

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101818\
 Data File : BG037442.D
 Acq On : 19 Oct 2018 6:27
 Operator : JU/SJ
 Sample : J5164-09
 Misc : MDL 5 PPM
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MDL-WATER-03-QT4-2018

Manual Integrations
 APPROVED

Sohil
 10/19/2018 3:57:54 PM

Quant Time: Oct 19 09:14:28 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 18 14:06:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	50576	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	238203	20.00	ng	0.00
39) Acenaphthene-d10	14.74	164	162086	20.00	ng	0.00
64) Phenanthrene-d10	17.48	188	395683	20.00	ng	0.00
76) Chrysene-d12	21.77	240	362178	20.00	ng	0.00
87) Perylene-d12	25.10	264	364857	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.64	112	424474	140.32	ng	0.00
7) Phenol-d6	7.25	99	622380	136.06	ng	0.00
23) Nitrobenzene-d5	9.28	82	343149	80.70	ng	0.00
42) 2,4,6-Tribromophenol	16.23	330	238142	134.71	ng	0.00
45) 2-Fluorobiphenyl	13.36	172	851217	79.19	ng	0.00
79) Terphenyl-d14	20.09	244	1294387	76.37	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	7442	4.851	ng	87
3) Pyridine	3.95	79	19492	5.041	ng	89
4) n-Nitrosodimethylamine	3.86	42	8419	6.113	ng	# 91
6) Aniline	7.43	93	19390	3.475	ng	96
8) 2-Chlorophenol	7.67	128	18072	5.107	ng	98
9) Benzaldehyde	7.25	77	14076	4.919	ng	97
10) Phenol	7.27	94	24063	4.986	ng	81
11) bis(2-Chloroethyl)ether	7.52	93	17544	4.786	ng	94
12) 1,3-Dichlorobenzene	7.99	146	18881	4.792	ng	98
13) 1,4-Dichlorobenzene	8.14	146	19813	4.916	ng	95
14) 1,2-Dichlorobenzene	8.46	146	18303	4.721	ng	92
15) Benzyl Alcohol	8.35	79	15471	4.577	ng	93
16) 2,2'-oxybis(1-Chloropropan	8.63	45	32538	4.961	ng	99
17) 2-Methylphenol	8.54	107	16165	5.066	ng	88
18) Hexachloroethane	9.19	117	7049	5.131	ng	95
19) n-Nitroso-di-n-propylamine	8.91	70	14662	4.655	ng	93
20) 3+4-Methylphenols	8.86	107	21959	4.813	ng	97
22) Acetophenone	8.93	105	30453	5.098	ng	# 94
24) Nitrobenzene	9.33	77	22110	5.005	ng	# 97
25) Isophorone	9.84	82	41217	5.305	ng	97
26) 2-Nitrophenol	10.04	139	8875	4.460	ng	98
27) 2,4-Dimethylphenol	10.07	122	19485	5.484	ng	91
28) bis(2-Chloroethoxy)methane	10.33	93	25518	5.013	ng	93
29) 2,4-Dichlorophenol	10.56	162	18286	4.622	ng	97
30) 1,2,4-Trichlorobenzene	10.79	180	19849	4.614	ng	97
31) Naphthalene	10.98	128	63134	5.396	ng	100
32) Benzoic acid	10.12	122	4674	2.025	ng	93
33) 4-Chloroaniline	11.08	127	19333	3.517	ng	94
34) Hexachlorobutadiene	11.24	225	11732	4.601	ng	95
35) Caprolactam	11.85	113	6105	3.848	ng	94
36) 4-Chloro-3-methylphenol	12.18	107	20069	4.498	ng	95
37) 2-Methylnaphthalene	12.58	142	46399	5.440	ng	93
38) 1-Methylnaphthalene	12.79	142	45603m	5.506	ng	
40) 1,2,4,5-Tetrachlorobenzene	12.94	216	24305	4.649	ng	98

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41) Hexachlorocyclopentadiene	12.91	237	11278	4.516	nq	94
43) 2,4,6-Trichlorophenol	13.17	196	15043	4.307	nq	100
44) 2,4,5-Trichlorophenol	13.24	196	17581	4.623	nq	96
46) 1,1'-Biphenyl	13.58	154	60404	4.500	nq	94
47) 2-Chloronaphthalene	13.62	162	48512	4.777	nq	99
48) 2-Nitroaniline	13.82	65	13610	4.419	nq	91
49) Acenaphthylene	14.47	152	80216	5.251	nq	94
50) Dimethylphthalate	14.19	163	63827	5.090	nq	97
51) 2,6-Dinitrotoluene	14.32	165	12889	4.646	nq	88
52) Acenaphthene	14.80	154	47617	5.061	nq	97
53) 3-Nitroaniline	14.65	138	11795	3.830	nq	91
54) 2,4-Dinitrophenol	14.85	184	658	7.145	nq	# 76
55) Dibenzofuran	15.13	168	82033	5.263	nq	97
56) 4-Nitrophenol	14.93	139	6793	2.980	nq	89
57) 2,4-Dinitrotoluene	15.10	165	15914	4.370	nq	97
58) Fluorene	15.79	166	64543	5.529	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.35	232	15053	4.623	nq	# 98
60) Diethylphthalate	15.54	149	65377	5.078	nq	97
61) 4-Chlorophenyl-phenylether	15.77	204	32989	4.849	nq	97
62) 4-Nitroaniline	15.80	138	13045	4.263	nq	# 82
63) Azobenzene	16.07	77	64843	5.279	nq	96
65) 4,6-Dinitro-2-methylphenol	15.86	198	3147	9.754	nq	# 73
66) n-Nitrosodiphenylamine	15.99	169	58285	4.897	nq	97
67) 4-Bromophenyl-phenylether	16.67	248	20655	4.753	nq	85
68) Hexachlorobenzene	16.78	284	21922	4.868	nq	95
69) Atrazine	16.93	200	18732	4.353	nq	95
70) Pentachlorophenol	17.12	266	6078	2.015	nq	94
71) Phenanthrene	17.52	178	111068	5.523	nq	98
72) Anthracene	17.62	178	109333	5.540	nq	98
73) Carbazole	17.89	167	93993	4.761	nq	95
74) Di-n-butylphthalate	18.43	149	106334	4.842	nq	98
75) Fluoranthene	19.53	202	123762	5.426	nq	99
77) Benzidine	19.72	184	15608	1.267	nq	96
78) Pyrene	19.89	202	125990	5.321	nq	96
80) Butylbenzylphthalate	20.77	149	42956	4.433	nq	90
81) Benzo(a)anthracene	21.75	228	114979	5.367	nq	98
82) 3,3'-Dichlorobenzidine	21.66	252	26735	3.220	nq	99
83) Chrysene	21.81	228	109431	5.312	nq	95
84) Bis(2-ethylhexyl)phthalate	21.63	149	60443	4.450	nq	97
85) Di-n-octyl phthalate	22.88	149	98548	4.450	nq	98
86) Indeno(1,2,3-cd)pyrene	28.92	276	92151	4.171	nq	# 92
88) Benzo(b)fluoranthene	24.03	252	103365	5.168	nq	100
89) Benzo(k)fluoranthene	24.10	252	106596m	5.439	nq	
90) Benzo(a)pyrene	24.94	252	98413	5.178	nq	96
91) Dibenzo(a,h)anthracene	29.01	278	74104	4.227	nq	98
92) Benzo(g,h,i)perylene	30.12	276	82577	4.655	nq	98

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

