

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
 Data File : BF091850.D
 Acq On : 21 Dec 2016 22:56
 Operator : UM/SJ
 Sample : H6103-06
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOQ-WATER-06-QT1-2017

Manual Integrations
 APPROVED

sohil
 12/22/2016 7:09:01 PM

Quant Time: Dec 22 14:17:09 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 21 16:17:51 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	332358	20.00	ng	0.00
21) Naphthalene-d8	8.03	136	1357798	20.00	ng	-0.01
38) Acenaphthene-d10	9.78	164	717786	20.00	ng	-0.01
63) Phenanthrene-d10	11.26	188	913667	20.00	ng	0.00
75) Chrysene-d12	13.89	240	667143	20.00	ng	-0.01
86) Perylene-d12	15.29	264	678485	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.37	112	2403153	120.73	ng	0.02
7) Phenol-d6	6.40	99	3212908	132.21	ng	-0.01
23) Nitrobenzene-d5	7.31	82	2578273	105.59	ng	-0.01
41) 2,4,6-Tribromophenol	10.58	330	701852	118.15	ng	0.00
44) 2-Fluorobiphenyl	9.12	172	3820179	113.35	ng	0.00
78) Terphenyl-d14	12.85	244	2359032	85.76	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.43	88	148259	15.13	ng	97
3) Pyridine	3.14	79	422816m	12.36	ng	
4) n-Nitrosodimethylamine	3.05	42	156852	12.16	ng	97
6) Aniline	6.41	93	318799	10.10	ng	# 81
8) 2-Chlorophenol	6.53	128	403493	17.82	ng	97
9) Benzaldehyde	6.29	77	267142	15.70	ng	93
10) Phenol	6.41	94	423557	15.29	ng	100
11) bis(2-Chloroethyl)ether	6.49	93	385500	17.50	ng	97
12) 1,3-Dichlorobenzene	6.68	146	399665m	16.54	ng	
13) 1,4-Dichlorobenzene	6.76	146	414373	16.84	ng	98
14) 1,2-Dichlorobenzene	6.92	146	342562	16.05	ng	95
15) Benzyl Alcohol	6.89	79	307231	16.27	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.02	45	599950	18.28	ng	88
17) 2-Methylphenol	7.01	107	319166	17.53	ng	98
18) Hexachloroethane	7.25	117	149754	16.42	ng	# 86
19) n-Nitroso-di-n-propylamine	7.16	70	270899	16.76	ng	99
20) 3+4-Methylphenols	7.17	107	413838	19.05	ng	# 73
22) Acetophenone	7.16	105	578822	17.28	ng	# 92
24) Nitrobenzene	7.33	77	448860	17.26	ng	92
25) Isophorone	7.57	82	746433	16.57	ng	98
26) 2-Nitrophenol	7.65	139	210967	16.45	ng	# 82
27) 2,4-Dimethylphenol	7.70	122	372930	17.53	ng	94
28) bis(2-Chloroethoxy)methane	7.79	93	478188	17.50	ng	97
29) 2,4-Dichlorophenol	7.89	162	297533	16.52	ng	89
30) 1,2,4-Trichlorobenzene	7.97	180	326504	16.54	ng	95
31) Naphthalene	8.05	128	1141746	18.37	ng	99
32) Benzoic acid	7.80	122	218200	13.83	ng	98
33) 4-Chloroaniline	8.11	127	316640	11.30	ng	# 93
34) Hexachlorobutadiene	8.18	225	178332	17.12	ng	99
35) Caprolactam	8.45	113	92069	15.55	ng	# 58
36) 4-Chloro-3-methylphenol	8.60	107	324059	15.63	ng	# 81
37) 2-Methylnaphthalene	8.75	142	720473m	15.27	ng	
39) 1,2,4,5-Tetrachlorobenzene	8.91	216	297405	16.40	ng	98
40) Hexachlorocyclopentadiene	8.90	237	93020	14.92	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.02	196	203300	16.03	nq	95
43) 2,4,5-Trichlorophenol	9.07	196	206531	15.82	nq	# 92
45) 1,1'-Biphenyl	9.21	154	868461	17.67	nq	95
46) 2-Chloronaphthalene	9.23	162	655348	16.84	nq	96
47) 2-Nitroaniline	9.33	65	233276	16.83	nq	# 76
48) Acenaphthylene	9.64	152	1025060	17.12	nq	99
49) Dimethylphthalate	9.52	163	764521	16.65	nq	99
50) 2,6-Dinitrotoluene	9.57	165	175762	17.49	nq	90
51) Acenaphthene	9.81	154	675208	16.64	nq	99
52) 3-Nitroaniline	9.73	138	147694	11.88	nq	92
53) 2,4-Dinitrophenol	9.85	184	68114	12.18	nq	# 73
54) Dibenzofuran	9.98	168	821682	14.99	nq	99
55) 4-Nitrophenol	9.92	139	132265	15.67	nq	# 75
56) 2,4-Dinitrotoluene	9.97	165	207819	16.48	nq	97
57) Fluorene	10.33	166	651337	16.34	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.11	232	156669	15.18	nq	99
59) Diethylphthalate	10.21	149	707600	15.87	nq	97
60) 4-Chlorophenyl-phenylether	10.33	204	303005	17.36	nq	# 85
61) 4-Nitroaniline	10.35	138	199813	15.51	nq	89
62) Azobenzene	10.49	77	735345	14.98	nq	89
64) 4,6-Dinitro-2-methylphenol	10.38	198	102884	18.62	nq	84
65) n-Nitrosodiphenylamine	10.44	169	533848	19.69	nq	99
66) 4-Bromophenyl-phenylether	10.81	248	152246	16.02	nq	# 88
67) Hexachlorobenzene	10.88	284	161292	15.88	nq	# 88
68) Atrazine	10.97	200	99991	11.43	nq	97
69) Pentachlorophenol	11.07	266	66557	13.35	nq	98
70) Phenanthrene	11.29	178	818914	16.53	nq	98
71) Anthracene	11.33	178	786159	14.22	nq	98
72) Carbazole	11.49	167	719539	13.49	nq	99
73) Di-n-butylphthalate	11.84	149	903909	16.67	nq	# 97
74) Fluoranthene	12.48	202	838079	16.89	nq	91
76) Benzidine	12.60	184	237948	10.50	nq	99
77) Pyrene	12.69	202	829003	16.94	nq	98
79) Butylbenzylphthalate	13.32	149	347871	15.63	nq	97
80) Benzo(a)anthracene	13.88	228	668600	17.75	nq	99
81) 3,3'-Dichlorobenzidine	13.85	252	167178	11.71	nq	99
82) Chrysene	13.92	228	671599	17.78	nq	99
83) Bis(2-ethylhexyl)phthalate	13.89	149	444310	16.95	nq	# 95
84) Di-n-octyl phthalate	14.50	149	810122	16.68	nq	99
85) Indeno(1,2,3-cd)pyrene	16.63	276	589778	18.02	nq	99
87) Benzo(b)fluoranthene	14.90	252	673446m	15.94	nq	
88) Benzo(k)fluoranthene	14.92	252	602456m	16.73	nq	
89) Benzo(a)pyrene	15.23	252	604502	16.12	nq	97
90) Dibenzo(a,h)anthracene	16.64	278	499844	16.22	nq	97
91) Benzo(g,h,i)perylene	17.03	276	500727	15.63	nq	98

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

