

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012120\
 Data File : BG044133.D
 Acq On : 21 Jan 2020 14:43
 Operator : JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 1/23/2020 7:58:16 AM

Quant Time: Jan 22 06:53:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 31 13:32:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	96349	20.00	ng	-0.03
21) Naphthalene-d8	10.73	136	418225	20.00	ng	-0.03
39) Acenaphthene-d10	14.57	164	270331	20.00	ng	-0.03
64) Phenanthrene-d10	17.31	188	549739	20.00	ng	-0.02
76) Chrysene-d12	21.56	240	449390	20.00	ng	-0.03
87) Perylene-d12	24.66	264	503133	20.00	ng	-0.06

System Monitoring Compounds

5) 2-Fluorophenol	5.51	112	469589	89.49	ng	-0.03
7) Phenol-d6	7.10	99	651350	88.80	ng	-0.02
23) Nitrobenzene-d5	9.09	82	605840	77.49	ng	-0.03
42) 2,4,6-Tribromophenol	16.06	330	238660	84.36	ng	-0.03
45) 2-Fluorobiphenyl	13.19	172	1311151	87.87	ng	-0.03
79) Terphenyl-d14	19.93	244	1722801	95.39	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.40	88	96576	42.210	ng	99
3) Pyridine	3.80	79	286008	42.315	ng	93
4) n-Nitrosodimethylamine	3.71	42	117576	39.071	ng	95
6) Aniline	7.25	93	391953	40.684	ng	99
8) 2-Chlorophenol	7.50	128	263287	42.739	ng	97
9) Benzaldehyde	7.06	77	160665	34.224	ng	96
10) Phenol	7.13	94	326244	43.169	ng	99
11) bis(2-Chloroethyl)ether	7.35	93	264386	40.528	ng	99
12) 1,3-Dichlorobenzene	7.82	146	301872	41.875	ng	98
13) 1,4-Dichlorobenzene	7.97	146	299779	42.146	ng	99
14) 1,2-Dichlorobenzene	8.29	146	286320	41.938	ng	98
15) Benzyl Alcohol	8.17	79	234413	40.493	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.46	45	380114	37.663	ng	98
17) 2-Methylphenol	8.38	107	231111	42.259	ng	97
18) Hexachloroethane	9.02	117	115825	41.849	ng	98
19) n-Nitroso-di-n-propylamine	8.74	70	217880	39.887	ng	98
20) 3+4-Methylphenols	8.71	107	324542	42.539	ng	96
22) Acetophenone	8.75	105	411966	39.622	ng	# 98
24) Nitrobenzene	9.13	77	300618	38.824	ng	98
25) Isophorone	9.66	82	575703	39.251	ng	99
26) 2-Nitrophenol	9.84	139	157996	43.129	ng	98
27) 2,4-Dimethylphenol	9.91	122	221820	40.225	ng	98
28) bis(2-Chloroethoxy)methane	10.14	93	382641	39.132	ng	99
29) 2,4-Dichlorophenol	10.39	162	271796	41.320	ng	98
30) 1,2,4-Trichlorobenzene	10.60	180	295908	40.121	ng	98
31) Naphthalene	10.79	128	849481	41.973	ng	99
32) Benzoic acid	10.06	122	136407m	32.674	ng	
33) 4-Chloroaniline	10.89	127	388040	40.199	ng	99
34) Hexachlorobutadiene	11.08	225	178469	39.633	ng	98
35) Caprolactam	11.66	113	96803	39.533	ng	97
36) 4-Chloro-3-methylphenol	12.02	107	294161	42.008	ng	99
37) 2-Methylnaphthalene	12.39	142	636966	41.804	ng	98
38) 1-Methylnaphthalene	12.61	142	601992	41.687	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.77	216	322487	40.701	ng	99

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41) Hexachlorocyclopentadiene	12.75	237	154169	37.531	nq	98
43) 2,4,6-Trichlorophenol	13.00	196	222526	40.816	nq	100
44) 2,4,5-Trichlorophenol	13.08	196	230360	40.967	nq	99
46) 1,1'-Biphenyl	13.40	154	814471	42.021	nq	99
47) 2-Chloronaphthalene	13.44	162	648126	41.382	nq	98
48) 2-Nitroaniline	13.64	65	201935	40.082	nq	96
49) Acenaphthylene	14.29	152	969204	43.054	nq	98
50) Dimethylphthalate	14.02	163	804823	41.930	nq	99
51) 2,6-Dinitrotoluene	14.13	165	176346	40.647	nq	99
52) Acenaphthene	14.63	154	620666	39.316	nq	99
53) 3-Nitroaniline	14.46	138	199963	40.588	nq	98
54) 2,4-Dinitrophenol	14.67	184	88092	42.608	nq	97
55) Dibenzofuran	14.97	168	935664	43.054	nq	97
56) 4-Nitrophenol	14.79	139	144291	38.220	nq	92
57) 2,4-Dinitrotoluene	14.93	165	246978	41.837	nq	100
58) Fluorene	15.62	166	733009	42.555	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.19	232	193403	39.999	nq	97
60) Diethylphthalate	15.39	149	816842	42.548	nq	98
61) 4-Chlorophenyl-phenylether	15.61	204	387101	41.387	nq	98
62) 4-Nitroaniline	15.63	138	192151	39.495	nq	96
63) Azobenzene	15.90	77	736915	41.389	nq	99
65) 4,6-Dinitro-2-methylphenol	15.69	198	122677	43.741	nq	99
66) n-Nitrosodiphenylamine	15.82	169	667419	41.226	nq	100
67) 4-Bromophenyl-phenylether	16.50	248	250348	40.490	nq	98
68) Hexachlorobenzene	16.63	284	272203	40.858	nq	99
69) Atrazine	16.77	200	199550	37.921	nq	99
70) Pentachlorophenol	16.97	266	130749	35.602	nq	98
71) Phenanthrene	17.35	178	1125829	42.696	nq	98
72) Anthracene	17.44	178	1119313	42.952	nq	97
73) Carbazole	17.71	167	1020070	42.377	nq	97
74) Di-n-butylphthalate	18.28	149	1306458	42.509	nq	97
75) Fluoranthene	19.36	202	1266091	41.985	nq	97
77) Benzidine	19.54	184	471184	41.714	nq	98
78) Pyrene	19.72	202	1266492	43.176	nq	97
80) Butylbenzylphthalate	20.62	149	607408	43.494	nq	96
81) Benzo(a)anthracene	21.54	228	1187889	42.630	nq	97
82) 3,3'-Dichlorobenzidine	21.46	252	446460	40.554	nq	99
83) Chrysene	21.60	228	1155388	43.637	nq	96
84) Bis(2-ethylhexyl)phthalate	21.46	149	819931	42.551	nq	98
85) Di-n-octyl phthalate	22.64	149	1417537	43.758	nq	97
86) Indeno(1,2,3-cd)pyrene	28.18	276	1426373	39.640	nq	99
88) Benzo(b)fluoranthene	23.69	252	1237007	41.589	nq	97
89) Benzo(k)fluoranthene	23.75	252	1200553	42.446	nq	97
90) Benzo(a)pyrene	24.52	252	1160795	41.349	nq	98
91) Dibenzo(a,h)anthracene	28.25	278	1147785	40.711	nq	98
92) Benzo(g,h,i)perylene	29.28	276	1146624	40.943	nq	98

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

