

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG020720\
 Data File : BG044331.D
 Acq On : 7 Feb 2020 10:54
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Feb 08 02:29:50 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 30 16:05:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	104802	20.00	ng	0.00
21) Naphthalene-d8	10.71	136	423533	20.00	ng	0.00
39) Acenaphthene-d10	14.55	164	267738	20.00	ng	0.00
64) Phenanthrene-d10	17.29	188	566065	20.00	ng	0.00
76) Chrysene-d12	21.55	240	497021	20.00	ng	0.00
87) Perylene-d12	24.64	264	579030	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	485341	81.73	ng	0.00
7) Phenol-d6	7.09	99	647243	81.17	ng	0.00
23) Nitrobenzene-d5	9.07	82	592224	78.94	ng	0.00
42) 2,4,6-Tribromophenol	16.04	330	254485	80.97	ng	0.00
45) 2-Fluorobiphenyl	13.18	172	1282801	80.23	ng	0.00
79) Terphenyl-d14	19.91	244	1842694	76.26	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.39	88	104011	38.984	ng	99
3) Pyridine	3.79	79	299254	42.099	ng	98
4) n-Nitrosodimethylamine	3.70	42	116681	39.282	ng	100
6) Aniline	7.24	93	394481	39.854	ng	100
8) 2-Chlorophenol	7.48	128	263892	40.435	ng	99
9) Benzaldehyde	7.05	77	173289	38.156	ng	97
10) Phenol	7.12	94	320501	40.611	ng	99
11) bis(2-Chloroethyl)ether	7.34	93	266416	39.487	ng	97
12) 1,3-Dichlorobenzene	7.81	146	307741	39.493	ng	97
13) 1,4-Dichlorobenzene	7.95	146	307882	39.893	ng	98
14) 1,2-Dichlorobenzene	8.27	146	295014	40.442	ng	100
15) Benzyl Alcohol	8.15	79	226295	40.084	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.45	45	363357	39.790	ng	99
17) 2-Methylphenol	8.36	107	225348	40.021	ng	98
18) Hexachloroethane	9.00	117	116459	40.212	ng	97
19) n-Nitroso-di-n-propylamine	8.72	70	207009	38.952	ng	98
20) 3+4-Methylphenols	8.69	107	310823	40.055	ng	99
22) Acetophenone	8.73	105	409495	39.744	ng	# 99
24) Nitrobenzene	9.12	77	288328	39.369	ng	99
25) Isophorone	9.64	82	549586	39.305	ng	99
26) 2-Nitrophenol	9.83	139	159182	41.888	ng	99
27) 2,4-Dimethylphenol	9.89	122	215330	40.141	ng	100
28) bis(2-Chloroethoxy)methane	10.12	93	367241	39.371	ng	99
29) 2,4-Dichlorophenol	10.37	162	261478	40.311	ng	99
30) 1,2,4-Trichlorobenzene	10.58	180	296949	39.925	ng	99
31) Naphthalene	10.77	128	821027	39.955	ng	100
32) Benzoic acid	10.06	122	137988	41.694	ng	93
33) 4-Chloroaniline	10.87	127	372270	39.834	ng	99
34) Hexachlorobutadiene	11.06	225	180541	39.449	ng	98
35) Caprolactam	11.64	113	96138	40.460	ng	98
36) 4-Chloro-3-methylphenol	12.01	107	268344	39.458	ng	100
37) 2-Methylnaphthalene	12.38	142	602366	39.482	ng	99
38) 1-Methylnaphthalene	12.59	142	569201	39.572	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.75	216	320280	39.990	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.74	237	178834	46.536	nq	99
43) 2,4,6-Trichlorophenol	12.99	196	217065	41.931	nq	99
44) 2,4,5-Trichlorophenol	13.06	196	218912	41.401	nq	99
46) 1,1'-Biphenyl	13.39	154	779342	40.355	nq	100
47) 2-Chloronaphthalene	13.43	162	615627	40.060	nq	99
48) 2-Nitroaniline	13.63	65	180374	39.771	nq	94
49) Acenaphthylene	14.27	152	912437	40.480	nq	100
50) Dimethylphthalate	14.01	163	762529	39.621	nq	100
51) 2,6-Dinitrotoluene	14.12	165	170101	39.640	nq	98
52) Acenaphthene	14.62	154	576496	39.459	nq	98
53) 3-Nitroaniline	14.45	138	192224	40.489	nq	100
54) 2,4-Dinitrophenol	14.66	184	98873	42.804	nq	100
55) Dibenzofuran	14.95	168	880607	39.810	nq	99
56) 4-Nitrophenol	14.77	139	148906	42.819	nq	97
57) 2,4-Dinitrotoluene	14.91	165	240565	39.998	nq	97
58) Fluorene	15.60	166	691699	39.701	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.18	232	197708	41.396	nq	98
60) Diethylphthalate	15.37	149	763094	39.792	nq	100
61) 4-Chlorophenyl-phenylether	15.60	204	369712	39.137	nq	97
62) 4-Nitroaniline	15.62	138	192027	40.479	nq	96
63) Azobenzene	15.88	77	661611	39.364	nq	99
65) 4,6-Dinitro-2-methylphenol	15.68	198	133089	42.593	nq	99
66) n-Nitrosodiphenylamine	15.81	169	637061	40.159	nq	99
67) 4-Bromophenyl-phenylether	16.49	248	247788	40.178	nq	98
68) Hexachlorobenzene	16.61	284	272017	39.369	nq	99
69) Atrazine	16.76	200	218976	40.273	nq	98
70) Pentachlorophenol	16.95	266	146905	43.536	nq	97
71) Phenanthrene	17.34	178	1108056	40.424	nq	100
72) Anthracene	17.43	178	1098139	40.522	nq	100
73) Carbazole	17.70	167	1036139	41.474	nq	99
74) Di-n-butylphthalate	18.26	149	1303282	42.214	nq	100
75) Fluoranthene	19.35	202	1304587	41.142	nq	99
77) Benzidine	19.53	184	590499	42.833	nq	99
78) Pyrene	19.71	202	1311411	41.086	nq	99
80) Butylbenzylphthalate	20.60	149	612642	42.118	nq	99
81) Benzo(a)anthracene	21.52	228	1266238	41.337	nq	100
82) 3,3'-Dichlorobenzidine	21.44	252	508771	42.795	nq	100
83) Chrysene	21.59	228	1215010	41.036	nq	99
84) Bis(2-ethylhexyl)phthalate	21.45	149	834356	41.674	nq	100
85) Di-n-octyl phthalate	22.62	149	1474666	42.570	nq	100
86) Indeno(1,2,3-cd)pyrene	28.16	276	1571894	40.692	nq	# 93
88) Benzo(b)fluoranthene	23.66	252	1364399	40.892	nq	100
89) Benzo(k)fluoranthene	23.73	252	1319138	40.565	nq	99
90) Benzo(a)pyrene	24.50	252	1292371	40.967	nq	98
91) Dibenzo(a,h)anthracene	28.21	278	1272906	40.771	nq	98
92) Benzo(g,h,i)perylene	29.25	276	1257611	40.238	nq	97

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

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