

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050919\
 Data File : BG040825.D
 Acq On : 9 May 2019 21:03
 Operator : HP/JU
 Sample : PB119594BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB119594BS

Quant Time: May 10 06:23:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG042319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 24 05:35:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	48567	20.00	ng	-0.02
21) Naphthalene-d8	10.96	136	200385	20.00	ng	-0.02
39) Acenaphthene-d10	14.77	164	146801	20.00	ng	-0.02
64) Phenanthrene-d10	17.51	188	415633	20.00	ng	-0.02
76) Chrysene-d12	21.80	240	422626	20.00	ng	-0.03
87) Perylene-d12	25.15	264	426480	20.00	ng	-0.05

System Monitoring Compounds

5) 2-Fluorophenol	5.67	112	369506	134.11	ng	-0.01
7) Phenol-d6	7.28	99	545695	128.64	ng	-0.02
23) Nitrobenzene-d5	9.32	82	376212	86.75	ng	-0.02
42) 2,4,6-Tribromophenol	16.26	330	307203	152.25	ng	-0.02
45) 2-Fluorobiphenyl	13.40	172	850897	83.71	ng	-0.02
79) Terphenyl-d14	20.11	244	1698467	89.03	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.54	88	42828	34.180	ng	93
3) Pyridine	3.95	79	123801	34.088	ng	99
4) n-Nitrosodimethylamine	3.87	42	70435	34.964	ng	96
6) Aniline	7.46	93	179027	31.840	ng	98
8) 2-Chlorophenol	7.70	128	135176	40.181	ng	99
9) Benzaldehyde	7.28	77	75076	26.524	ng	96
10) Phenol	7.31	94	189250	41.501	ng	97
11) bis(2-Chloroethyl)ether	7.56	93	124539	36.420	ng	99
12) 1,3-Dichlorobenzene	8.03	146	139064	37.872	ng	96
13) 1,4-Dichlorobenzene	8.18	146	140571	37.764	ng	99
14) 1,2-Dichlorobenzene	8.50	146	139283	38.800	ng	98
15) Benzyl Alcohol	8.38	79	156493	35.454	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.66	45	213649	31.381	ng	97
17) 2-Methylphenol	8.57	107	128890	39.939	ng	93
18) Hexachloroethane	9.23	117	55133	36.107	ng	98
19) n-Nitroso-di-n-propylamine	8.95	70	121818	31.900	ng	94
20) 3+4-Methylphenols	8.90	107	181394	40.349	ng	98
22) Acetophenone	8.97	105	227166	40.768	ng	# 96
24) Nitrobenzene	9.36	77	189787	41.017	ng	95
25) Isophorone	9.88	82	351361	36.372	ng	99
26) 2-Nitrophenol	10.07	139	78222	47.161	ng	95
27) 2,4-Dimethylphenol	10.11	122	145939	46.896	ng	98
28) bis(2-Chloroethoxy)methane	10.36	93	180429	37.227	ng	98
29) 2,4-Dichlorophenol	10.60	162	158570	44.820	ng	96
30) 1,2,4-Trichlorobenzene	10.82	180	169368	43.169	ng	96
31) Naphthalene	11.02	128	423270	39.759	ng	99
32) Benzoic acid	10.22	122	72217	33.898	ng	84
33) 4-Chloroaniline	11.12	127	142196	30.126	ng	99
34) Hexachlorobutadiene	11.28	225	120666	41.659	ng	98
35) Caprolactam	11.91	113	57903	41.454	ng	91
36) 4-Chloro-3-methylphenol	12.22	107	197902	44.341	ng	99
37) 2-Methylnaphthalene	12.61	142	326683	41.042	ng	99
38) 1-Methylnaphthalene	12.83	142	312826	40.208	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.97	216	229584	41.981	ng	99

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41) Hexachlorocyclopentadiene	12.94	237	280028	86.777	ng	97
43) 2,4,6-Trichlorophenol	13.20	196	159245	48.186	ng	99
44) 2,4,5-Trichlorophenol	13.27	196	177970	49.435	ng	99
46) 1,1'-Biphenyl	13.61	154	444724	39.019	ng	98
47) 2-Chloronaphthalene	13.65	162	374624	41.999	ng	97
48) 2-Nitroaniline	13.86	65	155453	48.960	ng	90
49) Acenaphthylene	14.50	152	604240	39.927	ng	99
50) Dimethylphthalate	14.22	163	521803	42.483	ng	98
51) 2,6-Dinitrotoluene	14.35	165	117169	48.878	ng	92
52) Acenaphthene	14.84	154	350141	39.729	ng	96
53) 3-Nitroaniline	14.68	138	101815	41.453	ng	# 92
54) 2,4-Dinitrophenol	14.88	184	64470	50.519	ng	90
55) Dibenzofuran	15.17	168	619888	42.942	ng	99
56) 4-Nitrophenol	14.96	139	179885	84.462	ng	95
57) 2,4-Dinitrotoluene	15.13	165	174309	53.512	ng	93
58) Fluorene	15.82	166	498386	42.716	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.39	232	185414	51.170	ng	97
60) Diethylphthalate	15.57	149	549105	42.372	ng	99
61) 4-Chlorophenyl-phenylether	15.80	204	306091	45.107	ng	97
62) 4-Nitroaniline	15.85	138	128618	46.812	ng	91
63) Azobenzene	16.09	77	526343	39.438	ng	98
65) 4,6-Dinitro-2-methylphenol	15.89	198	69149	35.871	ng	92
66) n-Nitrosodiphenylamine	16.02	169	465710	38.084	ng	99
67) 4-Bromophenyl-phenylether	16.70	248	208149	40.197	ng	94
68) Hexachlorobenzene	16.81	284	214230	40.011	ng	95
69) Atrazine	16.96	200	204017	42.110	ng	96
70) Pentachlorophenol	17.15	266	256256	81.788	ng	98
71) Phenanthrene	17.56	178	896552	40.048	ng	98
72) Anthracene	17.65	178	921042	40.930	ng	99
73) Carbazole	17.92	167	807680	39.396	ng	98
74) Di-n-butylphthalate	18.45	149	974188	39.369	ng	99
75) Fluoranthene	19.56	202	1190739	42.682	ng	99
77) Benzidine	19.74	184	449628	38.856	ng	99
78) Pyrene	19.92	202	1162688	43.894	ng	98
80) Butylbenzylphthalate	20.79	149	443007	42.911	ng	95
81) Benzo(a)anthracene	21.78	228	1133315	42.430	ng	99
82) 3,3'-Dichlorobenzidine	21.69	252	337623	35.464	ng	100
83) Chrysene	21.85	228	1126300	44.339	ng	99
84) Bis(2-ethylhexyl)phthalate	21.65	149	618111	41.477	ng	97
85) Di-n-octyl phthalate	22.90	149	1047343	41.730	ng	96
86) Indeno(1,2,3-cd)pyrene	29.00	276	1002246	32.633	ng	# 91
88) Benzo(b)fluoranthene	24.08	252	1110216	42.600	ng	99
89) Benzo(k)fluoranthene	24.15	252	1131454	46.487	ng	98
90) Benzo(a)pyrene	24.99	252	1068682	43.320	ng	99
91) Dibenzo(a,h)anthracene	29.08	278	796744	31.915	ng	97
92) Benzo(g,h,i)perylene	30.22	276	805495	32.420	ng	98

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

