

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021920\
 Data File : BG044493.D
 Acq On : 19 Feb 2020 17:27
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Feb 20 07:14:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 10 16:11:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	65241	20.00	ng	-0.01
21) Naphthalene-d8	10.69	136	266020	20.00	ng	0.00
39) Acenaphthene-d10	14.53	164	171388	20.00	ng	-0.01
64) Phenanthrene-d10	17.27	188	380811	20.00	ng	-0.01
76) Chrysene-d12	21.52	240	377049	20.00	ng	-0.01
87) Perylene-d12	24.60	264	412352	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.47	112	313078	84.69	ng	0.00
7) Phenol-d6	7.07	99	427989	86.22	ng	0.00
23) Nitrobenzene-d5	9.05	82	390650	82.91	ng	-0.01
42) 2,4,6-Tribromophenol	16.02	330	177230	88.09	ng	0.00
45) 2-Fluorobiphenyl	13.15	172	890928	87.05	ng	-0.01
79) Terphenyl-d14	19.89	244	1376144	75.07	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.36	88	69263	41.702	ng	98
3) Pyridine	3.75	79	191083	43.182	ng	97
4) n-Nitrosodimethylamine	3.67	42	73466	39.731	ng	94
6) Aniline	7.22	93	257900	41.855	ng	98
8) 2-Chlorophenol	7.46	128	171751	42.274	ng	99
9) Benzaldehyde	7.03	77	111477	39.430	ng	99
10) Phenol	7.09	94	211043	42.958	ng	99
11) bis(2-Chloroethyl)ether	7.31	93	172445	41.057	ng	98
12) 1,3-Dichlorobenzene	7.78	146	203129	41.875	ng	99
13) 1,4-Dichlorobenzene	7.93	146	200324	41.696	ng	99
14) 1,2-Dichlorobenzene	8.24	146	191538	42.178	ng	98
15) Benzyl Alcohol	8.13	79	149295	42.480	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.43	45	235609	41.445	ng	98
17) 2-Methylphenol	8.34	107	146328	41.746	ng	97
18) Hexachloroethane	8.97	117	74011	41.052	ng	95
19) n-Nitroso-di-n-propylamine	8.70	70	135406	40.929	ng	98
20) 3+4-Methylphenols	8.67	107	200656	41.538	ng	99
22) Acetophenone	8.71	105	265594	41.041	ng	# 98
24) Nitrobenzene	9.09	77	190731	41.463	ng	99
25) Isophorone	9.61	82	357579	40.715	ng	100
26) 2-Nitrophenol	9.80	139	103047	43.172	ng	98
27) 2,4-Dimethylphenol	9.87	122	142019	42.150	ng	99
28) bis(2-Chloroethoxy)methane	10.10	93	236811	40.420	ng	100
29) 2,4-Dichlorophenol	10.35	162	172938	42.448	ng	99
30) 1,2,4-Trichlorobenzene	10.56	180	194152	41.560	ng	98
31) Naphthalene	10.74	128	547297	42.405	ng	100
32) Benzoic acid	10.03	122	54439	31.458	ng	93
33) 4-Chloroaniline	10.85	127	244308	41.620	ng	99
34) Hexachlorobutadiene	11.03	225	121549	42.284	ng	99
35) Caprolactam	11.62	113	65447	43.853	ng	99
36) 4-Chloro-3-methylphenol	11.99	107	178199	41.718	ng	98
37) 2-Methylnaphthalene	12.35	142	403209	42.077	ng	99
38) 1-Methylnaphthalene	12.57	142	378875	41.936	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.73	216	216022	42.136	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.71	237	85804	34.880	ng	98
43) 2,4,6-Trichlorophenol	12.97	196	142509	43.004	ng	98
44) 2,4,5-Trichlorophenol	13.04	196	147084	43.455	ng	99
46) 1,1'-Biphenyl	13.36	154	519986	42.062	ng	99
47) 2-Chloronaphthalene	13.40	162	412162	41.897	ng	99
48) 2-Nitroaniline	13.60	65	122415	42.166	ng	98
49) Acenaphthylene	14.25	152	617784	42.816	ng	99
50) Dimethylphthalate	13.98	163	517656	42.018	ng	100
51) 2,6-Dinitrotoluene	14.10	165	114376	41.638	ng	99
52) Acenaphthene	14.60	154	391425	41.853	ng	99
53) 3-Nitroaniline	14.43	138	128533	42.294	ng	97
54) 2,4-Dinitrophenol	14.65	184	55382	38.407	ng	96
55) Dibenzofuran	14.93	168	607897	42.931	ng	98
56) 4-Nitrophenol	14.75	139	94693	42.537	ng	95
57) 2,4-Dinitrotoluene	14.89	165	164100	42.623	ng	96
58) Fluorene	15.58	166	480658	43.098	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.17	232	127188	41.601	ng	98
60) Diethylphthalate	15.35	149	520963	42.438	ng	99
61) 4-Chlorophenyl-phenylether	15.57	204	255160	42.195	ng	98
62) 4-Nitroaniline	15.60	138	134888	44.419	ng	95
63) Azobenzene	15.86	77	446489	41.499	ng	99
65) 4,6-Dinitro-2-methylphenol	15.66	198	85165	40.515	ng	95
66) n-Nitrosodiphenylamine	15.79	169	440508	41.277	ng	97
67) 4-Bromophenyl-phenylether	16.47	248	169810	40.929	ng	96
68) Hexachlorobenzene	16.59	284	192494	41.413	ng	99
69) Atrazine	16.74	200	142328	38.910	ng	99
70) Pentachlorophenol	16.93	266	89112	39.565	ng	98
71) Phenanthrene	17.32	178	791751	42.936	ng	98
72) Anthracene	17.41	178	779803	42.773	ng	98
73) Carbazole	17.68	167	730882	43.487	ng	98
74) Di-n-butylphthalate	18.24	149	916712	44.138	ng	98
75) Fluoranthene	19.33	202	944853	44.293	ng	98
77) Benzidine	19.51	184	459188	43.906	ng	98
78) Pyrene	19.69	202	944902	39.023	ng	97
80) Butylbenzylphthalate	20.58	149	435119	39.432	ng	98
81) Benzo(a)anthracene	21.50	228	929117	39.983	ng	97
82) 3,3'-Dichlorobenzidine	21.42	252	364409	40.405	ng	99
83) Chrysene	21.56	228	901120	40.118	ng	97
84) Bis(2-ethylhexyl)phthalate	21.42	149	597222	39.321	ng	100
85) Di-n-octyl phthalate	22.59	149	1076576	40.967	ng	99
86) Indeno(1,2,3-cd)pyrene	28.09	276	1158460	39.532	ng	100
88) Benzo(b)fluoranthene	23.63	252	995006	41.875	ng	99
89) Benzo(k)fluoranthene	23.70	252	967610	41.782	ng	99
90) Benzo(a)pyrene	24.46	252	944543	42.043	ng	99
91) Dibenzo(a,h)anthracene	28.15	278	947848	42.631	ng	99
92) Benzo(g,h,i)perylene	29.18	276	937761	42.132	ng	98

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

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