

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050818\
 Data File : BF105166.D
 Acq On : 9 May 2018 1:11
 Operator : JU/SJ
 Sample : J2133-09
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MDL-WATER-03-QT2-2018

Manual Integrations
 APPROVED

Sohil
 5/9/2018 6:51:07 PM

Quant Time: May 09 09:34:34 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 08 16:33:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	203464	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	805100	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	397800	20.00	ng	-0.01
63) Phenanthrene-d10	11.31	188	777973	20.00	ng	0.00
75) Chrysene-d12	14.00	240	667793	20.00	ng	-0.02
86) Perylene-d12	15.54	264	581926	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.40	112	1534843	124.28	ng	0.00
7) Phenol-d6	6.42	99	1823388	123.99	ng	0.00
23) Nitrobenzene-d5	7.36	82	1072362	90.27	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	575610	131.15	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1965315	76.68	ng	0.00
78) Terphenyl-d14	12.91	244	2156520	75.01	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.52	88	7124	1.091	ng	98
3) Pyridine	3.37	79	16094	0.919	ng	97
4) n-Nitrosodimethylamine	3.20	42	6605	0.744	ng	# 83
6) Aniline	6.46	93	20291	0.945	ng	# 89
8) 2-Chlorophenol	6.57	128	14454	1.103	ng	95
9) Benzaldehyde	6.35	77	6233	0.553	ng	92
10) Phenol	6.43	94	16941	1.035	ng	# 1
11) bis(2-Chloroethyl)ether	6.53	93	14776	1.096	ng	94
12) 1,3-Dichlorobenzene	6.73	146	18394	1.165	ng	96
13) 1,4-Dichlorobenzene	6.81	146	18108	1.147	ng	98
14) 1,2-Dichlorobenzene	6.96	146	16431	1.128	ng	96
15) Benzyl Alcohol	6.93	79	10814	0.961	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.06	45	26231	1.047	ng	98
17) 2-Methylphenol	7.03	107	12312	1.088	ng	# 94
18) Hexachloroethane	7.30	117	5683	1.118	ng	92
19) n-Nitroso-di-n-propylamine	7.20	70	9949	1.022	ng	93
20) 3+4-Methylphenols	7.19	107	15447	1.162	ng	# 87
22) Acetophenone	7.19	105	23774	1.267	ng	# 93
24) Nitrobenzene	7.37	77	18487	1.373	ng	# 96
25) Isophorone	7.36	82	1071634	40.337	ng	# 64
26) 2-Nitrophenol	7.69	139	4400	10.574	ng	# 88
27) 2,4-Dimethylphenol	7.72	122	9728	0.824	ng	98
28) bis(2-Chloroethoxy)methane	7.82	93	16475	1.005	ng	97
29) 2,4-Dichlorophenol	7.92	162	10574	0.957	ng	93
30) 1,2,4-Trichlorobenzene	8.01	180	15166	1.191	ng	99
31) Naphthalene	8.09	128	47235	1.218	ng	97
32) Benzoic acid	7.75	122	2828	9.540	ng	93
33) 4-Chloroaniline	8.14	127	14885	0.923	ng	96
34) Hexachlorobutadiene	8.21	225	8668	1.175	ng	99
35) Caprolactam	8.46	113	2426	0.723	ng	91
36) 4-Chloro-3-methylphenol	8.60	107	9762	0.849	ng	95
37) 2-Methylnaphthalene	8.78	142	29786	1.102	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.95	216	14382	1.136	ng	97
40) Hexachlorocyclopentadiene	8.93	237	5804	0.848	ng	94

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.06	196	7022	0.909	nq	94
43) 2,4,5-Trichlorophenol	9.09	196	7652	0.948	nq	89
45) 1,1'-Biphenyl	9.25	154	46900	1.447	nq	96
46) 2-Chloronaphthalene	9.27	162	29519	1.219	nq	96
47) 2-Nitroaniline	9.36	65	5116	6.444	nq	97
48) Acenaphthylene	9.68	152	42699	1.141	nq	100
49) Dimethylphthalate	9.54	163	31993	1.188	nq	99
50) 2,6-Dinitrotoluene	9.60	165	4631	0.793	nq	94
51) Acenaphthene	9.86	154	29503	1.224	nq	95
52) 3-Nitroaniline	9.77	138	5092	0.827	nq	88
53) 2,4-Dinitrophenol	9.87	184	486	9.621	nq #	16
54) Dibenzofuran	10.03	168	40981	1.162	nq	93
55) 4-Nitrophenol	9.92	139	2877	4.792	nq	90
56) 2,4-Dinitrotoluene	10.00	165	5511	4.264	nq #	94
57) Fluorene	10.37	166	31752	1.228	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.14	232	6205	0.861	nq #	79
59) Diethylphthalate	10.25	149	29457	1.113	nq #	93
60) 4-Chlorophenyl-phenylether	10.37	204	16347	1.358	nq #	87
61) 4-Nitroaniline	10.37	138	4835	0.789	nq	99
62) Azobenzene	10.53	77	27789	1.007	nq	96
64) 4,6-Dinitro-2-methylphenol	10.41	198	1383	9.263	nq	80
65) n-Nitrosodiphenylamine	10.48	169	26959	1.112	nq	97
66) 4-Bromophenyl-phenylether	10.86	248	9504	1.145	nq #	86
67) Hexachlorobenzene	10.92	284	11699	1.224	nq #	82
68) Atrazine	11.00	200	4811	0.633	nq	97
69) Pentachlorophenol	11.11	266	3611	3.840	nq	97
70) Phenanthrene	11.33	178	51053	1.245	nq	98
71) Anthracene	11.39	178	45733	1.097	nq	96
72) Carbazole	11.54	167	41132	1.095	nq	98
73) Di-n-butylphthalate	11.87	149	34288	0.854	nq	99
74) Fluoranthene	12.52	202	45628	1.068	nq	92
76) Benzidine	12.65	184	5050m	0.203	nq	
77) Pyrene	12.75	202	47647	0.981	nq	99
79) Butylbenzylphthalate	13.39	149	9550	7.365	nq	94
80) Benzo(a)anthracene	13.99	228	41478	1.024	nq	98
81) 3,3'-Dichlorobenzidine	13.95	252	7414	0.487	nq	96
82) Chrysene	14.02	228	46336	1.111	nq	93
83) Bis(2-ethylhexyl)phthalate	13.99	149	13040	0.649	nq	99
84) Di-n-octyl phthalate	14.67	149	13807	7.239	nq #	88
85) Indeno(1,2,3-cd)pyrene	16.97	276	25849	0.782	nq	94
87) Benzo(b)fluoranthene	15.11	252	32865	0.923	nq	95
88) Benzo(k)fluoranthene	15.13	252	34563	1.001	nq	97
89) Benzo(a)pyrene	15.47	252	27885	0.861	nq	97
90) Dibenzo(a,h)anthracene	16.99	278	20131	0.800	nq #	96
91) Benzo(g,h,i)perylene	17.40	276	20637	0.839	nq	94

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

