

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061719\
 Data File : BG041269.D
 Acq On : 17 Jun 2019 11:11
 Operator : HP/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICCC040

Quant Time: Jun 17 11:56:39 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 17 11:50:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	30922	20.00	ng	0.00
21) Naphthalene-d8	10.91	136	131241	20.00	ng	0.00
39) Acenaphthene-d10	14.72	164	96288	20.00	ng	0.00
64) Phenanthrene-d10	17.46	188	253605	20.00	ng	0.00
76) Chrysene-d12	21.74	240	257661	20.00	ng	0.00
87) Perylene-d12	25.04	264	276625	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	137021	81.65	ng	0.00
7) Phenol-d6	7.23	99	202034	78.75	ng	0.00
23) Nitrobenzene-d5	9.26	82	248449	82.16	ng	0.00
42) 2,4,6-Tribromophenol	16.21	330	118199	88.74	ng	0.00
45) 2-Fluorobiphenyl	13.34	172	573477	78.83	ng	0.00
79) Terphenyl-d14	20.06	244	1102118	83.87	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.47	88	31872	45.558	ng	100
3) Pyridine	3.89	79	91093	44.570	ng	100
4) n-Nitrosodimethylamine	3.81	42	48560	45.673	ng	100
6) Aniline	7.41	93	141135	42.173	ng	100
8) 2-Chlorophenol	7.64	128	81962	39.475	ng	100
9) Benzaldehyde	7.22	77	61988	37.994	ng	100
10) Phenol	7.26	94	109883	39.890	ng	100
11) bis(2-Chloroethyl)ether	7.50	93	83376	41.324	ng	100
12) 1,3-Dichlorobenzene	7.97	146	101891	40.835	ng	100
13) 1,4-Dichlorobenzene	8.12	146	103026	40.598	ng	100
14) 1,2-Dichlorobenzene	8.44	146	97873	40.261	ng	100
15) Benzyl Alcohol	8.33	79	96146	40.718	ng	100
16) 2,2'-oxybis(1-Chloropropan	8.61	45	137300	42.520	ng	100
17) 2-Methylphenol	8.52	107	77896	39.143	ng	100
18) Hexachloroethane	9.16	117	40243	40.470	ng	100
19) n-Nitroso-di-n-propylamine	8.89	70	83078	42.309	ng	100
20) 3+4-Methylphenols	8.85	107	106025	38.706	ng	100
22) Acetophenone	8.92	105	150998	41.573	ng	100
24) Nitrobenzene	9.30	77	126057	41.576	ng	100
25) Isophorone	9.82	82	229878	43.127	ng	100
26) 2-Nitrophenol	10.01	139	50512	39.370	ng	100
27) 2,4-Dimethylphenol	10.06	122	74796	38.350	ng	100
28) bis(2-Chloroethoxy)methane	10.30	93	122385	42.413	ng	100
29) 2,4-Dichlorophenol	10.55	162	96666	38.721	ng	100
30) 1,2,4-Trichlorobenzene	10.76	180	121695	40.373	ng	100
31) Naphthalene	10.95	128	287158	41.269	ng	100
32) Benzoic acid	10.21	122	54142	78.099	ng	100
33) 4-Chloroaniline	11.07	127	122545	40.328	ng	100
34) Hexachlorobutadiene	11.21	225	92679	39.908	ng	100
35) Caprolactam	11.86	113	31930	44.820	ng	100
36) 4-Chloro-3-methylphenol	12.18	107	109560	41.462	ng	100
37) 2-Methylnaphthalene	12.55	142	211688	40.563	ng	100
38) 1-Methylnaphthalene	12.77	142	201718	40.987	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.91	216	160008	38.192	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.88	237	109403	59.989	nq	100
43) 2,4,6-Trichlorophenol	13.16	196	97973	38.864	nq	100
44) 2,4,5-Trichlorophenol	13.23	196	103608	41.105	nq	100
46) 1,1'-Biphenyl	13.55	154	290666	39.037	nq	100
47) 2-Chloronaphthalene	13.60	162	249908	38.447	nq	100
48) 2-Nitroaniline	13.81	65	93169	41.466	nq	100
49) Acenaphthylene	14.44	152	379621	40.702	nq	100
50) Dimethylphthalate	14.17	163	341817	42.118	nq	100
51) 2,6-Dinitrotoluene	14.30	165	71543	41.087	nq	100
52) Acenaphthene	14.78	154	221943	39.752	nq	100
53) 3-Nitroaniline	14.63	138	71692	42.611	nq	100
54) 2,4-Dinitrophenol	14.84	184	46652	135.115	nq	100
55) Dibenzofuran	15.11	168	390094	41.288	nq	100
56) 4-Nitrophenol	14.93	139	52987	58.973	nq	100
57) 2,4-Dinitrotoluene	15.08	165	103761	42.954	nq	100
58) Fluorene	15.76	166	305705	41.674	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.34	232	107578	46.528	nq	100
60) Diethylphthalate	15.52	149	346598	43.243	nq	100
61) 4-Chlorophenyl-phenylether	15.75	204	198195	41.259	nq	100
62) 4-Nitroaniline	15.80	138	75371	46.061	nq	100
63) Azobenzene	16.04	77	311163	43.075	nq	100
65) 4,6-Dinitro-2-methylphenol	15.84	198	68078	73.610	nq	100
66) n-Nitrosodiphenylamine	15.97	169	273056	38.711	nq	100
67) 4-Bromophenyl-phenylether	16.64	248	130724	38.914	nq	100
68) Hexachlorobenzene	16.76	284	141833	38.832	nq	100
69) Atrazine	16.91	200	105856	47.172	nq	100
70) Pentachlorophenol	17.11	266	76086	60.146	nq	100
71) Phenanthrene	17.51	178	542849	40.635	nq	100
72) Anthracene	17.59	178	537573	40.599	nq	100
73) Carbazole	17.87	167	466564	41.264	nq	100
74) Di-n-butylphthalate	18.40	149	586451	41.748	nq	100
75) Fluoranthene	19.51	202	725218	42.692	nq	100
77) Benzidine	19.69	184	263784	65.940	nq	100
78) Pyrene	19.87	202	729317	41.233	nq	100
80) Butylbenzylphthalate	20.74	149	268730	39.683	nq	100
81) Benzo(a)anthracene	21.72	228	714452	41.932	nq	100
82) 3,3'-Dichlorobenzidine	21.64	252	251649	48.025	nq	100
83) Chrysene	21.79	228	688784	40.899	nq	100
84) Bis(2-ethylhexyl)phthalate	21.58	149	370592	39.301	nq	100
85) Di-n-octyl phthalate	22.82	149	618455	39.507	nq	100
86) Indeno(1,2,3-cd)pyrene	28.85	276	806710	79.781	nq	100
88) Benzo(b)fluoranthene	23.99	252	690651	40.866	nq	100
89) Benzo(k)fluoranthene	24.06	252	695598	34.475	nq	100
90) Benzo(a)pyrene	24.89	252	654867	39.775	nq	100
91) Dibenzo(a,h)anthracene	28.91	278	655899	69.430	nq	100
92) Benzo(g,h,i)perylene	30.06	276	648256	79.766	nq	100

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

