

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\APR1802\LBA.D
 Acq On : 19 Apr 2002 2:36 am
 Sample : GC/MS SCAN 3/26
 Misc :
 MS Integration Params: GOOFY.P
 Quant Time: Apr 19 14:52 2002

Vial: 16
 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	704639	20.00	ug/l	-0.02
10) Naphthalene-d8	15.84	136	2564417	20.00	ug/l	-0.03
17) Acenaphthene-d10	21.15	164	1436788	20.00	ug/l	-0.02
30) Phenanthrene-d10	25.59	188	2109694	20.00	ug/l	-0.02
39) Chrysene-d12	33.64	240	1386143	20.00	ug/l	-0.03
45) Perylene-d12	37.88	264	926274	20.00	ug/l	-0.04

System Monitoring Compounds

28) TCMX	23.29	209	37961	3.17	ug/L	-0.04
Spiked Amount	4.000		Recovery	=	79.25%	
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	15.90	128	404	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	21.52	165	183	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.	
29) Azobenzen/ 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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DataAcq Meth : GCMS0412

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
33) Hexachlorobenzene	0.00	284	0	N.D.		
34) Phenanthrene	25.58	178	675	N.D.		
35) Anthracene	25.58	178	675	N.D.		
36) Di-n-butylphthalate	27.55	149	501	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	33.64	228	3010	N.D.		
43) Bis(2-ethylhexyl)phthalate	33.86	149	2595	N.D.		
44) Chrysene	33.64	228	3009	N.D.		
46) Di-n-Octylphthalate	0.00	149	0	N.D.		
47) Benzo(b)fluoranthene	0.00	252	0	N.D.		
48) Benzo(k)fluoranthene	0.00	252	0	N.D.		
49) Benzo(a)pyrene	37.88	252	3646	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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LRA D GCMS0419 M

Fri Apr 19 14:53:25 2002

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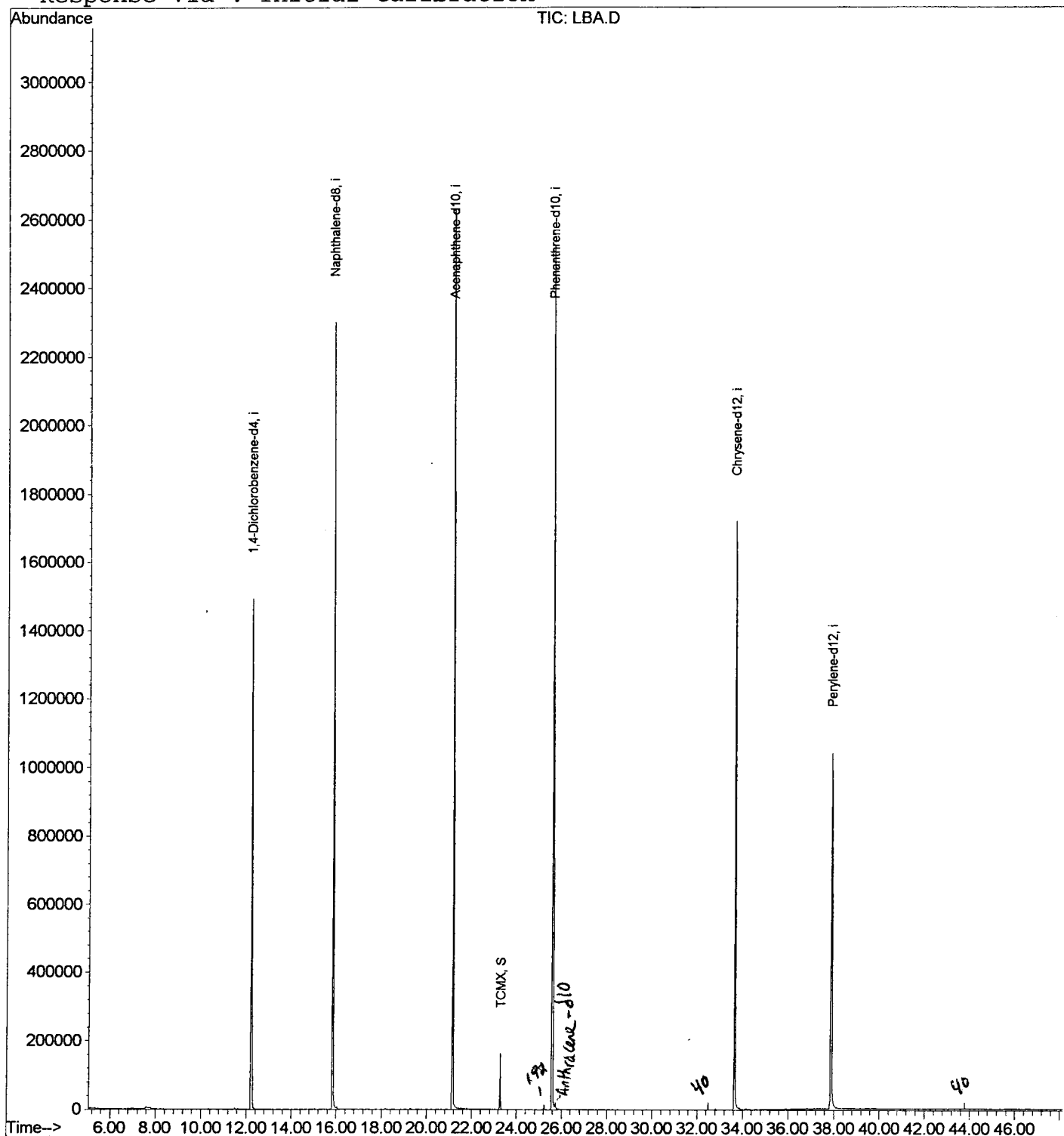
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 Last Update : Fri Apr 19 14:51:24 2002
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LSC Area Percent Report

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

Vial: 16

Acq On : 19 Apr 2002 2:36 am

Operator:

Sample : GC/MS SCAN 3/26

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	12.208	775	780	802	rBV	1492192	3748275	69.01%	14.077%
2	15.843	1170	1176	1193	rBV	2301681	4809038	88.55%	18.061%
3	21.150	1747	1754	1773	rBV	2631763	5431154	100.00%	20.398%
4	23.289	1980	1987	1993	rBV	163209	329836	6.07%	1.239%
5	25.593	2231	2238	2249	rBV	2523299	5213223	95.99%	19.579%
6	33.644	3108	3115	3134	rBV2	1721587	3979116	73.26%	14.944%
7	37.876	3567	3576	3595	rBV2	1040083	3115410	57.36%	11.701%

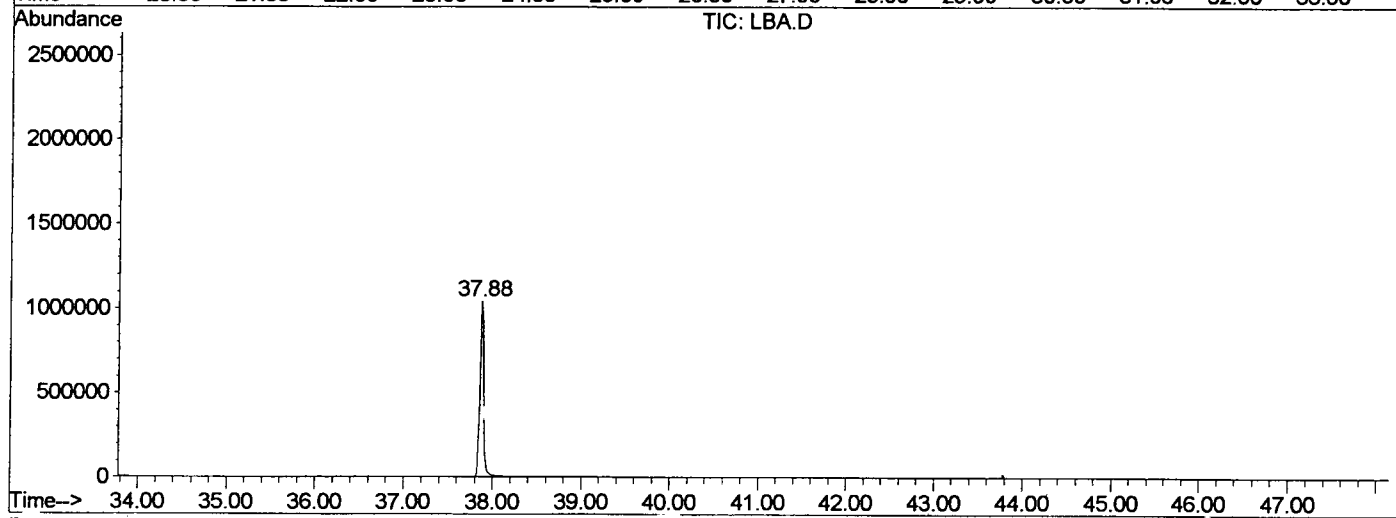
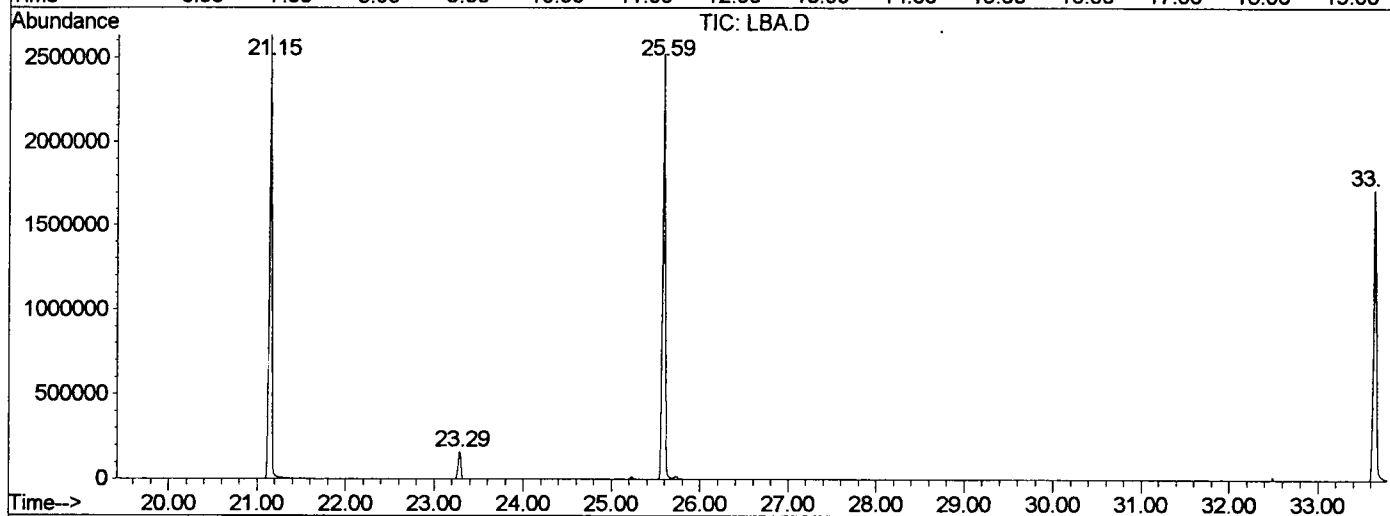
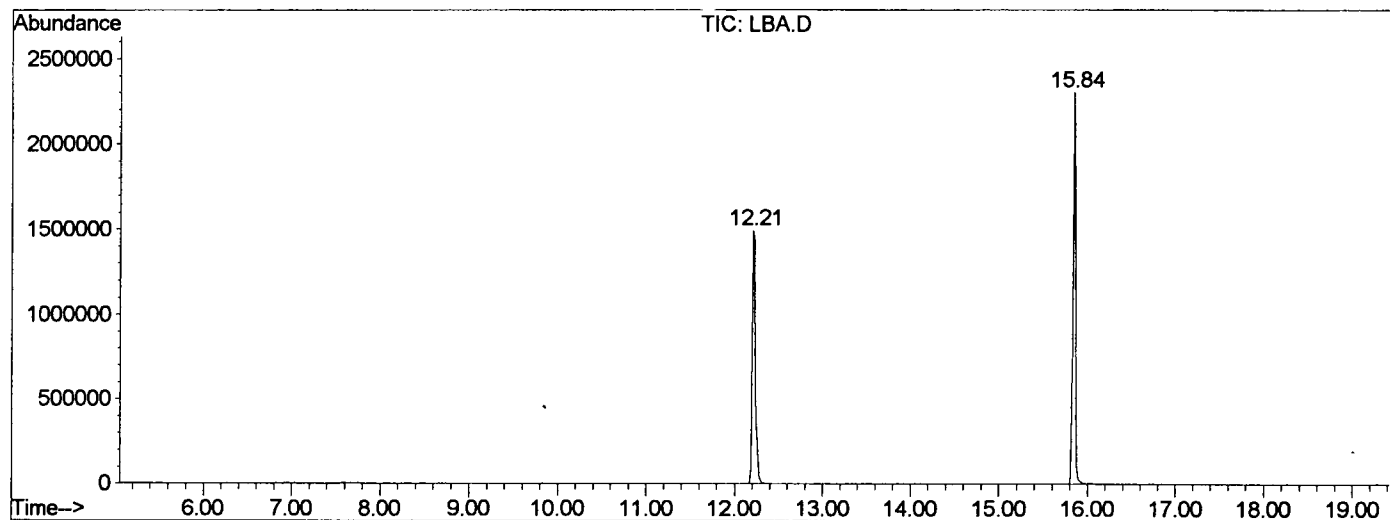
Sum of corrected areas: 26626052

LBA.D GCMS0419.M

Mon Apr 22 08:34:37 2002

LSC Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\APR1802\LBA.D
 Operator :
 Acquired : 19 Apr 2002 2:36 am using AcqMethod GCMS0412
 Instrument : 209
 Sample Name: GC/MS SCAN 3/26
 Misc Info :
 Vial Number: 16
 Quant File : GCMS0419.RES (RTE Integrator)



LBA.D GCMS0419.M Mon Apr 22 08:34:37 2002

Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 19 Apr 2002 2:36 am
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 Name: GC/MS SCAN 3/26
 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

LBA.D			GCMS0419.M	Mon Apr 22 08:34:38 2002							