

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082520\
 Data File : BG046449.D
 Acq On : 25 Aug 2020 15:28
 Operator : CG/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC050

Quant Time: Aug 25 16:23:19 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 25 15:31:23 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.94	152	80187	20.00	ng	0.00
21) Naphthalene-d8	10.73	136	310027	20.00	ng	0.00
39) Acenaphthene-d10	14.57	164	213013	20.00	ng	0.00
64) Phenanthrene-d10	17.31	188	519242	20.00	ng	0.00
76) Chrysene-d12	21.56	240	515409	20.00	ng	0.00
86) Perylene-d12	24.66	264	526581	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.52	112	425476	92.24	ng	0.00
7) Phenol-d6	7.11	99	591511	94.09	ng	0.00
23) Nitrobenzene-d5	9.09	82	512570	89.51	ng	0.00
42) 2,4,6-Tribromophenol	16.06	330	275014	84.05	ng	0.00
45) 2-Fluorobiphenyl	13.19	172	1276568	87.05	ng	0.00
79) Terphenyl-d14	19.92	244	2130384	84.18	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.43	88	85638	40.789	ng	98
3) Pyridine	3.81	79	262223	45.369	ng	100
4) n-Nitrosodimethylamine	3.73	42	100287	40.769	ng	97
6) Aniline	7.26	93	335175	44.789	ng	99
8) 2-Chlorophenol	7.50	128	220705	43.289	ng	99
9) Benzaldehyde	7.07	77	159773	42.830	ng	97
10) Phenol	7.13	94	272453	43.134	ng	99
11) bis(2-Chloroethyl)ether	7.36	93	221405	44.068	ng	99
12) 1,3-Dichlorobenzene	7.83	146	279081	45.160	ng	98
13) 1,4-Dichlorobenzene	7.97	146	281219	45.318	ng	98
14) 1,2-Dichlorobenzene	8.29	146	265848	45.425	ng	99
15) Benzyl Alcohol	8.17	79	198171	45.629	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.47	45	343184	41.664	ng	99
17) 2-Methylphenol	8.38	107	193028	45.266	ng	99
18) Hexachloroethane	9.02	117	95787	46.154	ng	96
19) n-Nitroso-di-n-propylamine	8.74	70	179613	43.900	ng	99
20) 3+4-Methylphenols	8.71	107	270916	46.416	ng	98
22) Acetophenone	8.75	105	354788	44.591	ng	99
24) Nitrobenzene	9.13	77	250684	41.062	ng	98
25) Isophorone	9.66	82	469329	41.191	ng	99
26) 2-Nitrophenol	9.84	139	137357	49.643	ng	94
27) 2,4-Dimethylphenol	9.91	122	185159	41.132	ng	99
28) bis(2-Chloroethoxy)methane	10.14	93	305117	49.249	ng	99
29) 2,4-Dichlorophenol	10.38	162	246159	46.486	ng	99
30) 1,2,4-Trichlorobenzene	10.60	180	286810	46.235	ng	98
31) Naphthalene	10.78	128	700127	42.680	ng	99
32) Benzoic acid	10.06	122	143639	51.997	ng	96
33) 4-Chloroaniline	10.89	127	320758	47.850	ng	98
34) Hexachlorobutadiene	11.08	225	188453	45.636	ng	100
35) Caprolactam	11.66	113	91363	52.713	ng	97
36) 4-Chloro-3-methylphenol	12.02	107	241453	46.464	ng	97
37) 2-Methylnaphthalene	12.39	142	551959	43.729	ng	100
38) 1-Methylnaphthalene	12.61	142	524190	43.886	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.76	216	330444	43.055	ng	100

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41) Hexachlorocyclopentadiene	12.74	237	164588	37.387	nq	100
43) 2,4,6-Trichlorophenol	13.00	196	232307	47.443	nq	98
44) 2,4,5-Trichlorophenol	13.08	196	240119	44.944	nq	99
46) 1,1'-Biphenyl	13.40	154	701570	41.985	nq	100
47) 2-Chloronaphthalene	13.44	162	589959	44.261	nq	100
48) 2-Nitroaniline	13.64	65	172842	49.470	nq	98
49) Acenaphthylene	14.29	152	897952	43.384	nq	100
50) Dimethylphthalate	14.02	163	770393	47.121	nq	99
51) 2,6-Dinitrotoluene	14.14	165	173868	45.695	nq	98
52) Acenaphthene	14.63	154	556838	40.764	nq	99
53) 3-Nitroaniline	14.47	138	191418	50.745	nq	99
54) 2,4-Dinitrophenol	14.68	184	111598	55.931	nq	98
55) Dibenzofuran	14.96	168	880372	42.711	nq	100
56) 4-Nitrophenol	14.78	139	148381	45.307	nq	99
57) 2,4-Dinitrotoluene	14.93	165	250251	47.836	nq	99
58) Fluorene	15.62	166	732169	43.862	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.19	232	234889	46.676	nq	98
60) Diethylphthalate	15.39	149	746123	46.392	nq	99
61) 4-Chlorophenyl-phenylether	15.61	204	409054	46.919	nq	98
62) 4-Nitroaniline	15.63	138	201862	53.209	nq	99
63) Azobenzene	15.90	77	609837	40.460	nq	98
65) 4,6-Dinitro-2-methylphenol	15.69	198	152533	44.418	nq	97
66) n-Nitrosodiphenylamine	15.82	169	660905	42.049	nq	99
67) 4-Bromophenyl-phenylether	16.50	248	287813	44.867	nq	98
68) Hexachlorobenzene	16.63	284	315284	43.832	nq	98
69) Atrazine	16.77	200	265483	43.865	nq	99
70) Pentachlorophenol	16.97	266	185922	39.836	nq	99
71) Phenanthrene	17.35	178	1216279	42.762	nq	99
72) Anthracene	17.44	178	1189515	42.045	nq	98
73) Carbazole	17.71	167	1135605	41.627	nq	99
74) Di-n-butylphthalate	18.28	149	1292824	45.337	nq	99
75) Fluoranthene	19.36	202	1520623	42.822	nq	98
77) Benzidine	19.54	184	581007	39.797	nq	98
78) Pyrene	19.72	202	1525099	44.039	nq	99
80) Butylbenzylphthalate	20.61	149	614126	48.955	nq	97
81) Benzo(a)anthracene	21.54	228	1477475	42.836	nq	99
82) 3,3'-Dichlorobenzidine	21.46	252	569320	39.749	nq	98
83) Chrysene	21.60	228	1418241	41.975	nq	99
84) Bis(2-ethylhexyl)phthalate	21.46	149	823134	45.896	nq	99
85) Di-n-octyl phthalate	22.63	149	1428722	47.049	nq	100
87) Indeno(1,2,3-cd)pyrene	28.19	276	1786289	44.479	nq	100
88) Benzo(b)fluoranthene	23.68	252	1551749	43.975	nq	97
89) Benzo(k)fluoranthene	23.75	252	1460959	42.211	nq	98
90) Benzo(a)pyrene	24.52	252	1445615	44.026	nq	99
91) Dibenzo(a,h)anthracene	28.25	278	1437786	42.826	nq	99
92) Benzo(g,h,i)perylene	29.30	276	1445394	42.630	nq	99

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

