

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111317\
 Data File : BG030974.D
 Acq On : 13 Nov 2017 15:46
 Operator : SJ/JU
 Sample : PB104186BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB104186BS

Quant Time: Nov 14 04:50:09 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 10 15:38:25 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	52908	20.00	ng	0.00
21) Naphthalene-d8	10.96	136	224447	20.00	ng	0.00
38) Acenaphthene-d10	14.76	164	157748	20.00	ng	-0.01
63) Phenanthrene-d10	17.50	188	409181	20.00	ng	-0.02
75) Chrysene-d12	21.77	240	485648	20.00	ng	-0.02
86) Perylene-d12	25.07	264	493777	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.71	112	281781	91.33	ng	0.00
7) Phenol-d6	7.31	99	364530	87.70	ng	0.00
23) Nitrobenzene-d5	9.31	82	213247	56.30	ng	-0.02
41) 2,4,6-Tribromophenol	16.25	330	242315	94.96	ng	0.00
44) 2-Fluorobiphenyl	13.38	172	653141	59.49	ng	-0.02
78) Terphenyl-d14	20.09	244	1261973	57.38	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.55	88	29285	23.389	ng	# 82
3) Pyridine	3.96	79	82621	22.729	ng	90
4) n-Nitrosodimethylamine	3.87	42	42299	24.553	ng	# 93
6) Aniline	7.46	93	76596	14.002	ng	97
8) 2-Chlorophenol	7.71	128	95348	28.003	ng	86
9) Benzaldehyde	7.27	77	43849	15.074	ng	91
10) Phenol	7.33	94	123759	28.669	ng	97
11) bis(2-Chloroethyl)ether	7.56	93	81821	23.739	ng	88
12) 1,3-Dichlorobenzene	8.03	146	103241	25.843	ng	96
13) 1,4-Dichlorobenzene	8.18	146	106099	26.209	ng	96
14) 1,2-Dichlorobenzene	8.50	146	100557	25.880	ng	97
15) Benzyl Alcohol	8.38	79	82832	24.843	ng	93
16) 2,2'-oxybis(1-Chloropropan	8.67	45	121959	32.034	ng	93
17) 2-Methylphenol	8.59	107	82709	27.514	ng	96
18) Hexachloroethane	9.23	117	36366	25.810	ng	92
19) n-Nitroso-di-n-propylamine	8.94	70	68174	21.570	ng	# 93
20) 3+4-Methylphenols	8.92	107	114340	26.750	ng	95
22) Acetophenone	8.97	105	138068	24.706	ng	# 93
24) Nitrobenzene	9.35	77	102424	26.472	ng	89
25) Isophorone	9.87	82	192950	26.120	ng	# 91
26) 2-Nitrophenol	10.06	139	54337	26.942	ng	# 87
27) 2,4-Dimethylphenol	10.12	122	95122	31.770	ng	95
28) bis(2-Chloroethoxy)methane	10.35	93	119089	25.597	ng	95
29) 2,4-Dichlorophenol	10.61	162	112004	31.152	ng	91
30) 1,2,4-Trichlorobenzene	10.82	180	120839	28.151	ng	95
31) Naphthalene	11.01	128	289198	26.211	ng	98
32) Benzoic acid	10.26	122	47598	22.497	ng	# 75
33) 4-Chloroaniline	11.12	127	53263	11.027	ng	98
34) Hexachlorobutadiene	11.29	225	84177	28.276	ng	97
35) Caprolactam	11.88	113	36753	25.024	ng	# 74
36) 4-Chloro-3-methylphenol	12.23	107	111549	28.506	ng	89
37) 2-Methylnaphthalene	12.60	142	229485	26.942	ng	92
39) 1,2,4,5-Tetrachlorobenzene	12.97	216	157961	25.230	ng	96
40) Hexachlorocyclopentadiene	12.94	237	130701	59.582	ng	90

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42) 2,4,6-Trichlorophenol	13.20	196	106557	29.771	nq	98
43) 2,4,5-Trichlorophenol	13.29	196	115008	29.368	nq #	92
45) 1,1'-Biphenyl	13.60	154	323817	24.755	nq	98
46) 2-Chloronaphthalene	13.64	162	261520	27.709	nq	95
47) 2-Nitroaniline	13.84	65	71841	25.681	nq #	82
48) Acenaphthylene	14.48	152	401546	25.995	nq	98
49) Dimethylphthalate	14.21	163	359317	26.344	nq	99
50) 2,6-Dinitrotoluene	14.33	165	78686	27.989	nq #	78
51) Acenaphthene	14.82	154	245669	25.465	nq	97
52) 3-Nitroaniline	14.67	138	40057	13.419	nq #	74
53) 2,4-Dinitrophenol	14.88	184	69221	47.340	nq #	47
54) Dibenzofuran	15.16	168	419507	27.300	nq	99
55) 4-Nitrophenol	14.98	139	111636	52.500	nq #	87
56) 2,4-Dinitrotoluene	15.12	165	113724	27.982	nq #	72
57) Fluorene	15.81	166	332666	26.665	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.39	232	116793	31.292	nq #	85
59) Diethylphthalate	15.56	149	352866	25.469	nq	97
60) 4-Chlorophenyl-phenylether	15.79	204	207730	27.570	nq	91
61) 4-Nitroaniline	15.82	138	77625	24.558	nq	89
62) Azobenzene	16.08	77	275849	26.107	nq	94
64) 4,6-Dinitro-2-methylphenol	15.89	198	60352	22.190	nq #	77
65) n-Nitrosodiphenylamine	16.01	169	310294	25.025	nq	98
66) 4-Bromophenyl-phenylether	16.68	248	148670	28.687	nq	93
67) Hexachlorobenzene	16.81	284	155709	28.277	nq #	92
68) Atrazine	16.94	200	133634	26.122	nq	96
69) Pentachlorophenol	17.16	266	154638	50.802	nq	98
70) Phenanthrene	17.54	178	580425	27.244	nq	99
71) Anthracene	17.63	178	588144	27.551	nq	99
72) Carbazole	17.90	167	523662	26.180	nq	99
73) Di-n-butylphthalate	18.44	149	606906	24.712	nq	99
74) Fluoranthene	19.54	202	769993	28.545	nq	97
76) Benzidine	19.71	184	248803	15.949	nq	98
77) Pyrene	19.90	202	788538	27.363	nq	99
79) Butylbenzylphthalate	20.76	149	283358	24.661	nq #	84
80) Benzo(a)anthracene	21.75	228	815494	27.033	nq	99
81) 3,3'-Dichlorobenzidine	21.66	252	165125	13.560	nq #	96
82) Chrysene	21.82	228	761409	27.286	nq	97
83) Bis(2-ethylhexyl)phthalate	21.63	149	396015	23.876	nq #	98
84) Di-n-octyl phthalate	22.86	149	666890	24.703	nq #	93
85) Indeno(1,2,3-cd)pyrene	28.84	276	922876	27.204	nq #	94
87) Benzo(b)fluoranthene	24.01	252	808064	27.054	nq #	98
88) Benzo(k)fluoranthene	24.08	252	794551	26.838	nq #	97
89) Benzo(a)pyrene	24.91	252	783126	27.599	nq #	95
90) Dibenzo(a,h)anthracene	28.90	278	773788	26.680	nq #	96
91) Benzo(g,h,i)perylene	30.02	276	772929	27.459	nq #	95

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

