

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010620\
 Data File : BG043946.D
 Acq On : 7 Jan 2020 00:47
 Operator : JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Jan 07 04:56:40 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	7.96	152	130781	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	558645	20.00	ng	0.00
39) Acenaphthene-d10	14.59	164	373128	20.00	ng	0.00
64) Phenanthrene-d10	17.34	188	793149	20.00	ng	0.00
76) Chrysene-d12	21.59	240	664031	20.00	ng	0.00
87) Perylene-d12	24.71	264	762826	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.54	112	605908	85.07	ng	0.00
7) Phenol-d6	7.13	99	845282	84.90	ng	0.00
23) Nitrobenzene-d5	9.12	82	797573	76.38	ng	0.00
42) 2,4,6-Tribromophenol	16.08	330	337151	86.34	ng	0.00
45) 2-Fluorobiphenyl	13.22	172	1710796	83.07	ng	0.00
79) Terphenyl-d14	19.95	244	2260119	81.05	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.43	88	126491	40.730	ng	97
3) Pyridine	3.83	79	377639	41.162	ng	97
4) n-Nitrosodimethylamine	3.75	42	157054	38.450	ng	93
6) Aniline	7.28	93	517370	39.563	ng	99
8) 2-Chlorophenol	7.53	128	340740	40.749	ng	96
9) Benzaldehyde	7.09	77	231856	36.385	ng	98
10) Phenol	7.15	94	429282	41.848	ng	98
11) bis(2-Chloroethyl)ether	7.38	93	349616	39.483	ng	98
12) 1,3-Dichlorobenzene	7.85	146	395642	40.433	ng	98
13) 1,4-Dichlorobenzene	8.00	146	390679	40.465	ng	99
14) 1,2-Dichlorobenzene	8.32	146	375914	40.565	ng	99
15) Benzyl Alcohol	8.19	79	317252	40.375	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.49	45	512675	37.424	ng	99
17) 2-Methylphenol	8.40	107	305110	41.102	ng	100
18) Hexachloroethane	9.04	117	152699	40.646	ng	97
19) n-Nitroso-di-n-propylamine	8.76	70	287401	38.762	ng	98
20) 3+4-Methylphenols	8.73	107	426610	41.195	ng	99
22) Acetophenone	8.78	105	541350	38.979	ng	# 98
24) Nitrobenzene	9.16	77	400092	38.683	ng	100
25) Isophorone	9.69	82	761102	38.848	ng	99
26) 2-Nitrophenol	9.87	139	202309	41.344	ng	99
27) 2,4-Dimethylphenol	9.93	122	287816	39.074	ng	99
28) bis(2-Chloroethoxy)methane	10.17	93	505183	38.678	ng	100
29) 2,4-Dichlorophenol	10.41	162	361747	41.172	ng	97
30) 1,2,4-Trichlorobenzene	10.63	180	390308	39.619	ng	98
31) Naphthalene	10.81	128	1100332	40.702	ng	99
32) Benzoic acid	10.07	122	154519	28.661	ng	95
33) 4-Chloroaniline	10.91	127	516304	40.042	ng	99
34) Hexachlorobutadiene	11.11	225	237329	39.457	ng	99
35) Caprolactam	11.68	113	136041	41.592	ng	99
36) 4-Chloro-3-methylphenol	12.04	107	390879	41.790	ng	99
37) 2-Methylnaphthalene	12.42	142	830419	40.802	ng	97
38) 1-Methylnaphthalene	12.63	142	789752	40.942	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.79	216	425904	38.944	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.78	237	206630	36.444	ng	99
43) 2,4,6-Trichlorophenol	13.02	196	302328	40.176	ng	99
44) 2,4,5-Trichlorophenol	13.10	196	314696	40.547	ng	98
46) 1,1'-Biphenyl	13.43	154	1072284	40.081	ng	98
47) 2-Chloronaphthalene	13.47	162	860923	39.825	ng	99
48) 2-Nitroaniline	13.66	65	279931	40.255	ng	96
49) Acenaphthylene	14.32	152	1270941	40.904	ng	100
50) Dimethylphthalate	14.05	163	1084813	40.946	ng	99
51) 2,6-Dinitrotoluene	14.16	165	245710	41.033	ng	94
52) Acenaphthene	14.66	154	877639	40.277	ng	97
53) 3-Nitroaniline	14.49	138	277554	40.816	ng	100
54) 2,4-Dinitrophenol	14.69	184	96471	34.938	ng	99
55) Dibenzofuran	14.99	168	1241565	41.391	ng	100
56) 4-Nitrophenol	14.80	139	208265	39.967	ng	96
57) 2,4-Dinitrotoluene	14.94	165	340473	41.786	ng	96
58) Fluorene	15.64	166	982002	41.304	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.22	232	278753	41.768	ng	97
60) Diethylphthalate	15.41	149	1106422	41.754	ng	99
61) 4-Chlorophenyl-phenylether	15.64	204	529914	41.047	ng	98
62) 4-Nitroaniline	15.65	138	283418	42.205	ng	96
63) Azobenzene	15.92	77	1012013	41.181	ng	99
65) 4,6-Dinitro-2-methylphenol	15.71	198	155097	38.872	ng	99
66) n-Nitrosodiphenylamine	15.85	169	911095	39.007	ng	99
67) 4-Bromophenyl-phenylether	16.52	248	349224	39.148	ng	98
68) Hexachlorobenzene	16.65	284	379676	39.500	ng	98
69) Atrazine	16.79	200	302837	39.887	ng	99
70) Pentachlorophenol	16.99	266	191080	36.062	ng	98
71) Phenanthrene	17.38	178	1531657	40.261	ng	100
72) Anthracene	17.47	178	1529698	40.686	ng	99
73) Carbazole	17.73	167	1411303	40.637	ng	99
74) Di-n-butylphthalate	18.30	149	1775783	40.048	ng	100
75) Fluoranthene	19.39	202	1715099	39.420	ng	99
77) Benzidine	19.56	184	744466	44.603	ng	99
78) Pyrene	19.74	202	1727496	39.856	ng	99
80) Butylbenzylphthalate	20.64	149	846970	41.044	ng	98
81) Benzo(a)anthracene	21.57	228	1665706	40.455	ng	99
82) 3,3'-Dichlorobenzidine	21.48	252	646268	39.729	ng	100
83) Chrysene	21.63	228	1605847	41.045	ng	99
84) Bis(2-ethylhexyl)phthalate	21.49	149	1155479	40.581	ng	99
85) Di-n-octyl phthalate	22.68	149	1985791	41.485	ng	99
86) Indeno(1,2,3-cd)pyrene	28.25	276	2046202	38.484	ng	99
88) Benzo(b)fluoranthene	23.73	252	1792810	39.755	ng	99
89) Benzo(k)fluoranthene	23.79	252	1747478	40.749	ng	99
90) Benzo(a)pyrene	24.57	252	1707616	40.120	ng	98
91) Dibenzo(a,h)anthracene	28.32	278	1616126	37.808	ng	97
92) Benzo(g,h,i)perylene	29.35	276	1593907	37.539	ng	99

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

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