User: Vinci, David L
Westchester Dept. of Labs & Research
Date: 03/29/2002 Time: 08:19:00

Sample: AE03265
Analysis: $C1GCMS CALI GCMS
Units: ug
Started: 3/25/02 14:44 Ended: 3/25/02 14:44 Analyst: MF
Result source: manual entry
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	n-Nitrosodimethylamine		9.76		2.0
2	bis(2-Chloroethyl)ether		11.55		2.0
3	1,3-Dichlorobenzene		11.66		2.0
4	1,4-Dichlorobenzene		11.77		2.0
5	1,2-Dichlorobenzene		11.82		2.0
6	2,2'-oxybis(1-Chloropropane)		11.85		2.0
7	n-Nitrosodi-n-propylamine		12.10		2.0
8	Hexachloroethane		12.08		2.0
9	Nitrobenzene		10.30		2.0
10	Isophorone		9.82		2.0
11	bis(2-Chloroethoxy)methane		10.43		2.0
12	1,2,4-Trichlorobenzene		10.94		2.0
13	Naphthalene		10.92		2.0
14	Hexachlorobutadiene		11.14		2.0
15	2-Chloronaphthalene		11.15		2.0
16	Dimethylphthalate		10.53		2.0
17	2,6-Dinitrotoluene		9.95		2.0
18	Acenaphthylene		10.85		2.0
19	Acenaphthene		10.97		2.0
20	2,4-Dinitrotoluene		9.58		2.0
21	Diethylphthalate		11.30		2.0
22	4-Chlorophenylphenylether		10.96		2.0
23	Fluorene		11.45		2.0
24	Azobenzene/1,2-Diphenylhydraz		10.25		2.0
25	n-Nitrosodiphenylamine		10.99		2.0
26	4-Bromophenylphenylether		11.54		2.0
27	Hexachlorobenzene		11.34		2.0
28	Phenanthrene		11.27		2.0
29	Anthracene		10.84		2.0
30	Di-n-butylphthalate		10.90		2.0
31	Fluoranthene		11.10		2.0
32	Pyrene		9.68		2.0
33	Butylbenzylphthalate		9.04		2.0
34	Benzo(a)anthracene		9.58		2.0
35	bis(2-Ethylhexyl)phthalate		8.88		2.0
36	Chrysene		10.54		2.0
37	Di-n-octylphthalate		8.96		2.0
38	Benzo(b)fluoranthene		10.31		2.0
39	Benzo(k)fluoranthene		11.53		2.0
40	Benzo(a)pyrene		9.84		2.0
41	Indeno(1,2,3-cd)pyrene		8.09		2.0
42	Dibenz(a,h)anthracene		8.87		2.0
43	Benzo(g,h,i)perylene		9.66		2.0

Quantitation Report (QT Reviewed)

Data File : C:\HPCHEM\1\DATA\MAR0102\CAL10B.D Vial: 2
Acq On : 2 Mar 2002 10:42 pm Operator:
Sample : Inst : GC/MS 209
Misc : Multiplr: 1.00
MS Integration Params: GOOFY.P
Quant Time: Mar 25 11:56 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	23.30	168	405157	19.25 ug/l	#	84
31)	4-Bromophenyl-phenylether	24.38	248	100700	7.40 ug/l		92
32)	Hexachlorobenzene	24.81	284	116812	7.33 ug/l		96
33)	Phenanthrene	25.79	178	542244	8.35 ug/l		99
34)	Anthracene	25.92	178	547206	8.60 ug/l		99
35)	Di-n-butylphthalate	27.68	149	700090	7.72 ug/l		98
36)	Fluoranthene	29.41	202	556039	8.58 ug/l	#	89
39)	Pyrene	30.09	202	568250	8.20 ug/l	#	88
40)	Butylbenzylphthalate	32.22	149	289273	7.53 ug/l		99
41)	Benzo(a)anthracene	33.73	228	420346	8.29 ug/l		99
42)	Bis(2-ethylhexyl)phthalate	33.99	149	363971	7.02 ug/l		95
43)	Chrysene	33.86	228	407042	8.16 ug/l		99
45)	Di-n-Octylphthalate	35.73	149	526612	6.47 ug/l	#	98
46)	Benzo(b)fluoranthene	36.85	252	375669	8.18 ug/l	#	97
47)	Benzo(k)fluoranthene	36.92	252	401482	8.69 ug/l	#	93
48)	Benzo(a)pyrene	37.85	252	310947	8.16 ug/l	#	95
49)	Indeno(1,2,3-cd)pyrene	42.24	276	305771	8.80 ug/l	#	94
50)	Dibenz(a,h)anthracene	42.31	278	239367	9.01 ug/l	#	94
51)	Benzo(g,h,i)perylene	43.50	276	248568	8.87 ug/l	#	93

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

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Sample : Inst : GC/MS 209
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MS Integration Params: GOOFY.P
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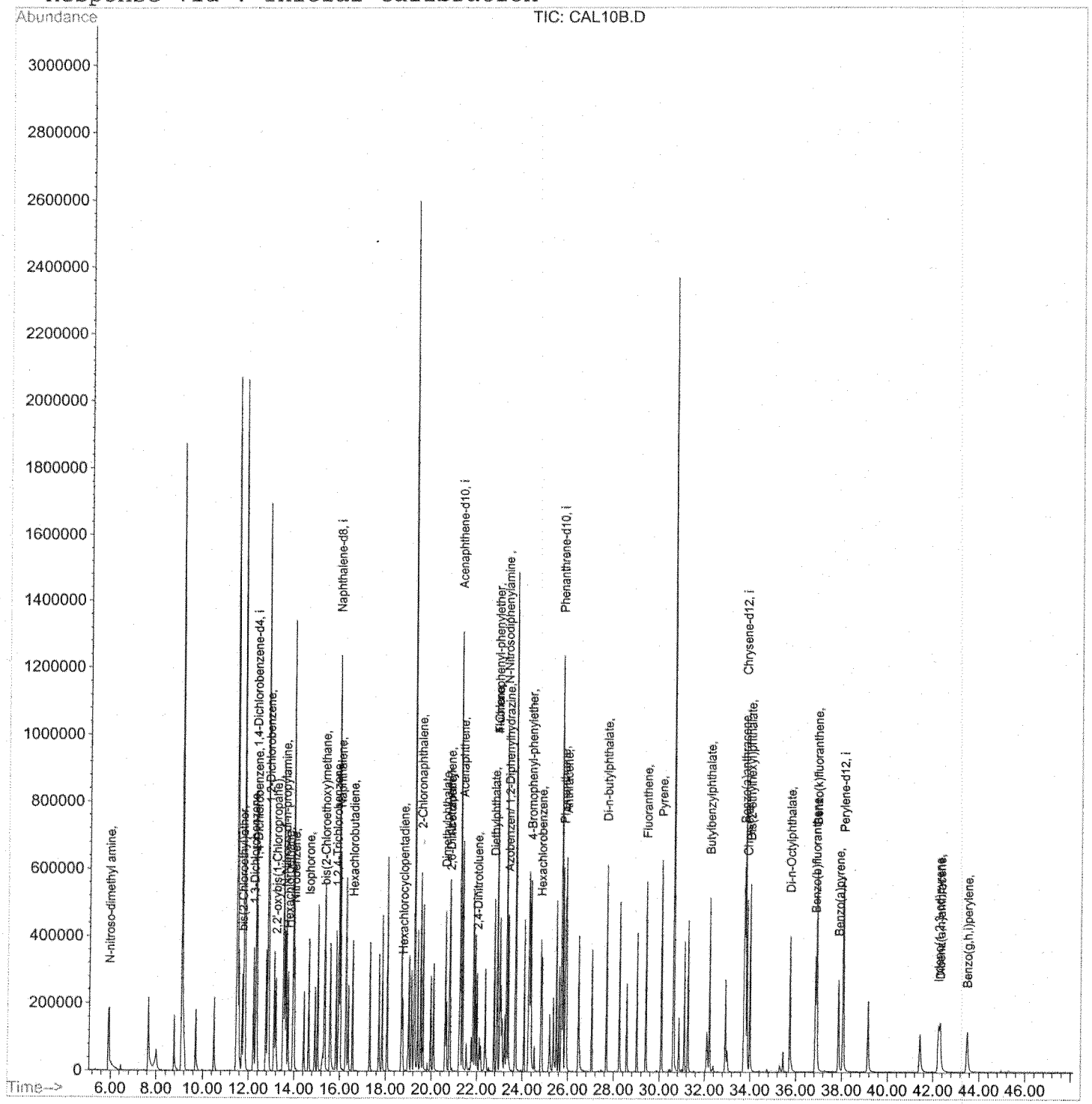

CAL10B.D GCMS0208.M Mon Mar 25 11:59:38 2002 Page 3