

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100516\
 Data File : BF090388.D
 Acq On : 5 Oct 2016 19:12
 Operator : UM/SJ
 Sample : H4818-06
 Misc : LOO 20PPM
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOQ-WATER2016-06-QT4

Manual Integrations
 APPROVED

SOHIL
 10/6/2016 5:28:38 PM

Quant Time: Oct 06 01:40:12 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF100516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 05 14:57:00 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	221456	20.00	ng	0.00
21) Naphthalene-d8	8.29	136	977236	20.00	ng	0.00
38) Acenaphthene-d10	10.05	164	479660	20.00	ng	0.00
63) Phenanthrene-d10	11.54	188	738907	20.00	ng	-0.01
75) Chrysene-d12	14.19	240	610808	20.00	ng	-0.01
86) Perylene-d12	15.70	264	419566	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	1717959	115.78	ng	0.01
7) Phenol-d6	6.62	99	2389578	122.96	ng	0.00
23) Nitrobenzene-d5	7.56	82	1811291	90.16	ng	-0.01
41) 2,4,6-Tribromophenol	10.85	330	545189	139.78	ng	0.00
44) 2-Fluorobiphenyl	9.38	172	2908997	101.05	ng	0.00
78) Terphenyl-d14	13.14	244	2573343	94.72	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.82	88	145483	19.89	ng	94
3) Pyridine	3.58	79	415649m	20.38	ng	
4) n-Nitrosodimethylamine	3.50	42	168044	19.97	ng	# 69
6) Aniline	6.66	93	444027	16.54	ng	98
8) 2-Chlorophenol	6.78	128	317102	19.42	ng	86
9) Benzaldehyde	6.54	77	192897	19.64	ng	88
10) Phenol	6.63	94	377320	16.70	ng	88
11) bis(2-Chloroethyl)ether	6.74	93	349714	20.21	ng	97
12) 1,3-Dichlorobenzene	6.94	146	342895m	21.07	ng	
13) 1,4-Dichlorobenzene	7.02	146	329409	19.92	ng	96
14) 1,2-Dichlorobenzene	7.17	146	308255	19.35	ng	95
15) Benzyl Alcohol	7.14	79	331069	22.41	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.27	45	605631	20.64	ng	97
17) 2-Methylphenol	7.25	107	271900	20.25	ng	97
18) Hexachloroethane	7.51	117	124942	19.17	ng	# 86
19) n-Nitroso-di-n-propylamine	7.41	70	302045	21.20	ng	# 85
20) 3+4-Methylphenols	7.40	107	366482	21.08	ng	# 86
22) Acetophenone	7.41	105	472392	20.41	ng	# 87
24) Nitrobenzene	7.58	77	437371	21.93	ng	99
25) Isophorone	7.82	82	742692	20.22	ng	99
26) 2-Nitrophenol	7.90	139	168283	19.60	ng	# 88
27) 2,4-Dimethylphenol	7.94	122	335027	20.07	ng	99
28) bis(2-Chloroethoxy)methane	8.03	93	408377	18.80	ng	# 95
29) 2,4-Dichlorophenol	8.14	162	245862	19.37	ng	94
30) 1,2,4-Trichlorobenzene	8.23	180	245888	18.80	ng	95
31) Naphthalene	8.31	128	970364	20.58	ng	97
32) Benzoic acid	8.02	122	218582	18.83	ng	91
33) 4-Chloroaniline	8.36	127	283145	14.53	ng	# 90
34) Hexachlorobutadiene	8.43	225	124197	16.99	ng	99
35) Caprolactam	8.70	113	76032	17.82	ng	92
36) 4-Chloro-3-methylphenol	8.83	107	284958	17.93	ng	91
37) 2-Methylnaphthalene	9.00	142	556112	19.49	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.17	216	235861	18.65	ng	# 96
40) Hexachlorocyclopentadiene	9.16	237	118895	17.15	ng	98

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42) 2,4,6-Trichlorophenol	9.29	196	161681	18.53	nq	98
43) 2,4,5-Trichlorophenol	9.32	196	169552	20.55	nq	94
45) 1,1'-Biphenyl	9.47	154	714996	20.12	nq	93
46) 2-Chloronaphthalene	9.49	162	517681	19.71	nq	96
47) 2-Nitroaniline	9.58	65	214668	20.72	nq	# 85
48) Acenaphthylene	9.91	152	911605	21.04	nq	97
49) Dimethylphthalate	9.77	163	607635	19.78	nq	98
50) 2,6-Dinitrotoluene	9.82	165	120432	17.66	nq	# 80
51) Acenaphthene	10.09	154	562086	20.80	nq	98
52) 3-Nitroaniline	9.99	138	143120	17.96	nq	98
53) 2,4-Dinitrophenol	10.10	184	53970	17.75	nq	# 43
54) Dibenzofuran	10.26	168	712758	20.38	nq	95
55) 4-Nitrophenol	10.14	139	134830	19.34	nq	92
56) 2,4-Dinitrotoluene	10.22	165	168133	18.52	nq	# 1
57) Fluorene	10.60	166	603711	20.54	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.37	232	116978	18.98	nq	# 75
59) Diethylphthalate	10.47	149	668131	21.02	nq	# 93
60) 4-Chlorophenyl-phenylether	10.59	204	253693	18.98	nq	# 81
61) 4-Nitroaniline	10.60	138	154773	19.28	nq	94
62) Azobenzene	10.75	77	685083	18.66	nq	87
64) 4,6-Dinitro-2-methylphenol	10.63	198	71518	18.15	nq	95
65) n-Nitrosodiphenylamine	10.71	169	503424	20.33	nq	99
66) 4-Bromophenyl-phenylether	11.08	248	134658	18.59	nq	# 67
67) Hexachlorobenzene	11.15	284	153428	19.02	nq	# 68
68) Atrazine	11.23	200	107978	17.69	nq	95
69) Pentachlorophenol	11.34	266	100773	20.99	nq	98
70) Phenanthrene	11.57	178	782726	20.69	nq	95
71) Anthracene	11.62	178	832307	21.54	nq	97
72) Carbazole	11.77	167	744176	20.73	nq	98
73) Di-n-butylphthalate	12.11	149	969902	22.20	nq	# 95
74) Fluoranthene	12.76	202	789708	21.28	nq	94
76) Benzidine	12.87	184	240385	14.97	nq	98
77) Pyrene	12.99	202	830895	20.40	nq	98
79) Butylbenzylphthalate	13.61	149	397832	19.04	nq	# 73
80) Benzo(a)anthracene	14.18	228	691217	20.17	nq	97
81) 3,3'-Dichlorobenzidine	14.14	252	195520	15.74	nq	# 97
82) Chrysene	14.21	228	601214	19.06	nq	97
83) Bis(2-ethylhexyl)phthalate	14.17	149	590348	19.87	nq	# 92
84) Di-n-octyl phthalate	14.80	149	916374	20.81	nq	95
85) Indeno(1,2,3-cd)pyrene	17.21	276	410816	18.27	nq	94
87) Benzo(b)fluoranthene	15.25	252	516630	19.33	nq	# 97
88) Benzo(k)fluoranthene	15.28	252	513290m	20.45	nq	
89) Benzo(a)pyrene	15.63	252	452079	19.63	nq	97
90) Dibenzo(a,h)anthracene	17.23	278	343069	17.51	nq	# 95
91) Benzo(g,h,i)perylene	17.66	276	322962	17.17	nq	94

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

