

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022619\
 Data File : BG039597.D
 Acq On : 26 Feb 2019 12:45
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Feb 26 23:55:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 29 15:41:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	53740	20.00	ng	-0.02
21) Naphthalene-d8	10.99	136	227913	20.00	ng	-0.02
39) Acenaphthene-d10	14.79	164	155824	20.00	ng	-0.02
64) Phenanthrene-d10	17.54	188	384579	20.00	ng	-0.02
76) Chrysene-d12	21.82	240	383142	20.00	ng	-0.02
87) Perylene-d12	25.20	264	399178	20.00	ng	-0.04

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.71	112	249928	79.60	ng	0.00
7) Phenol-d6	7.31	99	367343	79.48	ng	0.00
23) Nitrobenzene-d5	9.35	82	345195	94.62	ng	-0.02
42) 2,4,6-Tribromophenol	16.28	330	154655	92.17	ng	-0.01
45) 2-Fluorobiphenyl	13.42	172	841266	77.45	ng	-0.02
79) Terphenyl-d14	20.13	244	1397544	77.21	ng	-0.02

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.61	88	55081	40.103	ng	94
3) Pyridine	4.01	79	151565	41.679	ng	99
4) n-Nitrosodimethylamine	3.93	42	65964	36.271	ng	86
6) Aniline	7.50	93	223131	38.542	ng	96
8) 2-Chlorophenol	7.73	128	137865	40.168	ng	97
9) Benzaldehyde	7.31	77	90812	39.086	ng	96
10) Phenol	7.34	94	190964	39.636	ng	93
11) bis(2-Chloroethyl)ether	7.58	93	146321	38.434	ng	96
12) 1,3-Dichlorobenzene	8.05	146	162885	39.175	ng	97
13) 1,4-Dichlorobenzene	8.21	146	164764	39.757	ng	97
14) 1,2-Dichlorobenzene	8.52	146	157289	39.205	ng	98
15) Benzyl Alcohol	8.41	79	142535	39.281	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.69	45	288323	33.536	ng	99
17) 2-Methylphenol	8.60	107	121486	38.965	ng	96
18) Hexachloroethane	9.25	117	56903	39.910	ng	95
19) n-Nitroso-di-n-propylamine	8.97	70	123680	37.702	ng	93
20) 3+4-Methylphenols	8.93	107	170498	39.711	ng	97
22) Acetophenone	9.00	105	244742	39.977	ng	# 94
24) Nitrobenzene	9.39	77	176300	44.731	ng	95
25) Isophorone	9.90	82	310728	38.435	ng	100
26) 2-Nitrophenol	10.10	139	88284	54.696	ng	97
27) 2,4-Dimethylphenol	10.14	122	132188	40.420	ng	98
28) bis(2-Chloroethoxy)methane	10.38	93	199676	40.099	ng	95
29) 2,4-Dichlorophenol	10.63	162	153720	43.007	ng	96
30) 1,2,4-Trichlorobenzene	10.84	180	167416	41.720	ng	98
31) Naphthalene	11.04	128	446196	39.463	ng	99
32) Benzoic acid	10.26	122	98307	46.299	ng	93
33) 4-Chloroaniline	11.15	127	211665	40.375	ng	99
34) Hexachlorobutadiene	11.30	225	111621	41.826	ng	94
35) Caprolactam	11.94	113	66189	44.750	ng	# 88
36) 4-Chloro-3-methylphenol	12.25	107	173968	42.085	ng	96
37) 2-Methylnaphthalene	12.63	142	328487	39.226	ng	96
38) 1-Methylnaphthalene	12.85	142	315226	39.050	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.99	216	202833	40.911	ng	99

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41) Hexachlorocyclopentadiene	12.95	237	116869	43.259	ng	99
43) 2,4,6-Trichlorophenol	13.23	196	134263	44.044	ng	98
44) 2,4,5-Trichlorophenol	13.30	196	140061	43.839	ng	99
46) 1,1'-Biphenyl	13.63	154	494785	38.163	ng	98
47) 2-Chloronaphthalene	13.68	162	382798	39.248	ng	98
48) 2-Nitroaniline	13.88	65	133271	47.642	ng	94
49) Acenaphthylene	14.52	152	578899	39.336	ng	99
50) Dimethylphthalate	14.24	163	505694	40.182	ng	100
51) 2,6-Dinitrotoluene	14.37	165	113939	48.382	ng	95
52) Acenaphthene	14.86	154	367878	38.509	ng	99
53) 3-Nitroaniline	14.70	138	121124	47.759	ng	# 94
54) 2,4-Dinitrophenol	14.90	184	58393	62.804	ng	91
55) Dibenzofuran	15.19	168	603420	39.396	ng	98
56) 4-Nitrophenol	14.99	139	100705	46.845	ng	89
57) 2,4-Dinitrotoluene	15.15	165	161982	52.824	ng	# 88
58) Fluorene	15.84	166	454857	39.837	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.41	232	136282	45.557	ng	96
60) Diethylphthalate	15.59	149	518950	39.883	ng	97
61) 4-Chlorophenyl-phenylether	15.82	204	262865	40.659	ng	96
62) 4-Nitroaniline	15.87	138	127163	48.169	ng	90
63) Azobenzene	16.11	77	466854	36.708	ng	97
65) 4,6-Dinitro-2-methylphenol	15.91	198	100579	60.275	ng	96
66) n-Nitrosodiphenylamine	16.04	169	452859	38.727	ng	99
67) 4-Bromophenyl-phenylether	16.72	248	175778	41.200	ng	90
68) Hexachlorobenzene	16.83	284	177771	41.530	ng	97
69) Atrazine	16.98	200	148367	34.719	ng	97
70) Pentachlorophenol	17.17	266	125028	42.025	ng	97
71) Phenanthrene	17.58	178	770788	38.724	ng	97
72) Anthracene	17.67	178	767130	39.127	ng	98
73) Carbazole	17.94	167	763174	39.004	ng	98
74) Di-n-butylphthalate	18.47	149	885838	38.806	ng	99
75) Fluoranthene	19.58	202	918332	39.749	ng	98
77) Benzidine	19.76	184	442080	35.193	ng	98
78) Pyrene	19.94	202	941542	39.475	ng	99
80) Butylbenzylphthalate	20.81	149	403870	40.140	ng	95
81) Benzo(a)anthracene	21.81	228	895184	39.541	ng	99
82) 3,3'-Dichlorobenzidine	21.72	252	349621	41.648	ng	98
83) Chrysene	21.88	228	853588	39.607	ng	99
84) Bis(2-ethylhexyl)phthalate	21.67	149	573658	39.965	ng	99
85) Di-n-octyl phthalate	22.93	149	945103	39.853	ng	96
86) Indeno(1,2,3-cd)pyrene	29.08	276	980942	42.423	ng	# 95
88) Benzo(b)fluoranthene	24.12	252	856408	39.340	ng	99
89) Benzo(k)fluoranthene	24.19	252	836391	38.268	ng	98
90) Benzo(a)pyrene	25.04	252	834015	39.780	ng	98
91) Dibenzo(a,h)anthracene	29.16	278	798741	40.681	ng	100
92) Benzo(g,h,i)perylene	30.31	276	805606	41.165	ng	99

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

