

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012919\
 Data File : BG039311.D
 Acq On : 29 Jan 2019 10:15
 Operator : JU/SJ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC010

Quant Time: Jan 29 12:19:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 29 12:14:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.18	152	38155	20.00	ng	0.00
21) Naphthalene-d8	11.01	136	171498	20.00	ng	0.00
39) Acenaphthene-d10	14.81	164	114076	20.00	ng	0.00
64) Phenanthrene-d10	17.56	188	276045	20.00	ng	0.00
76) Chrysene-d12	21.85	240	287980	20.00	ng	0.00
87) Perylene-d12	25.25	264	282418	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.72	112	44941	22.34	ng	0.00
7) Phenol-d6	7.31	99	68310	23.60	ng	0.00
23) Nitrobenzene-d5	9.36	82	41614	13.28	ng	0.00
42) 2,4,6-Tribromophenol	16.29	330	20738	10.84	ng	0.00
45) 2-Fluorobiphenyl	13.43	172	165224	19.82	ng	0.00
79) Terphenyl-d14	20.15	244	295145	18.96	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.62	88	10318	13.100	ng	95
3) Pyridine	4.03	79	25281	11.812	ng	94
4) n-Nitrosodimethylamine	3.95	42	12317	12.789	ng	# 89
6) Aniline	7.51	93	41609	11.969	ng	97
8) 2-Chlorophenol	7.74	128	25562	10.781	ng	99
9) Benzaldehyde	7.32	77	23100	13.838	ng	93
10) Phenol	7.34	94	35788	12.332	ng	96
11) bis(2-Chloroethyl)ether	7.60	93	28952	13.053	ng	99
12) 1,3-Dichlorobenzene	8.07	146	31422	11.061	ng	96
13) 1,4-Dichlorobenzene	8.22	146	32157	11.177	ng	99
14) 1,2-Dichlorobenzene	8.54	146	30726	11.201	ng	95
15) Benzyl Alcohol	8.42	79	26084	11.966	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.71	45	65992	15.516	ng	96
17) 2-Methylphenol	8.60	107	22213	11.023	ng	# 89
18) Hexachloroethane	9.27	117	10524	10.767	ng	93
19) n-Nitroso-di-n-propylamine	8.99	70	24794	13.044	ng	97
20) 3+4-Methylphenols	8.93	107	30740	10.699	ng	94
22) Acetophenone	9.01	105	46053	10.283	ng	98
24) Nitrobenzene	9.40	77	23700	7.481	ng	98
25) Isophorone	9.92	82	60725	11.248	ng	99
26) 2-Nitrophenol	10.11	139	7101	4.144	ng	90
27) 2,4-Dimethylphenol	10.15	122	24834	10.302	ng	96
28) bis(2-Chloroethoxy)methane	10.40	93	37517	11.563	ng	97
29) 2,4-Dichlorophenol	10.64	162	25221	8.422	ng	94
30) 1,2,4-Trichlorobenzene	10.87	180	29913	8.313	ng	97
31) Naphthalene	11.06	128	85491	9.939	ng	98
32) Benzoic acid	10.21	122	7408	5.801	ng	96
33) 4-Chloroaniline	11.16	127	40287	10.468	ng	95
34) Hexachlorobutadiene	11.32	225	19924	8.047	ng	93
35) Caprolactam	11.93	113	10792	8.986	ng	95
36) 4-Chloro-3-methylphenol	12.26	107	29948	9.350	ng	97
37) 2-Methylnaphthalene	12.65	142	63779	9.890	ng	97
38) 1-Methylnaphthalene	12.87	142	61258	9.897	ng	100
40) 1,2,4,5-Tetrachlorobenzene	13.00	216	35514	8.290	ng	96

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41) Hexachlorocyclopentadiene	12.98	237	17877	8.487	nq	96
43) 2,4,6-Trichlorophenol	13.24	196	20611	7.644	nq	92
44) 2,4,5-Trichlorophenol	13.31	196	21504	7.546	nq	95
46) 1,1'-Biphenyl	13.65	154	94968	10.506	nq	97
47) 2-Chloronaphthalene	13.69	162	71669	9.798	nq	97
48) 2-Nitroaniline	13.90	65	10779	5.260	nq	98
49) Acenaphthylene	14.54	152	111020	10.098	nq	98
50) Dimethylphthalate	14.25	163	93583	9.708	nq	98
51) 2,6-Dinitrotoluene	14.38	165	9797	4.546	nq	94
52) Acenaphthene	14.87	154	69662	10.546	nq	92
53) 3-Nitroaniline	14.71	138	11500	5.260	nq	# 88
54) 2,4-Dinitrophenol	14.91	184	4068	3.652	nq	# 72
55) Dibenzofuran	15.21	168	114636	9.826	nq	98
56) 4-Nitrophenol	14.99	139	9826	6.246	nq	97
57) 2,4-Dinitrotoluene	15.16	165	11666	3.773	nq	95
58) Fluorene	15.85	166	88250	10.049	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.42	232	20388	7.579	nq	99
60) Diethylphthalate	15.61	149	98485	9.916	nq	99
61) 4-Chlorophenyl-phenylether	15.84	204	49155	8.885	nq	98
62) 4-Nitroaniline	15.87	138	12708	5.580	nq	94
63) Azobenzene	16.13	77	96727	12.453	nq	98
65) 4,6-Dinitro-2-methylphenol	15.91	198	6166	3.015	nq	91
66) n-Nitrosodiphenylamine	16.05	169	86640	10.587	nq	98
67) 4-Bromophenyl-phenylether	16.73	248	30785	8.676	nq	90
68) Hexachlorobenzene	16.85	284	29709	7.673	nq	97
69) Atrazine	16.99	200	33446	9.621	nq	96
70) Pentachlorophenol	17.19	266	17092	8.251	nq	97
71) Phenanthrene	17.60	178	150051	10.284	nq	98
72) Anthracene	17.69	178	146424	10.150	nq	97
73) Carbazole	17.96	167	149091	10.469	nq	99
74) Di-n-butylphthalate	18.50	149	172391	10.881	nq	99
75) Fluoranthene	19.59	202	177948	9.838	nq	97
77) Benzidine	19.78	184	101414	11.514	nq	99
78) Pyrene	19.96	202	184646	10.061	nq	98
80) Butylbenzylphthalate	20.83	149	70845	10.149	nq	99
81) Benzo(a)anthracene	21.83	228	171577	9.787	nq	97
82) 3,3'-Dichlorobenzidine	21.74	252	63738	8.923	nq	# 92
83) Chrysene	21.90	228	163724	9.820	nq	97
84) Bis(2-ethylhexyl)phthalate	21.72	149	106214	10.959	nq	98
85) Di-n-octyl phthalate	22.99	149	173259	10.572	nq	100
86) Indeno(1,2,3-cd)pyrene	29.14	276	165253	8.295	nq	99
88) Benzo(b)fluoranthene	24.15	252	153624	9.865	nq	98
89) Benzo(k)fluoranthene	24.23	252	158854	10.317	nq	98
90) Benzo(a)pyrene	25.08	252	148475	9.864	nq	97
91) Dibenzo(a,h)anthracene	29.22	278	139357	9.308	nq	97
92) Benzo(g,h,i)perylene	30.37	276	134681	9.087	nq	97

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

