

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091619\
 Data File : BG042850.D
 Acq On : 16 Sep 2019 15:58
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Sep 16 16:34:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG091219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 12 14:06:09 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	74995	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	314402	20.00	ng	0.00
39) Acenaphthene-d10	14.74	164	218818	20.00	ng	0.00
64) Phenanthrene-d10	17.48	188	563963	20.00	ng	0.00
76) Chrysene-d12	21.76	240	613590	20.00	ng	0.00
87) Perylene-d12	25.03	264	686114	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.69	112	363526	84.27	ng	0.00
7) Phenol-d6	7.27	99	508646	82.15	ng	0.00
23) Nitrobenzene-d5	9.28	82	496524	82.22	ng	0.00
42) 2,4,6-Tribromophenol	16.22	330	245617	84.37	ng	0.00
45) 2-Fluorobiphenyl	13.36	172	1109998	80.30	ng	0.00
79) Terphenyl-d14	20.09	244	2180862	78.77	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.59	88	82014	41.660	ng	98
3) Pyridine	3.99	79	248177	41.870	ng	98
4) n-Nitrosodimethylamine	3.89	42	111230	38.565	ng	81
6) Aniline	7.44	93	337288	40.377	ng	100
8) 2-Chlorophenol	7.68	128	190880	40.937	ng	97
9) Benzaldehyde	7.25	77	153995	38.220	ng	98
10) Phenol	7.30	94	262715	41.375	ng	99
11) bis(2-Chloroethyl)ether	7.53	93	196757	40.376	ng	97
12) 1,3-Dichlorobenzene	8.01	146	234198	41.180	ng	98
13) 1,4-Dichlorobenzene	8.15	146	229888	40.375	ng	96
14) 1,2-Dichlorobenzene	8.48	146	216610	39.965	ng	99
15) Benzyl Alcohol	8.35	79	213070	40.100	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.65	45	411065	38.412	ng	97
17) 2-Methylphenol	8.55	107	175207	41.245	ng	97
18) Hexachloroethane	9.21	117	85876	40.852	ng	95
19) n-Nitroso-di-n-propylamine	8.92	70	183138	40.142	ng	94
20) 3+4-Methylphenols	8.88	107	236670	40.417	ng	96
22) Acetophenone	8.93	105	311396	39.217	ng	96
24) Nitrobenzene	9.32	77	260608	41.114	ng	98
25) Isophorone	9.85	82	498441	39.598	ng	99
26) 2-Nitrophenol	10.03	139	105774	42.436	ng	93
27) 2,4-Dimethylphenol	10.09	122	175018	39.694	ng	97
28) bis(2-Chloroethoxy)methane	10.32	93	280022	39.890	ng	# 98
29) 2,4-Dichlorophenol	10.57	162	204934	39.997	ng	97
30) 1,2,4-Trichlorobenzene	10.79	180	246049	39.333	ng	99
31) Naphthalene	10.98	128	619404	39.902	ng	99
32) Benzoic acid	10.20	122	134679	39.037	ng	90
33) 4-Chloroaniline	11.07	127	292133	40.264	ng	98
34) Hexachlorobutadiene	11.27	225	169194	38.603	ng	97
35) Caprolactam	11.84	113	84878	44.547	ng	88
36) 4-Chloro-3-methylphenol	12.18	107	227671	40.924	ng	98
37) 2-Methylnaphthalene	12.58	142	462120	39.741	ng	95
38) 1-Methylnaphthalene	12.79	142	438511	40.208	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.94	216	284005	39.277	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.92	237	166958	40.306	nq	99
43) 2,4,6-Trichlorophenol	13.17	196	192185	39.923	nq	95
44) 2,4,5-Trichlorophenol	13.24	196	200152	40.592	nq	95
46) 1,1'-Biphenyl	13.58	154	611806	39.939	nq	98
47) 2-Chloronaphthalene	13.62	162	506949	40.006	nq	98
48) 2-Nitroaniline	13.81	65	199721	45.434	nq	98
49) Acenaphthylene	14.46	152	790595	40.826	nq	99
50) Dimethylphthalate	14.19	163	682645	41.645	nq	99
51) 2,6-Dinitrotoluene	14.30	165	150923	44.870	nq	87
52) Acenaphthene	14.80	154	516505	40.827	nq	98
53) 3-Nitroaniline	14.63	138	169648	44.940	nq	93
54) 2,4-Dinitrophenol	14.83	184	79087	47.763	nq	85
55) Dibenzofuran	15.13	168	768266	40.856	nq	99
56) 4-Nitrophenol	14.92	139	138203	47.350	nq	95
57) 2,4-Dinitrotoluene	15.08	165	217165	48.439	nq	# 99
58) Fluorene	15.79	166	643449	41.288	nq	97
59) 2,3,4,6-Tetrachlorophenol	15.36	232	205531	42.999	nq	97
60) Diethylphthalate	15.55	149	696836	41.879	nq	99
61) 4-Chlorophenyl-phenylether	15.77	204	373237	40.505	nq	96
62) 4-Nitroaniline	15.79	138	190184	46.739	nq	96
63) Azobenzene	16.07	77	666753	41.612	nq	98
65) 4,6-Dinitro-2-methylphenol	15.85	198	110261	45.693	nq	97
66) n-Nitrosodiphenylamine	15.99	169	581426	38.313	nq	100
67) 4-Bromophenyl-phenylether	16.67	248	256689	37.604	nq	98
68) Hexachlorobenzene	16.79	284	274417	37.623	nq	99
69) Atrazine	16.93	200	189061	35.775	nq	97
70) Pentachlorophenol	17.12	266	176983	40.903	nq	91
71) Phenanthrene	17.52	178	1088892	40.373	nq	100
72) Anthracene	17.61	178	1091593	40.511	nq	98
73) Carbazole	17.88	167	1075886	42.010	nq	97
74) Di-n-butylphthalate	18.44	149	1295624	42.232	nq	99
75) Fluoranthene	19.53	202	1505428	43.068	nq	99
77) Benzidine	19.70	184	645739	37.668	nq	99
78) Pyrene	19.89	202	1526319	39.733	nq	99
80) Butylbenzylphthalate	20.77	149	638123	41.045	nq	100
81) Benzo(a)anthracene	21.74	228	1544941	39.601	nq	98
82) 3,3'-Dichlorobenzidine	21.65	252	600506	39.608	nq	99
83) Chrysene	21.81	228	1471999	39.833	nq	99
84) Bis(2-ethylhexyl)phthalate	21.65	149	867935	40.858	nq	98
85) Di-n-octyl phthalate	22.89	149	1472347	40.657	nq	99
86) Indeno(1,2,3-cd)pyrene	28.77	276	1825348	39.887	nq	# 89
88) Benzo(b)fluoranthene	24.00	252	1596462	39.627	nq	99
89) Benzo(k)fluoranthene	24.06	252	1546907	40.648	nq	98
90) Benzo(a)pyrene	24.88	252	1511268	40.062	nq	99
91) Dibenzo(a,h)anthracene	28.84	278	1486730	40.071	nq	98
92) Benzo(g,h,i)perylene	29.95	276	1447268	39.991	nq	98

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

