

Data Path : Z:\HPCHEM1\BNA F\DATA\BF090315\
 Data File : BF081399.D
 Acq On : 4 Sep 2015 3:15
 Operator : UM/IZ
 Sample : G3419-06
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HX2

Quant Time: Sep 04 06:01:35 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 04 02:03:29 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.52	152	231814	20.00	ng	0.15
21) Naphthalene-d8	7.82	136	472	20.00	ng	0.15
38) Acenaphthene-d10	9.61	164	276	20.00	ng	0.21
63) Phenanthrene-d10	11.13	188	214	20.00	ng	0.26
75) Chrysene-d12	0.00	240	0	0.00	ng	-13.47
86) Perylene-d12	0.00	264	0	0.00	ng	-14.79
System Monitoring Compounds						
5) 2-Fluorophenol	4.91	112	191744	12.83	ng	0.00
7) Phenol-d6	6.03	99	170661	8.63	ng	-0.01
23) Nitrobenzene-d5	6.95	82	431776	44135.12	ng	0.00
41) 2,4,6-Tribromophenol	10.19	330	115281	46909.64	ng	0.00
44) 2-Fluorobiphenyl	8.74	172	622049	28882.10	ng	-0.01
78) Terphenyl-d14	12.46	244	637557	0.00	ng	0.00
Target Compounds						Qvalue
18) Hexachloroethane	6.95	117	89316	12.82	ng	# 6
19) n-Nitroso-di-n-propylamine	6.95	70	61262	4.55	ng	# 63
22) Acetophenone	6.92	105	7499	581.67	ng	# 22
24) Nitrobenzene	7.07	77	2984	293.72	ng	# 7
25) Isophorone	7.39	82	1406	77.27	ng	# 19
26) 2-Nitrophenol	7.66	139	209	47.20	ng	# 1
27) 2,4-Dimethylphenol	7.46	122	1409	184.23	ng	# 12
28) bis(2-Chloroethoxy)methane	7.66	93	246	23.40	ng	# 1
29) 2,4-Dichlorophenol	7.71	162	230	34.68	ng	# 1
31) Naphthalene	7.86	128	5118	203.32	ng	# 1
32) Benzoic acid	7.66	122	322	58.39	ng	# 1
33) 4-Chloroaniline	7.91	127	1561	152.04	ng	# 1
35) Caprolactam	8.32	113	248	121.93	ng	# 1
36) 4-Chloro-3-methylphenol	8.41	107	4692	632.06	ng	# 12
37) 2-Methylnaphthalene	8.47	142	73912	4512.02	ng	# 82
45) 1,1'-Biphenyl	9.06	154	6695	263.66	ng	# 1
46) 2-Chloronaphthalene	9.06	162	6257	341.21	ng	# 1
47) 2-Nitroaniline	9.14	65	3682	492.67	ng	# 1
48) Acenaphthylene	9.51	152	4277	131.47	ng	# 1
49) Dimethylphthalate	9.15	163	142684	6653.82	ng	# 97
50) 2,6-Dinitrotoluene	9.45	165	2028	416.69	ng	90
51) Acenaphthene	9.43	154	16098	869.84	ng	91
52) 3-Nitroaniline	9.60	138	981	177.05	ng	# 59
53) 2,4-Dinitrophenol	9.70	184	217	105.86	ng	# 1
54) Dibenzofuran	10.00	168	37147	1374.66	ng	# 70
55) 4-Nitrophenol	9.80	139	2288	657.30	ng	# 1
56) 2,4-Dinitrotoluene	9.84	165	3209	517.82	ng	# 14
57) Fluorene	10.18	166	6345	309.41	ng	# 77
59) Diethylphthalate	10.02	149	1043	46.24	ng	# 10
60) 4-Chlorophenyl-phenylether	10.18	204	532	55.66	ng	# 1
61) 4-Nitroaniline	10.19	138	629	112.37	ng	# 1
62) Azobenzene	10.31	77	791	31.54	ng	# 15
64) 4,6-Dinitro-2-methylphenol	10.31	198	258	215.55	ng	# 1
65) n-Nitrosodiphenylamine	10.31	169	10977	1409.65	ng	# 1

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
68) Atrazine	10.88	200	697	309.31	ng	# 1
70) Phenanthrene	11.11	178	1247	104.81	ng	# 1
71) Anthracene	11.22	178	469	38.37	ng	# 1
72) Carbazole	11.32	167	1197	104.36	ng	# 1
73) Di-n-butylphthalate	11.46	149	14025	1035.50	ng	# 94
74) Fluoranthene	12.28	202	4473	347.30	ng	# 88

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

