

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062220\
 Data File : BG045829.D
 Acq On : 22 Jun 2020 16:35
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Jun 22 17:13:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 15 18:34:19 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	45944	20.00	ng	0.00
21) Naphthalene-d8	10.73	136	178641	20.00	ng	0.00
39) Acenaphthene-d10	14.57	164	128380	20.00	ng	0.00
64) Phenanthrene-d10	17.32	188	298813	20.00	ng	0.00
76) Chrysene-d12	21.58	240	290477	20.00	ng	0.00
86) Perylene-d12	24.71	264	326819	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.53	112	240139	88.29	ng	-0.01
7) Phenol-d6	7.15	99	364805	88.18	ng	-0.02
23) Nitrobenzene-d5	9.08	82	338022	86.44	ng	0.00
42) 2,4,6-Tribromophenol	16.07	330	163456	84.33	ng	0.00
45) 2-Fluorobiphenyl	13.19	172	764518	87.51	ng	0.00
79) Terphenyl-d14	19.94	244	1259237	87.85	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.37	88	52972	42.524	ng	100
3) Pyridine	3.78	79	155952	42.248	ng	99
4) n-Nitrosodimethylamine	3.69	42	57581	46.471	ng	# 89
6) Aniline	7.25	93	206696	42.432	ng	98
8) 2-Chlorophenol	7.51	128	128454	44.123	ng	93
9) Benzaldehyde	7.05	77	103700	41.733	ng	100
10) Phenol	7.17	94	170900	42.634	ng	95
11) bis(2-Chloroethyl)ether	7.34	93	127956	42.025	ng	95
12) 1,3-Dichlorobenzene	7.81	146	151242	43.594	ng	99
13) 1,4-Dichlorobenzene	7.95	146	154141	44.481	ng	99
14) 1,2-Dichlorobenzene	8.28	146	146061	44.027	ng	98
15) Benzyl Alcohol	8.18	79	145185	45.349	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.45	45	119923	41.791	ng	81
17) 2-Methylphenol	8.41	107	130885	43.874	ng	96
18) Hexachloroethane	9.00	117	59112	44.913	ng	94
19) n-Nitroso-di-n-propylamine	8.73	70	117738	43.288	ng	# 95
20) 3+4-Methylphenols	8.74	107	171530	43.720	ng	# 77
22) Acetophenone	8.75	105	208405	42.504	ng	98
24) Nitrobenzene	9.13	77	177645	42.606	ng	99
25) Isophorone	9.65	82	323505	42.734	ng	98
26) 2-Nitrophenol	9.84	139	77775	44.774	ng	94
27) 2,4-Dimethylphenol	9.93	122	114451	43.156	ng	98
28) bis(2-Chloroethoxy)methane	10.13	93	167153	42.271	ng	98
29) 2,4-Dichlorophenol	10.41	162	136179	44.012	ng	96
30) 1,2,4-Trichlorobenzene	10.59	180	161780	44.166	ng	99
31) Naphthalene	10.78	128	418312	42.831	ng	99
32) Benzoic acid	10.13	122	79033	46.334	ng	94
33) 4-Chloroaniline	10.90	127	176952	43.265	ng	99
34) Hexachlorobutadiene	11.07	225	116856	44.857	ng	97
35) Caprolactam	11.67	113	45507	45.449	ng	92
36) 4-Chloro-3-methylphenol	12.06	107	152856	42.787	ng	97
37) 2-Methylnaphthalene	12.39	142	315262	43.349	ng	100
38) 1-Methylnaphthalene	12.60	142	295142	42.885	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.76	216	182826	43.250	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.74	237	83845	40.548	ng	93
43) 2,4,6-Trichlorophenol	13.02	196	119583	41.724	ng	97
44) 2,4,5-Trichlorophenol	13.11	196	135699	41.542	ng	98
46) 1,1'-Biphenyl	13.40	154	390299	42.464	ng	100
47) 2-Chloronaphthalene	13.45	162	315489	42.493	ng	98
48) 2-Nitroaniline	13.66	65	105952	42.901	ng	99
49) Acenaphthylene	14.29	152	508586	42.766	ng	99
50) Dimethylphthalate	14.03	163	424947	42.253	ng	99
51) 2,6-Dinitrotoluene	14.14	165	96007	42.502	ng	94
52) Acenaphthene	14.63	154	306284	42.489	ng	99
53) 3-Nitroaniline	14.49	138	93481	41.432	ng	97
54) 2,4-Dinitrophenol	14.70	184	48296	43.544	ng	97
55) Dibenzofuran	14.97	168	503063	42.959	ng	99
56) 4-Nitrophenol	14.87	139	59158	38.751	ng	97
57) 2,4-Dinitrotoluene	14.94	165	137605	42.517	ng	# 97
58) Fluorene	15.62	166	415226	42.642	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.21	232	114385	40.346	ng	95
60) Diethylphthalate	15.39	149	440811	42.875	ng	98
61) 4-Chlorophenyl-phenylether	15.61	204	239627	42.697	ng	99
62) 4-Nitroaniline	15.66	138	97416	42.069	ng	91
63) Azobenzene	15.91	77	416483	42.189	ng	99
65) 4,6-Dinitro-2-methylphenol	15.71	198	84929	45.991	ng	98
66) n-Nitrosodiphenylamine	15.83	169	358179	42.320	ng	100
67) 4-Bromophenyl-phenylether	16.51	248	167204	43.064	ng	95
68) Hexachlorobenzene	16.63	284	175549	43.602	ng	96
69) Atrazine	16.78	200	139511	42.479	ng	97
70) Pentachlorophenol	16.99	266	67069	38.188	ng	96
71) Phenanthrene	17.36	178	653739	42.420	ng	99
72) Anthracene	17.45	178	662032	43.140	ng	99
73) Carbazole	17.73	167	632274	43.270	ng	99
74) Di-n-butylphthalate	18.29	149	758428	43.836	ng	100
75) Fluoranthene	19.37	202	832722	43.942	ng	100
77) Benzidine	19.56	184	273038	36.157	ng	99
78) Pyrene	19.74	202	843032	43.340	ng	100
80) Butylbenzylphthalate	20.62	149	352351	43.277	ng	95
81) Benzo(a)anthracene	21.56	228	820391	43.562	ng	100
82) 3,3'-Dichlorobenzidine	21.48	252	302481	42.462	ng	97
83) Chrysene	21.62	228	785352	43.115	ng	100
84) Bis(2-ethylhexyl)phthalate	21.47	149	488947	43.174	ng	98
85) Di-n-octyl phthalate	22.65	149	841686	43.087	ng	100
87) Indeno(1,2,3-cd)pyrene	28.27	276	1014589	42.822	ng	# 95
88) Benzo(b)fluoranthene	23.72	252	874023	42.653	ng	98
89) Benzo(k)fluoranthene	23.78	252	848489	43.264	ng	98
90) Benzo(a)pyrene	24.56	252	824998	43.100	ng	100
91) Dibenzo(a,h)anthracene	28.32	278	831066	43.741	ng	98
92) Benzo(g,h,i)perylene	29.36	276	818774	43.055	ng	97

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

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