

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\FEB2502\1962.D
 Acq On : 26 Feb 2002 10:28 pm
 Sample : gc/ms scan 1/29
 Misc :
 MS Integration Params: GOOFY.P
 Quant Time: Mar 19 10:45 2002

Vial: 34
 Operator:
 Inst : GC/MS 209
 Multiplr: 1.00

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	12.32	152	263580	20.00	ug/l	-0.03
10) Naphthalene-d8	15.96	136	1007762	20.00	ug/l	-0.04
17) Acenaphthene-d10	21.27	164	522569	20.00	ug/l	-0.04
29) Phenanthrene-d10	25.71	188	708376	20.00	ug/l	-0.05
38) Chrysene-d12	33.78	240	435641	20.00	ug/l	-0.06
44) Perylene-d12	38.05	264	292785	20.00	ug/l	-0.07

System Monitoring Compounds

37) 2,4-DDD 30.79 235 24113 3.26 ug/mL 0.29
 Spiked Amount 5.000 Recovery = 65.20%

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	0.00	165	0	N.D.	
25) Diethylphthalate	22.75	149	199	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

962.D GCMS0208.M Tue Mar 19 10:47:07 2002

Quantitation Report

(QT Reviewed)

Data File : C:\HPCHEM\1\DATA\FEB2502\1962.D

Vial: 34

Acq On : 26 Feb 2002 10:28 pm

Operator:

Sample : gc/ms scan 1/29

Inst : GC/MS 209

Misc :

Multiplr: 1.00

MS Integration Params: GOOFY.P

Quant Time: Mar 19 10:45 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

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	27.69	149	1969	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	1039	N.D.		
42) Bis(2-ethylhexyl)phthalate	33.99	149	875	N.D.		
43) Chrysene	33.78	228	1039	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	895	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

1962.D GCMS0208.M Tue Mar 19 10:47:10 2002

Page 2

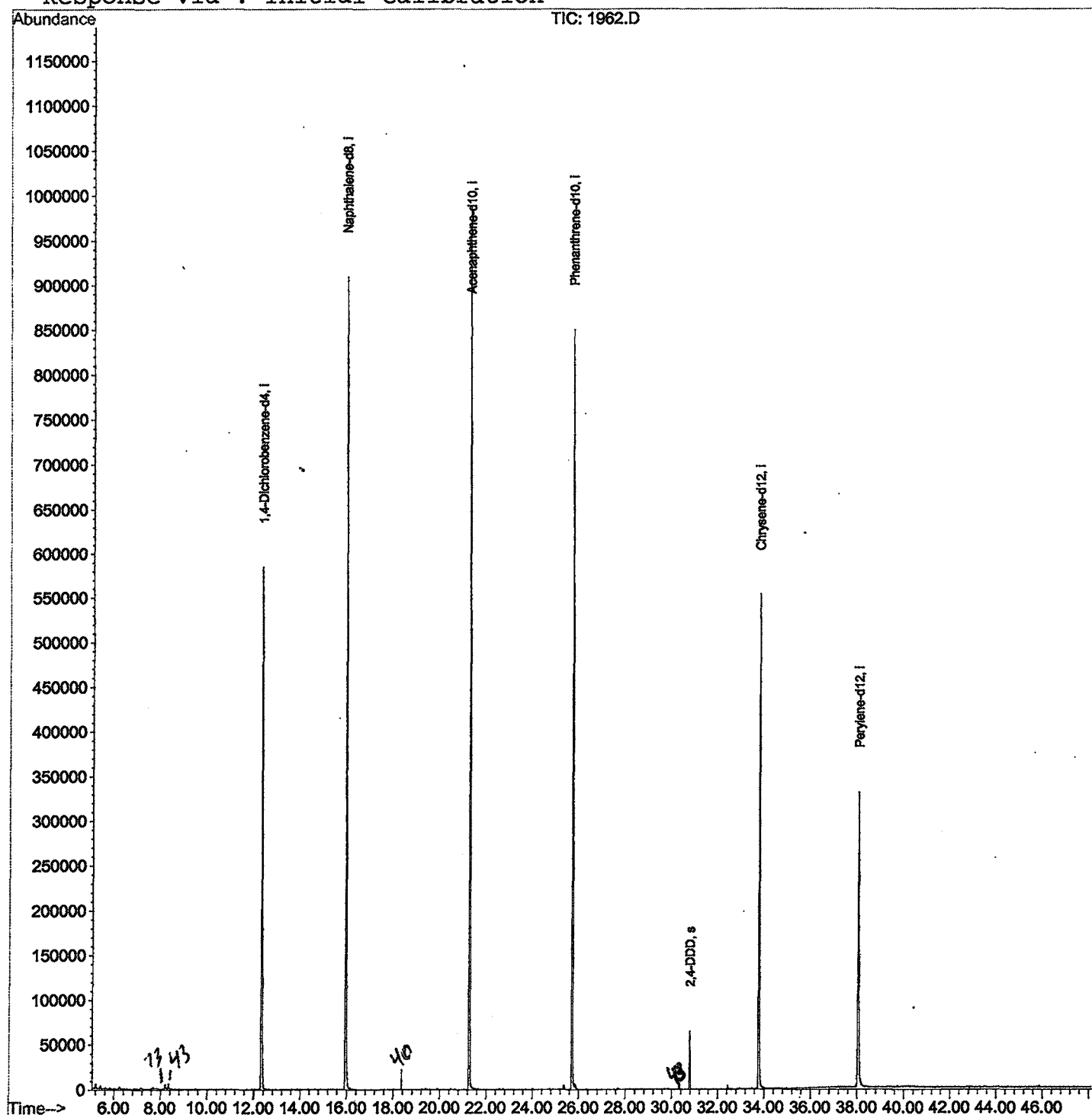
Quantitation Report

Data File : C:\HPCHEM\1\DATA\FEB2502\1962.D
 Acq On : 26 Feb 2002 10:28 pm
 Sample : gc/ms scan 1/29
 Misc :
 MS Integration Params: GOOFY.P
 Quant Time: Mar 19 10:45 2002

Vial: 34
 Operator:
 Inst : GC/MS 209
 Multiplr: 1.00

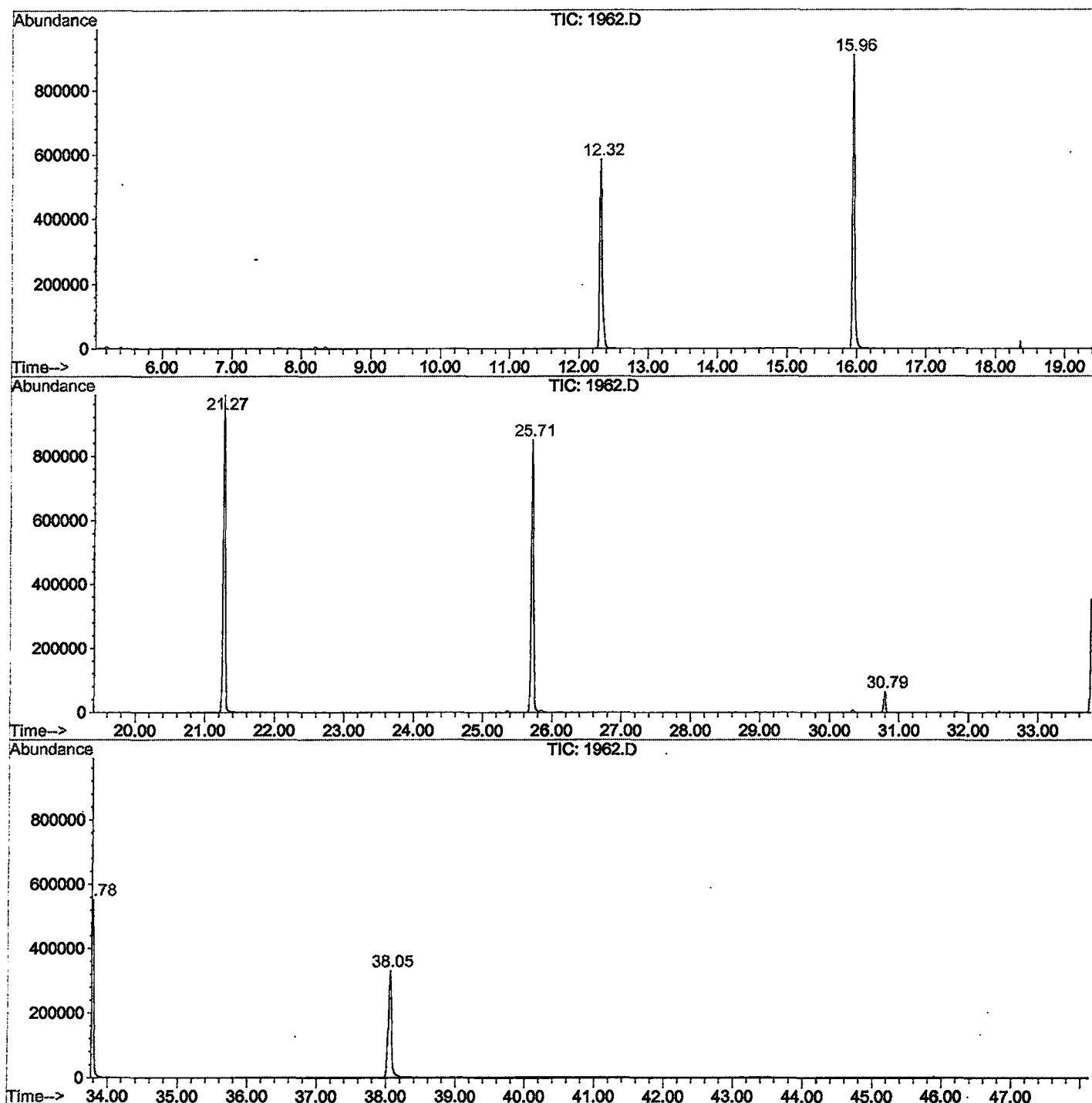
Quant Results File: GCMS0208.RES

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 Report - Integrated Chromatogram

File : C:\HPCHEM\1\DATA\FEB2502\1962.D
 Operator :
 Acquired : 26 Feb 2002 10:28 pm using AcqMethod GCMS0208
 Instrument : GC/MS 209
 Sample Name: gc/ms scan 1/29
 Misc Info :
 Vial Number: 34
 Quant File :GCMS0208.RES (RTE Integrator)



1962.D GCMS0208.M

Tue Mar 19 11:01:21 2002

LSC Area Percent Report

Data File : C:\HPCHEM\1\DATA\FEB2502\1962.D
 Acq On : 26 Feb 2002 10:28 pm
 Sample : gc/ms scan 1/29
 Misc :
 MS Integration Params: LSCINT.P

Vial: 34
 Operator:
 Inst : GC/MS 209
 Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\GCMS0208.M (RTE Integrator)
 Title : 209 625/TCLP calibration table
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 1 % of largest Peak
 Max Peaks: 100
 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	12.320	787	793	806	rBV	584230	1489940	72.15%	15.200%
2	15.956	1183	1189	1204	rBV	908722	1957575	94.80%	19.970%
3	21.271	1761	1768	1787	rBV	989413	2064976	100.00%	21.066%
4	25.714	2246	2252	2263	rBV2	849783	1849066	89.54%	18.863%
5	30.791	2800	2805	2811	rVB	64266	121530	5.89%	1.240%
6	33.784	3124	3131	3146	rBV2	552460	1303036	63.10%	13.293%
7	38.053	3588	3596	3614	rBV	328795	1016357	49.22%	10.368%

Sum of corrected areas: 9802480

1962.D GCMS0208.M Tue Mar 19 11:01:15 2002

Tentatively Identified Compound (LSC) summary

Operator ID: Date Acquired: 26 Feb 2002 10:28 pm
 Data File: C:\HPCHEM\1\DATA\FEB2502\1962.D
 Name: gc/ms scan 1/29
 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
1962.D	GCMS0208.M			Tue Mar 19 11:01:25 2002							