

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG032919\
 Data File : BG040244.D
 Acq On : 30 Mar 2019 3:16
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Mar 30 04:12:22 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 28 05:47:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	53756	20.00	ng	-0.01
21) Naphthalene-d8	10.87	136	261635	20.00	ng	-0.01
39) Acenaphthene-d10	14.69	164	170877	20.00	ng	-0.01
64) Phenanthrene-d10	17.43	188	411587	20.00	ng	-0.01
76) Chrysene-d12	21.72	240	385808	20.00	ng	0.00
87) Perylene-d12	25.03	264	393078	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	273831	84.68	ng	-0.01
7) Phenol-d6	7.21	99	400341	89.58	ng	-0.02
23) Nitrobenzene-d5	9.23	82	391031	79.44	ng	-0.01
42) 2,4,6-Tribromophenol	16.18	330	166543	90.18	ng	-0.01
45) 2-Fluorobiphenyl	13.31	172	927884	80.46	ng	-0.01
79) Terphenyl-d14	20.04	244	1368614	79.80	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.51	88	61463	40.423	ng	96
3) Pyridine	3.91	79	180879	44.183	ng	98
4) n-Nitrosodimethylamine	3.83	42	76183	40.230	ng	# 89
6) Aniline	7.38	93	255489	44.004	ng	98
8) 2-Chlorophenol	7.62	128	161727	42.725	ng	97
9) Benzaldehyde	7.20	77	133861	43.668	ng	93
10) Phenol	7.23	94	216275	44.587	ng	97
11) bis(2-Chloroethyl)ether	7.48	93	167570	43.132	ng	97
12) 1,3-Dichlorobenzene	7.95	146	171081	41.577	ng	96
13) 1,4-Dichlorobenzene	8.09	146	173267	42.278	ng	98
14) 1,2-Dichlorobenzene	8.41	146	164347	41.934	ng	97
15) Benzyl Alcohol	8.30	79	153359	42.611	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.58	45	309526	41.513	ng	99
17) 2-Methylphenol	8.49	107	147399	43.914	ng	97
18) Hexachloroethane	9.14	117	63050	40.445	ng	94
19) n-Nitroso-di-n-propylamine	8.86	70	141974	44.310	ng	95
20) 3+4-Methylphenols	8.82	107	203052	44.655	ng	97
22) Acetophenone	8.89	105	261561	40.077	ng	# 97
24) Nitrobenzene	9.27	77	202518	39.732	ng	99
25) Isophorone	9.78	82	414825	40.959	ng	100
26) 2-Nitrophenol	9.98	139	101819	42.157	ng	96
27) 2,4-Dimethylphenol	10.03	122	158556	41.329	ng	95
28) bis(2-Chloroethoxy)methane	10.27	93	242315	40.440	ng	98
29) 2,4-Dichlorophenol	10.51	162	175325	42.103	ng	98
30) 1,2,4-Trichlorobenzene	10.72	180	184673	40.388	ng	96
31) Naphthalene	10.92	128	549014	40.939	ng	99
32) Benzoic acid	10.16	122	61453	26.512	ng	91
33) 4-Chloroaniline	11.04	127	240868	42.107	ng	99
34) Hexachlorobutadiene	11.18	225	114239	39.066	ng	97
35) Caprolactam	11.82	113	73122	44.249	ng	98
36) 4-Chloro-3-methylphenol	12.14	107	207247	43.292	ng	97
37) 2-Methylnaphthalene	12.52	142	420095	42.355	ng	97
38) 1-Methylnaphthalene	12.74	142	408831	42.508	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.88	216	219840	40.478	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.85	237	107890	36.180	ng	97
43) 2,4,6-Trichlorophenol	13.12	196	144306	41.212	ng	99
44) 2,4,5-Trichlorophenol	13.20	196	161179	42.231	ng	98
46) 1,1'-Biphenyl	13.52	154	546786	40.498	ng	99
47) 2-Chloronaphthalene	13.57	162	421140	40.664	ng	98
48) 2-Nitroaniline	13.78	65	142370	40.755	ng	95
49) Acenaphthylene	14.41	152	724535	41.262	ng	100
50) Dimethylphthalate	14.14	163	551336	41.226	ng	99
51) 2,6-Dinitrotoluene	14.27	165	127597	42.492	ng	96
52) Acenaphthene	14.76	154	414815	41.227	ng	99
53) 3-Nitroaniline	14.61	138	132140	41.572	ng	96
54) 2,4-Dinitrophenol	14.81	184	80259	50.257	ng	90
55) Dibenzofuran	15.08	168	674788	41.737	ng	100
56) 4-Nitrophenol	14.90	139	108669	42.360	ng	99
57) 2,4-Dinitrotoluene	15.06	165	178372	42.750	ng	99
58) Fluorene	15.73	166	533049	41.935	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.31	232	151318	43.691	ng	99
60) Diethylphthalate	15.49	149	564682	41.263	ng	100
61) 4-Chlorophenyl-phenylether	15.72	204	286377	42.148	ng	98
62) 4-Nitroaniline	15.77	138	145679	42.707	ng	98
63) Azobenzene	16.01	77	529518	41.021	ng	99
65) 4,6-Dinitro-2-methylphenol	15.81	198	99753	43.139	ng	97
66) n-Nitrosodiphenylamine	15.94	169	485903	40.343	ng	99
67) 4-Bromophenyl-phenylether	16.62	248	188185	40.629	ng	95
68) Hexachlorobenzene	16.73	284	192853	41.411	ng	96
69) Atrazine	16.88	200	181134	40.429	ng	99
70) Pentachlorophenol	17.08	266	109442	40.648	ng	96
71) Phenanthrene	17.48	178	883169	40.824	ng	99
72) Anthracene	17.57	178	877855	40.541	ng	99
73) Carbazole	17.84	167	820863	40.056	ng	99
74) Di-n-butylphthalate	18.37	149	950953	39.148	ng	100
75) Fluoranthene	19.49	202	1032812	40.317	ng	99
77) Benzidine	19.67	184	539980	38.758	ng	99
78) Pyrene	19.84	202	1029520	40.578	ng	99
80) Butylbenzylphthalate	20.71	149	429096	39.524	ng	98
81) Benzo(a)anthracene	21.69	228	958371	40.329	ng	98
82) 3,3'-Dichlorobenzidine	21.61	252	365764	40.461	ng	98
83) Chrysene	21.77	228	951505	41.663	ng	97
84) Bis(2-ethylhexyl)phthalate	21.56	149	603980	38.918	ng	100
85) Di-n-octyl phthalate	22.79	149	1028781	38.908	ng	99
86) Indeno(1,2,3-cd)pyrene	28.80	276	1126998	40.347	ng	# 87
88) Benzo(b)fluoranthene	23.95	252	1013015	42.094	ng	99
89) Benzo(k)fluoranthene	24.03	252	918546	41.505	ng	99
90) Benzo(a)pyrene	24.86	252	941197	41.683	ng	99
91) Dibenzo(a,h)anthracene	28.89	278	904074	43.110	ng	99
92) Benzo(g,h,i)perylene	30.07	276	914567	42.435	ng	97

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

