

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG011119\
 Data File : BG039089.D
 Acq On : 11 Jan 2019 17:30
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Jan 11 18:51:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 09 14:23:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	25754	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	99549	20.00	ng	0.00
39) Acenaphthene-d10	14.67	164	66673	20.00	ng	0.00
64) Phenanthrene-d10	17.42	188	164242	20.00	ng	0.00
76) Chrysene-d12	21.69	240	170418	20.00	ng	-0.01
87) Perylene-d12	24.96	264	183187	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.58	112	119937	88.34	ng	0.00
7) Phenol-d6	7.18	99	170317	87.17	ng	0.00
23) Nitrobenzene-d5	9.21	82	159307	87.57	ng	0.00
42) 2,4,6-Tribromophenol	16.16	330	94507	84.56	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	427250	87.70	ng	0.00
79) Terphenyl-d14	20.02	244	773199	83.93	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.46	88	23494	44.191	ng	99
3) Pyridine	3.87	79	62827	43.490	ng	95
4) n-Nitrosodimethylamine	3.79	42	29517	45.406	ng	97
6) Aniline	7.36	93	98791	42.102	ng	97
8) 2-Chlorophenol	7.59	128	70593	44.111	ng	96
9) Benzaldehyde	7.17	77	43996	39.047	ng	98
10) Phenol	7.21	94	87111	44.470	ng	96
11) bis(2-Chloroethyl)ether	7.45	93	62437	41.705	ng	99
12) 1,3-Dichlorobenzene	7.91	146	80031	41.739	ng	95
13) 1,4-Dichlorobenzene	8.06	146	82485	42.476	ng	98
14) 1,2-Dichlorobenzene	8.38	146	78944	42.637	ng	98
15) Benzyl Alcohol	8.27	79	63638	43.251	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.54	45	117535	40.942	ng	99
17) 2-Methylphenol	8.47	107	57916	42.580	ng	96
18) Hexachloroethane	9.10	117	27908	42.302	ng	98
19) n-Nitroso-di-n-propylamine	8.83	70	52809	41.160	ng	94
20) 3+4-Methylphenols	8.80	107	82249	42.413	ng	96
22) Acetophenone	8.86	105	113943	43.829	ng	# 95
24) Nitrobenzene	9.25	77	78653	42.773	ng	100
25) Isophorone	9.76	82	132955	42.427	ng	100
26) 2-Nitrophenol	9.96	139	43741	43.977	ng	95
27) 2,4-Dimethylphenol	10.01	122	62008	44.313	ng	90
28) bis(2-Chloroethoxy)methane	10.24	93	82142	43.614	ng	99
29) 2,4-Dichlorophenol	10.49	162	78465	45.136	ng	92
30) 1,2,4-Trichlorobenzene	10.70	180	89702	42.945	ng	95
31) Naphthalene	10.89	128	217895	43.642	ng	99
32) Benzoic acid	10.16	122	25657	33.392	ng	98
33) 4-Chloroaniline	11.01	127	96410	43.154	ng	98
34) Hexachlorobutadiene	11.15	225	63368	44.090	ng	98
35) Caprolactam	11.80	113	28299	40.595	ng	97
36) 4-Chloro-3-methylphenol	12.13	107	79182	42.586	ng	93
37) 2-Methylnaphthalene	12.50	142	157919	42.187	ng	97
38) 1-Methylnaphthalene	12.71	142	150780	41.965	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.86	216	115011	45.932	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.83	237	51852	39.867	ng	99
43) 2,4,6-Trichlorophenol	13.10	196	70555	44.770	ng	96
44) 2,4,5-Trichlorophenol	13.18	196	74419	44.679	ng	98
46) 1,1'-Biphenyl	13.50	154	235976	44.667	ng	99
47) 2-Chloronaphthalene	13.55	162	189981	44.438	ng	98
48) 2-Nitroaniline	13.77	65	52975	44.231	ng	92
49) Acenaphthylene	14.39	152	280330	43.625	ng	99
50) Dimethylphthalate	14.12	163	237240	42.107	ng	99
51) 2,6-Dinitrotoluene	14.25	165	52578	41.740	ng	97
52) Acenaphthene	14.73	154	167106	43.283	ng	94
53) 3-Nitroaniline	14.59	138	55080	43.104	ng	96
54) 2,4-Dinitrophenol	14.80	184	44134	60.043	ng	95
55) Dibenzofuran	15.06	168	292115	42.841	ng	100
56) 4-Nitrophenol	14.90	139	36903	40.136	ng	90
57) 2,4-Dinitrotoluene	15.04	165	76080	42.098	ng	96
58) Fluorene	15.72	166	216046	42.094	ng	94
59) 2,3,4,6-Tetrachlorophenol	15.29	232	68190	43.370	ng	96
60) Diethylphthalate	15.47	149	237680	40.944	ng	99
61) 4-Chlorophenyl-phenylether	15.70	204	139156	43.035	ng	99
62) 4-Nitroaniline	15.75	138	57529	43.216	ng	99
63) Azobenzene	15.99	77	191207	42.120	ng	98
65) 4,6-Dinitro-2-methylphenol	15.80	198	43725	35.934	ng	97
66) n-Nitrosodiphenylamine	15.92	169	213676	43.885	ng	98
67) 4-Bromophenyl-phenylether	16.60	248	92085	43.616	ng	98
68) Hexachlorobenzene	16.71	284	98510	42.760	ng	94
69) Atrazine	16.86	200	89405	43.224	ng	97
70) Pentachlorophenol	17.07	266	53480	40.648	ng	95
71) Phenanthrene	17.46	178	372500	42.908	ng	99
72) Anthracene	17.55	178	370472	43.161	ng	98
73) Carbazole	17.83	167	363142	42.856	ng	98
74) Di-n-butylphthalate	18.35	149	403458	42.800	ng	98
75) Fluoranthene	19.47	202	454142	42.198	ng	99
77) Benzidine	19.65	184	185514	35.591	ng	99
78) Pyrene	19.83	202	464114	42.734	ng	98
80) Butylbenzylphthalate	20.70	149	178103	43.117	ng	98
81) Benzo(a)anthracene	21.67	228	445061	42.901	ng	98
82) 3,3'-Dichlorobenzidine	21.59	252	181821	43.015	ng	96
83) Chrysene	21.74	228	421793	42.749	ng	98
84) Bis(2-ethylhexyl)phthalate	21.54	149	251159	43.791	ng	96
85) Di-n-octyl phthalate	22.75	149	420698	43.379	ng	100
86) Indeno(1,2,3-cd)pyrene	28.72	276	509164	43.187	ng	99
88) Benzo(b)fluoranthene	23.92	252	444550	44.011	ng	99
89) Benzo(k)fluoranthene	23.99	252	436475	43.704	ng	98
90) Benzo(a)pyrene	24.81	252	425872	43.620	ng	98
91) Dibenzo(a,h)anthracene	28.78	278	421702	43.424	ng	99
92) Benzo(g,h,i)perylene	29.91	276	416669	43.340	ng	97

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

