

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG073020\
 Data File : BG046116.D
 Acq On : 30 Jul 2020 18:36
 Operator : CG/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ICVBG073020

Quant Time: Jul 30 19:16:33 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 30 18:43:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.96	152	112767	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	414337	20.00	ng	0.00
39) Acenaphthene-d10	14.58	164	264744	20.00	ng	0.00
64) Phenanthrene-d10	17.32	188	606145	20.00	ng	0.00
76) Chrysene-d12	21.57	240	617006	20.00	ng	0.00
86) Perylene-d12	24.68	264	671415	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.54	112	535898	82.61	ng	0.00
7) Phenol-d6	7.12	99	726598	82.19	ng	0.00
23) Nitrobenzene-d5	9.11	82	643061	84.03	ng	0.00
42) 2,4,6-Tribromophenol	16.07	330	346536	85.22	ng	0.00
45) 2-Fluorobiphenyl	13.21	172	1454733	79.82	ng	0.00
79) Terphenyl-d14	19.94	244	2298769	75.88	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.46	88	121969	41.309	ng	97
3) Pyridine	3.85	79	329324	40.516	ng	98
4) n-Nitrosodimethylamine	3.76	42	144285	41.709	ng	98
6) Aniline	7.28	93	429305	40.794	ng	98
8) 2-Chlorophenol	7.52	128	291777	40.695	ng	99
9) Benzaldehyde	7.09	77	225948	43.070	ng	97
10) Phenol	7.15	94	361174	40.660	ng	99
11) bis(2-Chloroethyl)ether	7.38	93	284978	40.334	ng	98
12) 1,3-Dichlorobenzene	7.85	146	347567	39.993	ng	97
13) 1,4-Dichlorobenzene	7.99	146	345716	39.616	ng	98
14) 1,2-Dichlorobenzene	8.31	146	328486	39.912	ng	99
15) Benzyl Alcohol	8.19	79	265199	43.421	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.49	45	463167	39.984	ng	99
17) 2-Methylphenol	8.40	107	250261	41.732	ng	100
18) Hexachloroethane	9.05	117	117826	40.371	ng	99
19) n-Nitroso-di-n-propylamine	8.76	70	236394	41.085	ng	99
20) 3+4-Methylphenols	8.72	107	345379	42.078	ng	98
22) Acetophenone	8.78	105	434170	40.831	ng	# 99
24) Nitrobenzene	9.15	77	336171	41.202	ng	98
25) Isophorone	9.68	82	641794	42.147	ng	99
26) 2-Nitrophenol	9.86	139	166139	44.929	ng	95
27) 2,4-Dimethylphenol	9.93	122	254105	42.237	ng	99
28) bis(2-Chloroethoxy)methane	10.16	93	334827	40.439	ng	98
29) 2,4-Dichlorophenol	10.40	162	303187	42.841	ng	98
30) 1,2,4-Trichlorobenzene	10.62	180	341602	41.204	ng	99
31) Naphthalene	10.80	128	882130	40.237	ng	99
32) Benzoic acid	10.07	122	177861	41.740	ng	97
33) 4-Chloroaniline	10.90	127	377182	42.102	ng	97
34) Hexachlorobutadiene	11.10	225	228163	41.342	ng	98
35) Caprolactam	11.67	113	102181	44.113	ng	# 85
36) 4-Chloro-3-methylphenol	12.03	107	298580	42.992	ng	97
37) 2-Methylnaphthalene	12.41	142	688955	40.841	ng	99
38) 1-Methylnaphthalene	12.63	142	648851	40.647	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.78	216	387438	40.617	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.77	237	228150	41.699	ng	99
43) 2,4,6-Trichlorophenol	13.01	196	264175	43.409	ng	99
44) 2,4,5-Trichlorophenol	13.09	196	286728	43.181	ng	97
46) 1,1'-Biphenyl	13.42	154	840984	40.494	ng	98
47) 2-Chloronaphthalene	13.46	162	671503	40.534	ng	98
48) 2-Nitroaniline	13.65	65	193435	44.546	ng	97
49) Acenaphthylene	14.30	152	1046710	40.690	ng	98
50) Dimethylphthalate	14.04	163	829848	40.840	ng	100
51) 2,6-Dinitrotoluene	14.15	165	204172	43.175	ng	93
52) Acenaphthene	14.64	154	703355	41.429	ng	98
53) 3-Nitroaniline	14.48	138	199154	42.479	ng	99
54) 2,4-Dinitrophenol	14.68	184	111146	40.887	ng	91
55) Dibenzofuran	14.98	168	1030423	40.222	ng	98
56) 4-Nitrophenol	14.79	139	176022	43.245	ng	99
57) 2,4-Dinitrotoluene	14.94	165	280798	43.187	ng	96
58) Fluorene	15.63	166	830105	40.012	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.21	232	266907	42.675	ng	99
60) Diethylphthalate	15.40	149	819176	40.982	ng	99
61) 4-Chlorophenyl-phenylether	15.63	204	437948	40.417	ng	97
62) 4-Nitroaniline	15.64	138	205447	43.572	ng	99
63) Azobenzene	15.92	77	764155	40.791	ng	98
65) 4,6-Dinitro-2-methylphenol	15.70	198	176892	44.127	ng	93
66) n-Nitrosodiphenylamine	15.84	169	754920	41.144	ng	98
67) 4-Bromophenyl-phenylether	16.52	248	307478	41.060	ng	97
68) Hexachlorobenzene	16.64	284	340985	40.609	ng	99
69) Atrazine	16.78	200	298709	42.279	ng	99
70) Pentachlorophenol	16.98	266	234880	43.111	ng	97
71) Phenanthrene	17.37	178	1337741	40.289	ng	100
72) Anthracene	17.46	178	1337514	40.498	ng	99
73) Carbazole	17.72	167	1306594	41.028	ng	99
74) Di-n-butylphthalate	18.29	149	1396236	41.943	ng	99
75) Fluoranthene	19.37	202	1670646	40.302	ng	98
77) Benzidine	19.55	184	732112	41.890	ng	99
78) Pyrene	19.73	202	1691988	40.813	ng	99
80) Butylbenzylphthalate	20.63	149	658151	43.825	ng	99
81) Benzo(a)anthracene	21.55	228	1674179	40.546	ng	99
82) 3,3'-Dichlorobenzidine	21.47	252	731443	42.659	ng	96
83) Chrysene	21.62	228	1641480	40.582	ng	98
84) Bis(2-ethylhexyl)phthalate	21.48	149	895923	41.729	ng	99
85) Di-n-octyl phthalate	22.66	149	1575736	43.346	ng	99
87) Indeno(1,2,3-cd)pyrene	28.20	276	2128292	41.563	ng	# 95
88) Benzo(b)fluoranthene	23.70	252	1871051	41.585	ng	99
89) Benzo(k)fluoranthene	23.77	252	1766578	40.030	ng	100
90) Benzo(a)pyrene	24.53	252	1749208	41.780	ng	99
91) Dibenzo(a,h)anthracene	28.27	278	1746295	40.795	ng	97
92) Benzo(g,h,i)perylene	29.29	276	1756007	40.619	ng	98

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

