

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050118\
 Data File : BG034100.D
 Acq On : 2 May 2018 4:46
 Operator : SJ/JU
 Sample : J2133-08
 Misc : 20PPM L00
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOQ-WATER-02-QT2-2018

Manual Integrations
 APPROVED

Sohil
 5/2/2018 4:50:54 PM

Quant Time: May 02 08:03:14 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 01 19:08:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	59089	20.00	ng	0.00
21) Naphthalene-d8	10.89	136	272138	20.00	ng	0.00
38) Acenaphthene-d10	14.72	164	207683	20.00	ng	0.00
63) Phenanthrene-d10	17.47	188	565579	20.00	ng	0.00
75) Chrysene-d12	21.76	240	598657	20.00	ng	0.00
86) Perylene-d12	25.05	264	600386	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.63	112	499012	136.44	ng	0.00
7) Phenol-d6	7.22	99	749707	131.67	ng	0.00
23) Nitrobenzene-d5	9.24	82	578606	81.58	ng	0.00
41) 2,4,6-Tribromophenol	16.20	330	371635	129.23	ng	0.00
44) 2-Fluorobiphenyl	13.34	172	1192732	83.69	ng	0.00
78) Terphenyl-d14	20.09	244	2222539	81.33	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.56	88	35440	23.076	ng	# 78
3) Pyridine	3.95	79	110266	22.882	ng	# 84
4) n-Nitrosodimethylamine	3.86	42	56258	21.411	ng	# 68
6) Aniline	7.39	93	163161	21.068	ng	97
8) 2-Chlorophenol	7.64	128	75825	20.777	ng	93
9) Benzaldehyde	7.21	77	69076	18.249	ng	92
10) Phenol	7.24	94	122294	21.295	ng	# 71
11) bis(2-Chloroethyl)ether	7.49	93	94532	20.909	ng	89
12) 1,3-Dichlorobenzene	7.96	146	97082	21.027	ng	# 87
13) 1,4-Dichlorobenzene	8.11	146	99330	21.662	ng	92
14) 1,2-Dichlorobenzene	8.43	146	94447	21.685	ng	96
15) Benzyl Alcohol	8.30	79	130360	21.463	ng	# 83
16) 2,2'-oxybis(1-Chloropropan	8.60	45	135057	21.712	ng	90
17) 2-Methylphenol	8.50	107	78652	20.814	ng	# 67
18) Hexachloroethane	9.16	117	38755	19.868	ng	86
19) n-Nitroso-di-n-propylamine	8.87	70	109703	21.455	ng	97
20) 3+4-Methylphenols	8.83	107	111156	19.862	ng	# 71
22) Acetophenone	8.89	105	176555	19.046	ng	# 97
24) Nitrobenzene	9.28	77	150845	19.429	ng	# 85
25) Isophorone	9.80	82	273398	18.895	ng	95
26) 2-Nitrophenol	9.99	139	43165	20.084	ng	# 62
27) 2,4-Dimethylphenol	10.04	122	77689	18.645	ng	91
28) bis(2-Chloroethoxy)methane	10.29	93	137624	19.364	ng	99
29) 2,4-Dichlorophenol	10.52	162	92058	18.484	ng	92
30) 1,2,4-Trichlorobenzene	10.75	180	113405	18.910	ng	97
31) Naphthalene	10.94	128	277137	19.375	ng	96
32) Benzoic acid	10.14	122	56707	18.366	ng	# 67
33) 4-Chloroaniline	11.04	127	124437	19.292	ng	# 84
34) Hexachlorobutadiene	11.23	225	100480	19.795	ng	95
35) Caprolactam	11.79	113	34725	19.629	ng	94
36) 4-Chloro-3-methylphenol	12.15	107	125943	19.502	ng	95
37) 2-Methylnaphthalene	12.55	142	217327	18.738	ng	# 90
39) 1,2,4,5-Tetrachlorobenzene	12.91	216	162248	20.189	ng	98
40) Hexachlorocyclopentadiene	12.89	237	90912	19.092	ng	90

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.14	196	91554	19.608	nq	98
43) 2,4,5-Trichlorophenol	13.21	196	107492	20.824	nq	94
45) 1,1'-Biphenyl	13.55	154	323710	19.952	nq	99
46) 2-Chloronaphthalene	13.59	162	251578	20.482	nq	96
47) 2-Nitroaniline	13.79	65	99725	20.307	nq	87
48) Acenaphthylene	14.44	152	391316	19.954	nq	97
49) Dimethylphthalate	14.17	163	360328	20.008	nq #	99
50) 2,6-Dinitrotoluene	14.28	165	69874	20.202	nq #	81
51) Acenaphthene	14.78	154	270437	20.290	nq	100
52) 3-Nitroaniline	14.61	138	70770	20.300	nq #	80
53) 2,4-Dinitrophenol	14.81	184	28968	21.409	nq #	77
54) Dibenzofuran	15.11	168	389406	20.153	nq #	91
55) 4-Nitrophenol	14.90	139	49789	18.711	nq #	48
56) 2,4-Dinitrotoluene	15.07	165	101304	21.570	nq	98
57) Fluorene	15.77	166	336902	20.004	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.34	232	102208	19.140	nq	95
59) Diethylphthalate	15.54	149	383229	20.109	nq	95
60) 4-Chlorophenyl-phenylether	15.76	204	205645	20.525	nq	93
61) 4-Nitroaniline	15.77	138	76562	20.704	nq #	52
62) Azobenzene	16.05	77	412360	20.030	nq	93
64) 4,6-Dinitro-2-methylphenol	15.83	198	53657	20.178	nq	80
65) n-Nitrosodiphenylamine	15.97	169	297240	19.274	nq	98
66) 4-Bromophenyl-phenylether	16.66	248	132585	18.881	nq	93
67) Hexachlorobenzene	16.77	284	139971	19.637	nq	93
68) Atrazine	16.92	200	90885	11.190	nq	96
69) Pentachlorophenol	17.11	266	95149	19.152	nq	94
70) Phenanthrene	17.51	178	576497	19.837	nq	100
71) Anthracene	17.60	178	573066	19.456	nq	98
72) Carbazole	17.87	167	531424	19.709	nq	98
73) Di-n-butylphthalate	18.44	149	630523	19.428	nq #	96
74) Fluoranthene	19.52	202	735533	20.029	nq	96
76) Benzidine	19.70	184	264985	17.184	nq #	96
77) Pyrene	19.88	202	749678	19.987	nq	94
79) Butylbenzylphthalate	20.78	149	285565	19.503	nq	93
80) Benzo(a)anthracene	21.74	228	756232	19.711	nq	100
81) 3,3'-Dichlorobenzidine	21.65	252	277679	19.139	nq	98
82) Chrysene	21.81	228	722758	19.682	nq	98
83) Bis(2-ethylhexyl)phthalate	21.67	149	393376	19.521	nq	97
84) Di-n-octyl phthalate	22.92	149	627257	19.489	nq	99
85) Indeno(1,2,3-cd)pyrene	28.80	276	799808m	19.791	nq	
87) Benzo(b)fluoranthene	24.00	252	729788	19.243	nq #	96
88) Benzo(k)fluoranthene	24.07	252	716351	19.763	nq #	97
89) Benzo(a)pyrene	24.89	252	701203	19.762	nq #	95
90) Dibenzo(a,h)anthracene	28.88	278	654911	19.580	nq #	92
91) Benzo(g,h,i)perylene	29.97	276	635936	19.290	nq #	91

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

