

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG123019\
 Data File : BG043863.D
 Acq On : 30 Dec 2019 12:04
 Operator : JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC060

Quant Time: Dec 30 13:03:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 30 11:43:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.97	152	165245	20.00	ng	0.00
21) Naphthalene-d8	10.77	136	681835	20.00	ng	0.00
39) Acenaphthene-d10	14.60	164	431527	20.00	ng	0.00
64) Phenanthrene-d10	17.34	188	838782	20.00	ng	0.00
76) Chrysene-d12	21.59	240	704267	20.00	ng	0.00
87) Perylene-d12	24.72	264	814764	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.54	112	1035607	114.33	ng	0.00
7) Phenol-d6	7.13	99	1430912	111.91	ng	0.00
23) Nitrobenzene-d5	9.12	82	1413133	115.23	ng	0.00
42) 2,4,6-Tribromophenol	16.09	330	528272	116.19	ng	0.00
45) 2-Fluorobiphenyl	13.23	172	2621808	108.49	ng	0.00
79) Terphenyl-d14	19.95	244	2984187	94.23	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.44	88	223071	57.158	ng	99
3) Pyridine	3.83	79	655637	56.223	ng	98
4) n-Nitrosodimethylamine	3.74	42	300915	58.291	ng	96
6) Aniline	7.29	93	951374	56.593	ng	99
8) 2-Chlorophenol	7.53	128	604869	57.011	ng	99
9) Benzaldehyde	7.10	77	389958	46.540	ng	98
10) Phenol	7.16	94	741213	56.246	ng	97
11) bis(2-Chloroethyl)ether	7.39	93	636077	55.863	ng	98
12) 1,3-Dichlorobenzene	7.86	146	706327	57.000	ng	99
13) 1,4-Dichlorobenzene	8.00	146	695079	56.293	ng	98
14) 1,2-Dichlorobenzene	8.32	146	666996	56.340	ng	99
15) Benzyl Alcohol	8.20	79	578620	57.492	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.50	45	984642	55.708	ng	99
17) 2-Methylphenol	8.41	107	540926	56.645	ng	96
18) Hexachloroethane	9.05	117	273983	57.682	ng	96
19) n-Nitroso-di-n-propylamine	8.78	70	531247	55.634	ng	97
20) 3+4-Methylphenols	8.74	107	757015	57.225	ng	99
22) Acetophenone	8.78	105	971168	56.903	ng	# 97
24) Nitrobenzene	9.17	77	723296	57.602	ng	100
25) Isophorone	9.69	82	1369415	57.048	ng	99
26) 2-Nitrophenol	9.88	139	358336	61.520	ng	98
27) 2,4-Dimethylphenol	9.94	122	520424	58.035	ng	98
28) bis(2-Chloroethoxy)methane	10.18	93	906064	56.730	ng	99
29) 2,4-Dichlorophenol	10.41	162	628078	58.729	ng	99
30) 1,2,4-Trichlorobenzene	10.63	180	693267	58.080	ng	98
31) Naphthalene	10.82	128	1872431	56.687	ng	97
32) Benzoic acid	10.10	122	454517	65.182	ng	98
33) 4-Chloroaniline	10.92	127	900273	56.833	ng	98
34) Hexachlorobutadiene	11.12	225	424741	58.793	ng	99
35) Caprolactam	11.70	113	231790	58.020	ng	97
36) 4-Chloro-3-methylphenol	12.04	107	654272	56.799	ng	99
37) 2-Methylnaphthalene	12.43	142	1404277	56.328	ng	98
38) 1-Methylnaphthalene	12.64	142	1322637	55.891	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.80	216	722881	57.632	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.78	237	402210	60.054	nq	99
43) 2,4,6-Trichlorophenol	13.03	196	513840	59.187	nq	99
44) 2,4,5-Trichlorophenol	13.10	196	524735	59.001	nq	98
46) 1,1'-Biphenyl	13.43	154	1742311	56.241	nq	96
47) 2-Chloronaphthalene	13.47	162	1419121	56.815	nq	97
48) 2-Nitroaniline	13.67	65	470722	58.839	nq	99
49) Acenaphthylene	14.32	152	2004301	55.350	nq	96
50) Dimethylphthalate	14.05	163	1703247	55.239	nq	98
51) 2,6-Dinitrotoluene	14.17	165	393577	56.473	nq	99
52) Acenaphthene	14.66	154	1441602	56.984	nq	99
53) 3-Nitroaniline	14.49	138	452631	57.438	nq	99
54) 2,4-Dinitrophenol	14.69	184	199914	64.787	nq	98
55) Dibenzofuran	14.99	168	1922387	54.999	nq	95
56) 4-Nitrophenol	14.80	139	354462	58.655	nq	98
57) 2,4-Dinitrotoluene	14.95	165	546849	58.217	nq	97
58) Fluorene	15.65	166	1513562	54.602	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.22	232	450398	57.987	nq	99
60) Diethylphthalate	15.42	149	1703066	55.322	nq	97
61) 4-Chlorophenyl-phenylether	15.64	204	838186	56.233	nq	98
62) 4-Nitroaniline	15.66	138	439962	56.924	nq	98
63) Azobenzene	15.93	77	1580884	55.269	nq	96
65) 4,6-Dinitro-2-methylphenol	15.72	198	253449	63.661	nq	97
66) n-Nitrosodiphenylamine	15.85	169	1406529	57.029	nq	98
67) 4-Bromophenyl-phenylether	16.53	248	544867	58.212	nq	99
68) Hexachlorobenzene	16.66	284	584054	57.604	nq	99
69) Atrazine	16.80	200	463643	57.324	nq	98
70) Pentachlorophenol	16.99	266	347772	62.645	nq	99
71) Phenanthrene	17.39	178	2243937	55.754	nq	94
72) Anthracene	17.47	178	2209844	55.460	nq	94
73) Carbazole	17.74	167	2041969	56.033	nq	96
74) Di-n-butylphthalate	18.31	149	2468664	56.328	nq	96
75) Fluoranthene	19.39	202	2442953	56.285	nq	94
77) Benzidine	19.56	184	1137067	60.489	nq	95
78) Pyrene	19.75	202	2435517	55.026	nq	94
80) Butylbenzylphthalate	20.65	149	1271516	58.877	nq	99
81) Benzo(a)anthracene	21.57	228	2426210	55.583	nq	95
82) 3,3'-Dichlorobenzidine	21.49	252	1009747	58.298	nq	98
83) Chrysene	21.64	228	2333453	56.201	nq	94
84) Bis(2-ethylhexyl)phthalate	21.50	149	1718130	57.315	nq	96
85) Di-n-octyl phthalate	22.69	149	2901919	57.556	nq	97
86) Indeno(1,2,3-cd)pyrene	28.28	276	3310400	58.509	nq	# 85
88) Benzo(b)fluoranthene	23.74	252	2775021	58.013	nq	97
89) Benzo(k)fluoranthene	23.81	252	2563643	55.874	nq	96
90) Benzo(a)pyrene	24.58	252	2605294	57.561	nq	98
91) Dibenzo(a,h)anthracene	28.35	278	2640671	57.675	nq	99
92) Benzo(g,h,i)perylene	29.39	276	2625939	58.030	nq	99

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

