

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060819\
 Data File : BG041131.D
 Acq On : 8 Jun 2019 13:36
 Operator : HP/JU
 Sample : K3153-02MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 HR-05-060519-AMS

Quant Time: Jun 09 23:18:19 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jun 09 23:08:34 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	40747	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	168036	20.00	ng	0.00
39) Acenaphthene-d10	14.74	164	129404	20.00	ng	0.00
64) Phenanthrene-d10	17.48	188	334786	20.00	ng	0.00
76) Chrysene-d12	21.76	240	342440	20.00	ng	0.00
87) Perylene-d12	25.09	264	364370	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.64	112	307239	135.97	ng	0.00
7) Phenol-d6	7.25	99	418101	119.80	ng	0.00
23) Nitrobenzene-d5	9.28	82	343201	90.67	ng	0.00
42) 2,4,6-Tribromophenol	16.23	330	282761	147.19	ng	0.00
45) 2-Fluorobiphenyl	13.36	172	818972	91.14	ng	0.00
79) Terphenyl-d14	20.08	244	1508469	93.49	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.51	88	34178	33.331	ng	# 15
3) Pyridine	3.92	79	92587	30.515	ng	99
4) n-Nitrosodimethylamine	3.84	42	58060	41.231	ng	85
6) Aniline	7.43	93	65221	14.001	ng	97
8) 2-Chlorophenol	7.67	128	119332	44.296	ng	93
9) Benzaldehyde	7.24	77	40245	19.332	ng	90
10) Phenol	7.28	94	139569	37.300	ng	98
11) bis(2-Chloroethyl)ether	7.53	93	109344	40.281	ng	98
12) 1,3-Dichlorobenzene	7.99	146	119338	37.089	ng	97
13) 1,4-Dichlorobenzene	8.14	146	120423	36.975	ng	95
14) 1,2-Dichlorobenzene	8.46	146	118814	37.572	ng	99
15) Benzyl Alcohol	8.35	79	139686	43.270	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.62	45	169024	38.106	ng	98
17) 2-Methylphenol	8.54	107	110859	40.919	ng	98
18) Hexachloroethane	9.19	117	47028	37.464	ng	96
19) n-Nitroso-di-n-propylamine	8.91	70	106785	39.762	ng	94
20) 3+4-Methylphenols	8.87	107	149827	39.924	ng	96
22) Acetophenone	8.94	105	209870	44.523	ng	# 98
24) Nitrobenzene	9.33	77	171072	44.349	ng	97
25) Isophorone	9.84	82	313969	43.586	ng	100
26) 2-Nitrophenol	10.03	139	72035	45.299	ng	86
27) 2,4-Dimethylphenol	10.08	122	121782	46.370	ng	95
28) bis(2-Chloroethoxy)methane	10.32	93	161459	41.529	ng	98
29) 2,4-Dichlorophenol	10.57	162	149204	44.737	ng	94
30) 1,2,4-Trichlorobenzene	10.79	180	152164	40.107	ng	98
31) Naphthalene	10.98	128	381383	42.272	ng	97
32) Benzoic acid	10.22	122	66949	35.361	ng	94
33) 4-Chloroaniline	11.10	127	15740	3.760	ng	97
34) Hexachlorobutadiene	11.24	225	108931	37.524	ng	96
35) Caprolactam	11.88	113	37221m	33.796	ng	
36) 4-Chloro-3-methylphenol	12.20	107	175717	46.133	ng	97
37) 2-Methylnaphthalene	12.57	142	306463	44.104	ng	98
38) 1-Methylnaphthalene	12.79	142	289816m	44.261	ng	
40) 1,2,4,5-Tetrachlorobenzene	12.93	216	220069	42.194	ng	99

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41) Hexachlorocyclopentadiene	12.90	237	247770	87.511	nq	100
43) 2,4,6-Trichlorophenol	13.17	196	151631	44.940	nq	97
44) 2,4,5-Trichlorophenol	13.25	196	158434	44.523	nq	97
46) 1,1'-Biphenyl	13.57	154	420317	43.880	nq	97
47) 2-Chloronaphthalene	13.62	162	346694	42.160	nq	99
48) 2-Nitroaniline	13.83	65	128873	43.685	nq	97
49) Acenaphthylene	14.46	152	554576	44.809	nq	99
50) Dimethylphthalate	14.19	163	480628	43.877	nq	98
51) 2,6-Dinitrotoluene	14.32	165	107237	45.405	nq	97
52) Acenaphthene	14.80	154	324073	43.224	nq	96
53) 3-Nitroaniline	14.65	138	37726	15.974	nq	95
54) 2,4-Dinitrophenol	14.85	184	124582	94.088	nq	95
55) Dibenzofuran	15.13	168	567319	44.969	nq	99
56) 4-Nitrophenol	14.95	139	143919	88.843	nq	94
57) 2,4-Dinitrotoluene	15.10	165	157446	47.193	nq	96
58) Fluorene	15.78	166	455568	45.344	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.35	232	166624	47.647	nq	99
60) Diethylphthalate	15.54	149	488269	43.677	nq	99
61) 4-Chlorophenyl-phenylether	15.77	204	284810	43.613	nq	94
62) 4-Nitroaniline	15.82	138	102444	40.973	nq	95
63) Azobenzene	16.06	77	452887	44.574	nq	99
65) 4,6-Dinitro-2-methylphenol	15.86	198	90525	50.020	nq	93
66) n-Nitrosodiphenylamine	15.99	169	419180	44.540	nq	98
67) 4-Bromophenyl-phenylether	16.66	248	189853	42.021	nq	99
68) Hexachlorobenzene	16.77	284	196400	40.823	nq	98
69) Atrazine	16.93	200	182429	50.237	nq	98
70) Pentachlorophenol	17.12	266	226733	87.550	nq	97
71) Phenanthrene	17.53	178	781960	44.841	nq	98
72) Anthracene	17.61	178	800934	46.569	nq	99
73) Carbazole	17.89	167	699396	47.393	nq	100
74) Di-n-butylphthalate	18.42	149	817683	44.580	nq	99
75) Fluoranthene	19.53	202	1022360	47.465	nq	99
77) Benzidine	19.71	184	31099	3.807	nq	93
78) Pyrene	19.89	202	1026000	46.090	nq	99
80) Butylbenzylphthalate	20.75	149	391868	45.235	nq	100
81) Benzo(a)anthracene	21.74	228	1011074	44.355	nq	99
82) 3,3'-Dichlorobenzidine	21.65	252	143313	17.875	nq	99
83) Chrysene	21.81	228	985281	45.168	nq	98
84) Bis(2-ethylhexyl)phthalate	21.61	149	536918	44.028	nq	99
85) Di-n-octyl phthalate	22.84	149	926752	45.630	nq	99
86) Indeno(1,2,3-cd)pyrene	28.91	276	1210087	45.285	nq	# 99
88) Benzo(b)fluoranthene	24.02	252	1023079	46.293	nq	98
89) Benzo(k)fluoranthene	24.09	252	1005842	45.992	nq	99
90) Benzo(a)pyrene	24.93	252	1000089	47.431	nq	96
91) Dibenzo(a,h)anthracene	28.98	278	996161	47.282	nq	100
92) Benzo(g,h,i)perylene	30.12	276	1002464	48.975	nq	98

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

