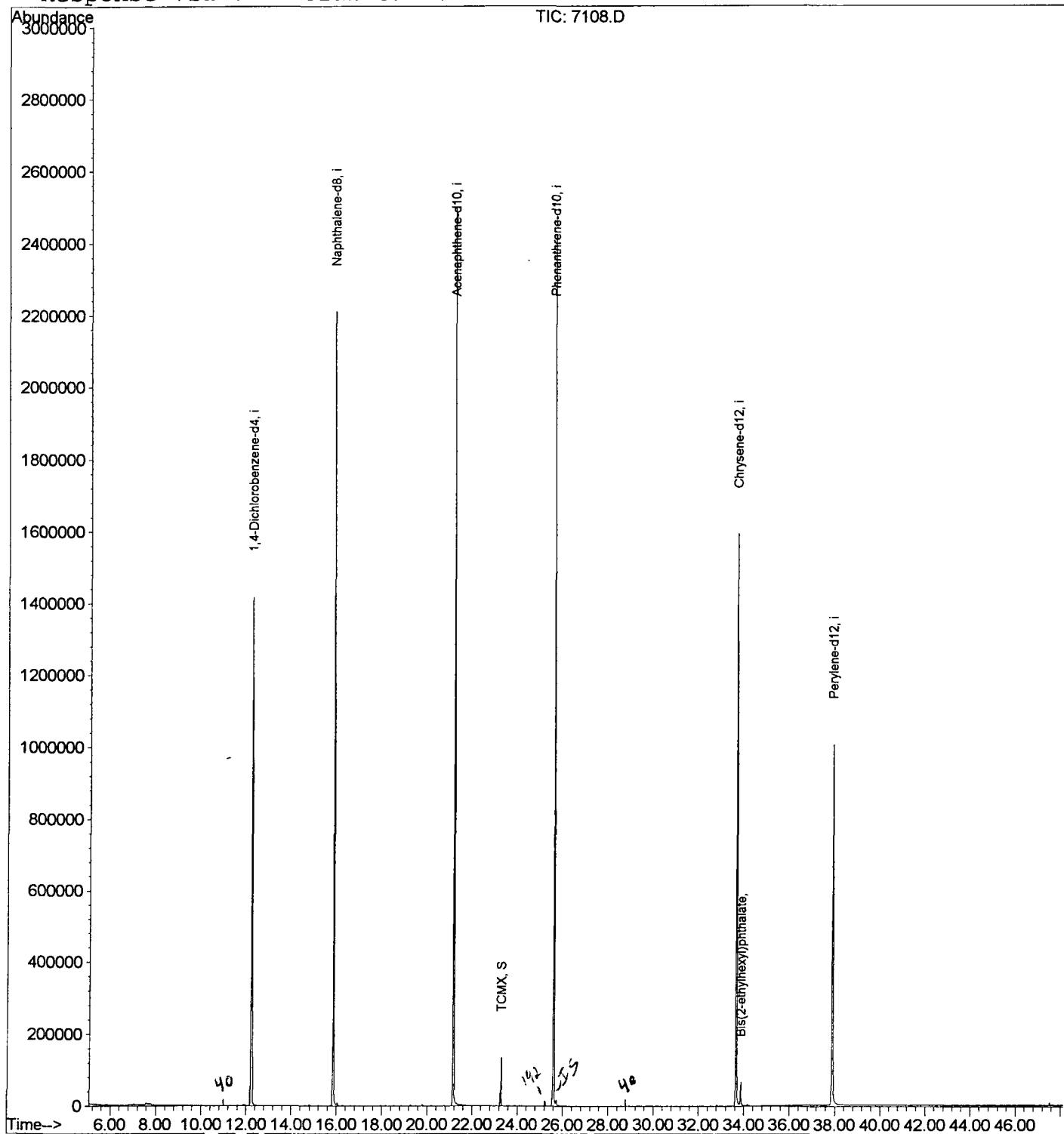
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Quant Time: Apr 22 8:29 2002

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Title : 209 625/TCLP calibration table
Last Update : Fri Apr 19 14:51:24 2002
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LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\APR1802\7108.D

Vial: 27

Acq On : 19 Apr 2002 3:38 pm

Operator:

Sample : GC/MS SCAN 3/26 FB

Inst : 209

Misc :

Multiplr: 1.00

MS Integration Params: LSCINT.P

Method : C:\HPCHEM\1\METHODS\GCMS0419.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	12.217	775	781	797	rVB	1414939	3598896	68.98%	13.946%
2	15.843	1170	1176	1193	rBV	2210530	4634300	88.83%	17.958%
3	21.149	1748	1754	1767	rBV	2501070	5216993	100.00%	20.216%
4	23.288	1981	1987	1992	rBV	135727	263338	5.05%	1.020%
5	25.592	2231	2238	2250	rBV2	2422021	5074987	97.28%	19.666%
6	33.653	3108	3116	3129	rBV2	1594238	3882093	74.41%	15.043%
7	33.864	3134	3139	3151	rVB	66298	151086	2.90%	0.585%
8	37.876	3568	3576	3588	rBV2	1002967	2984416	57.21%	11.565%

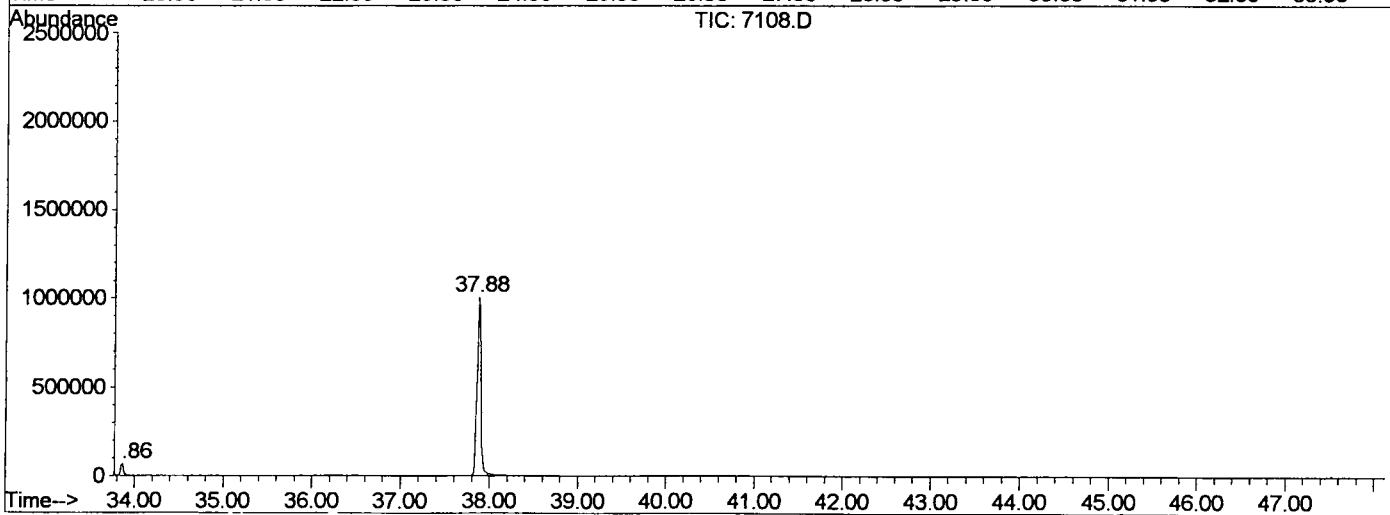
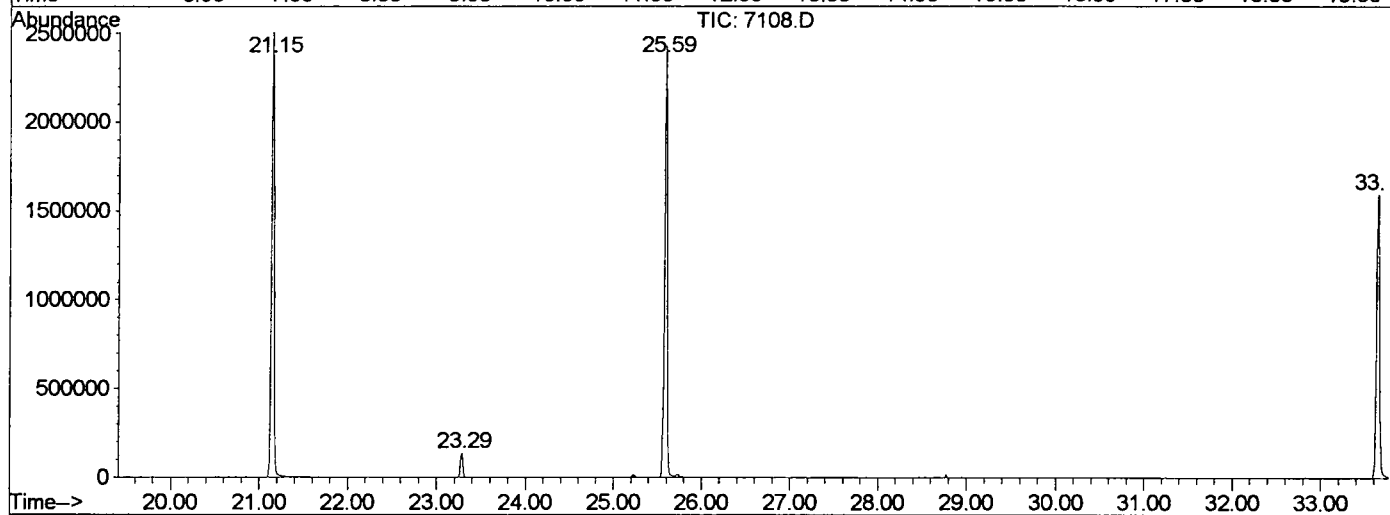
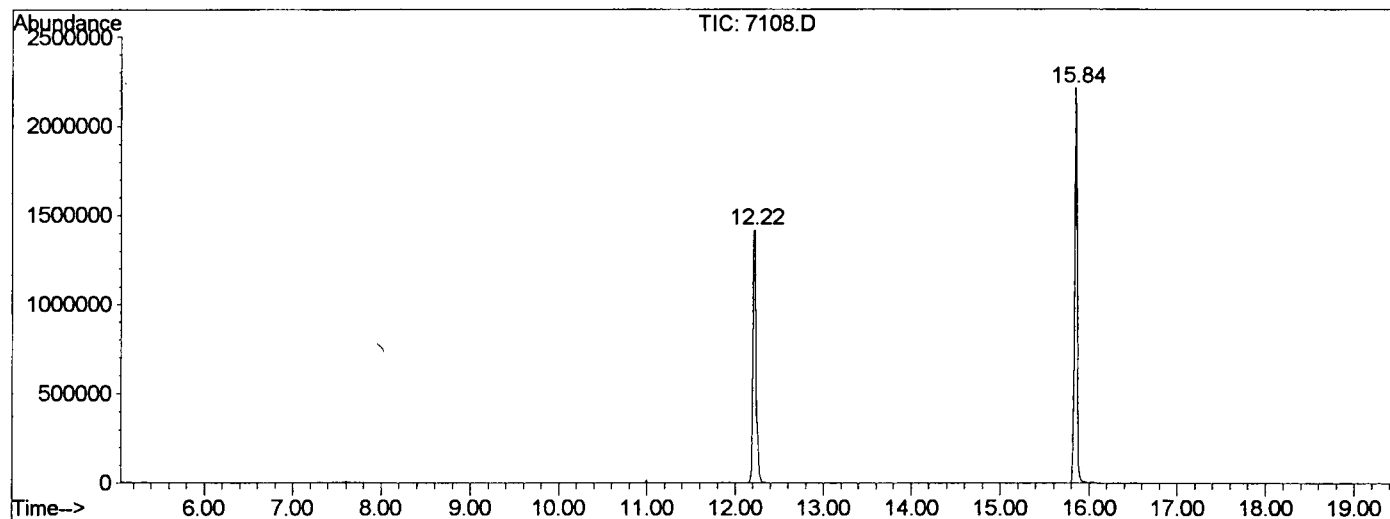
Sum of corrected areas: 25806109

7108.D GCMS0419.M

Mon Apr 22 08:34:21 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\APR1802\7108.D
 Operator :
 Acquired : 19 Apr 2002 3:38 pm using AcqMethod GCMS0412
 Instrument : 209
 Sample Name: GC/MS SCAN 3/26 FB
 Misc Info :
 Vial Number: 27
 Quant File :GCMS0419.RES (RTE Integrator)



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Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 19 Apr 2002 3:38 pm
 Data File: C:\HPCHEM\1\DATA\APR1802\7108.D
 Name: GC/MS SCAN 3/26 FB
 Misc:
 Method: C:\HPCHEM\1\METHODS\GCMS0419.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
7108.D	GCMS0419.M			Mon Apr 22 08:34:22 2002							