

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021419\
 Data File : BG039506.D
 Acq On : 14 Feb 2019 9:57
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Feb 14 15:59:53 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 15:41:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	56160	20.00	ng	-0.01
21) Naphthalene-d8	10.99	136	269468	20.00	ng	-0.02
39) Acenaphthene-d10	14.80	164	194033	20.00	ng	-0.01
64) Phenanthrene-d10	17.54	188	432264	20.00	ng	-0.01
76) Chrysene-d12	21.83	240	417991	20.00	ng	-0.02
87) Perylene-d12	25.21	264	430173	20.00	ng	-0.03

System Monitoring Compounds						
5) 2-Fluorophenol	5.72	112	278846	84.98	ng	0.00
7) Phenol-d6	7.32	99	457724	94.77	ng	0.00
23) Nitrobenzene-d5	9.35	82	402286	93.26	ng	-0.01
42) 2,4,6-Tribromophenol	16.28	330	182060	87.13	ng	0.00
45) 2-Fluorobiphenyl	13.42	172	1023557	75.67	ng	-0.01
79) Terphenyl-d14	20.13	244	1489660	75.44	ng	-0.02

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.61	88	50465	35.159	ng	94
3) Pyridine	4.01	79	148546	39.089	ng	93
4) n-Nitrosodimethylamine	3.93	42	68054	35.808	ng	90
6) Aniline	7.50	93	254786	42.114	ng	96
8) 2-Chlorophenol	7.73	128	161757	45.098	ng	97
9) Benzaldehyde	7.31	77	104068	44.553	ng	96
10) Phenol	7.34	94	238171	47.304	ng	97
11) bis(2-Chloroethyl)ether	7.58	93	162053	40.732	ng	97
12) 1,3-Dichlorobenzene	8.06	146	171828	39.545	ng	96
13) 1,4-Dichlorobenzene	8.21	146	176172	40.678	ng	98
14) 1,2-Dichlorobenzene	8.53	146	170379	40.638	ng	98
15) Benzyl Alcohol	8.41	79	169527	44.707	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.68	45	303754	33.809	ng	100
17) 2-Methylphenol	8.61	107	153287	47.046	ng	97
18) Hexachloroethane	9.26	117	61358	41.180	ng	95
19) n-Nitroso-di-n-propylamine	8.98	70	139221	40.611	ng	93
20) 3+4-Methylphenols	8.93	107	222832	49.664	ng	97
22) Acetophenone	9.00	105	278017	38.409	ng	# 94
24) Nitrobenzene	9.39	77	204471	43.878	ng	95
25) Isophorone	9.90	82	357784	37.431	ng	97
26) 2-Nitrophenol	10.10	139	108220	56.165	ng	96
27) 2,4-Dimethylphenol	10.14	122	169929	43.947	ng	94
28) bis(2-Chloroethoxy)methane	10.39	93	232778	39.537	ng	99
29) 2,4-Dichlorophenol	10.63	162	208039	49.228	ng	96
30) 1,2,4-Trichlorobenzene	10.85	180	196662	41.450	ng	99
31) Naphthalene	11.04	128	522048	39.052	ng	99
32) Benzoic acid	10.28	122	129071	50.417	ng	95
33) 4-Chloroaniline	11.16	127	265304	42.802	ng	97
34) Hexachlorobutadiene	11.31	225	126661	40.143	ng	93
35) Caprolactam	11.95	113	80195	45.858	ng	# 91
36) 4-Chloro-3-methylphenol	12.25	107	223479	45.726	ng	98
37) 2-Methylnaphthalene	12.64	142	396051	40.000	ng	97
38) 1-Methylnaphthalene	12.85	142	388226	40.676	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.99	216	246408	39.913	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.96	237	142848	42.463	nq	98
43) 2,4,6-Trichlorophenol	13.23	196	175858	46.329	nq	97
44) 2,4,5-Trichlorophenol	13.31	196	183889	46.223	nq	98
46) 1,1'-Biphenyl	13.63	154	607705	37.643	nq	99
47) 2-Chloronaphthalene	13.68	162	475316	39.137	nq	98
48) 2-Nitroaniline	13.89	65	161952	46.694	nq	93
49) Acenaphthylene	14.52	152	701955	38.305	nq	100
50) Dimethylphthalate	14.25	163	595818	38.021	nq	100
51) 2,6-Dinitrotoluene	14.37	165	137994	47.237	nq	94
52) Acenaphthene	14.86	154	443154	37.254	nq	97
53) 3-Nitroaniline	14.71	138	141556	45.180	nq	# 91
54) 2,4-Dinitrophenol	14.91	184	47464	46.376	nq	89
55) Dibenzofuran	15.19	168	733045	38.435	nq	99
56) 4-Nitrophenol	15.00	139	107078	40.848	nq	# 87
57) 2,4-Dinitrotoluene	15.16	165	188409	49.858	nq	# 86
58) Fluorene	15.84	166	539046	37.914	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.41	232	160000	42.953	nq	95
60) Diethylphthalate	15.60	149	589197	36.364	nq	97
61) 4-Chlorophenyl-phenylether	15.83	204	322768	40.093	nq	94
62) 4-Nitroaniline	15.87	138	137955	42.508	nq	95
63) Azobenzene	16.12	77	547722	34.586	nq	95
65) 4,6-Dinitro-2-methylphenol	15.91	198	83365	48.675	nq	90
66) n-Nitrosodiphenylamine	16.04	169	527871	40.162	nq	99
67) 4-Bromophenyl-phenylether	16.72	248	203351	42.405	nq	93
68) Hexachlorobenzene	16.83	284	210160	43.680	nq	96
69) Atrazine	16.98	200	179248	37.318	nq	96
70) Pentachlorophenol	17.18	266	128836	38.528	nq	98
71) Phenanthrene	17.58	178	857185	38.314	nq	99
72) Anthracene	17.68	178	849036	38.527	nq	99
73) Carbazole	17.95	167	814729	37.045	nq	99
74) Di-n-butylphthalate	18.47	149	955594	37.244	nq	99
75) Fluoranthene	19.59	202	990863	38.157	nq	99
77) Benzidine	19.76	184	510887	37.280	nq	98
78) Pyrene	19.94	202	1000749	38.459	nq	99
80) Butylbenzylphthalate	20.81	149	440127	40.097	nq	94
81) Benzo(a)anthracene	21.81	228	957399	38.763	nq	99
82) 3,3'-Dichlorobenzidine	21.72	252	371613	40.577	nq	99
83) Chrysene	21.88	228	909981	38.703	nq	99
84) Bis(2-ethylhexyl)phthalate	21.68	149	614782	39.259	nq	# 99
85) Di-n-octyl phthalate	22.95	149	1028247	39.745	nq	96
86) Indeno(1,2,3-cd)pyrene	29.10	276	1059591	42.004	nq	99
88) Benzo(b)fluoranthene	24.13	252	940974	40.110	nq	98
89) Benzo(k)fluoranthene	24.20	252	895961	38.040	nq	98
90) Benzo(a)pyrene	25.05	252	906074	40.103	nq	98
91) Dibenzo(a,h)anthracene	29.17	278	858321	40.566	nq	99
92) Benzo(g,h,i)perylene	30.33	276	867852	41.151	nq	97

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

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