

Data Path : Z:\HPCHEM1\BNA G\DATA\BG041316\
 Data File : BG021625.D
 Acq On : 13 Apr 2016 17:34
 Operator : UM/SJ
 Sample : H1739-03DL 5X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PT-BN-WPDL

Quant Time: Apr 14 02:18:17 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 15:48:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.22	152	27626	20.00	ng	0.00
21) Naphthalene-d8	11.04	136	129847	20.00	ng	0.00
38) Acenaphthene-d10	14.83	164	86021	20.00	ng	0.00
63) Phenanthrene-d10	17.56	188	210455	20.00	ng	0.00
75) Chrysene-d12	21.83	240	244869	20.00	ng	0.00
86) Perylene-d12	25.16	264	237751	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.76	112	42294	25.49	ng	0.00
7) Phenol-d6	7.36	99	65958	26.62	ng	0.00
23) Nitrobenzene-d5	9.38	82	40157	15.90	ng	0.00
41) 2,4,6-Tribromophenol	16.30	330	21326	23.00	ng	0.00
44) 2-Fluorobiphenyl	13.46	172	104328	17.77	ng	0.00
78) Terphenyl-d14	20.14	244	175756	17.91	ng	0.00

Target Compounds						Qvalue
4) n-Nitrosodimethylamine	3.96	42	24381	22.20	ng	# 90
11) bis(2-Chloroethyl)ether	7.63	93	41614	19.51	ng	95
12) 1,3-Dichlorobenzene	8.11	146	61384	27.64	ng	94
18) Hexachloroethane	9.31	117	16733	20.33	ng	# 84
24) Nitrobenzene	9.43	77	35597	13.16	ng	95
25) Isophorone	9.95	82	138713	25.09	ng	98
30) 1,2,4-Trichlorobenzene	10.90	180	16524	7.68	ng	93
31) Naphthalene	11.08	128	28842	4.15	ng	# 95
37) 2-Methylnaphthalene	12.67	142	56990	11.95	ng	99
46) 2-Chloronaphthalene	13.71	162	155908	29.10	ng	96
49) Dimethylphthalate	14.27	163	218542	29.38	ng	100
50) 2,6-Dinitrotoluene	14.39	165	32162	22.68	ng	94
51) Acenaphthene	14.89	154	109194	19.28	ng	98
56) 2,4-Dinitrotoluene	15.17	165	16768	8.75	ng	# 99
57) Fluorene	15.87	166	149249	21.61	ng	96
59) Diethylphthalate	15.62	149	132300	17.04	ng	99
60) 4-Chlorophenyl-phenylether	15.85	204	76256	22.42	ng	99
66) 4-Bromophenyl-phenylether	16.75	248	8241	3.65	ng	92
67) Hexachlorobenzene	16.86	284	21731	9.13	ng	99
70) Phenanthrene	17.61	178	369059	31.82	ng	98
71) Anthracene	17.69	178	74378	6.32	ng	98
73) Di-n-butylphthalate	18.50	149	238178	17.04	ng	100
74) Fluoranthene	19.60	202	215108	15.53	ng	99
79) Butylbenzylphthalate	20.83	149	78556	11.99	ng	100
83) Bis(2-ethylhexyl)phthalate	21.71	149	314735	32.83	ng	99
84) Di-n-octyl phthalate	22.96	149	511833	31.87	ng	100
85) Indeno(1,2,3-cd)pyrene	28.97	276	106276	6.60	ng	# 83
87) Benzo(b)fluoranthene	24.10	252	350002	25.37	ng	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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