

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111618\
 Data File : BG038048.D
 Acq On : 17 Nov 2018 00:31
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 11/19/2018 2:13:49 PM

Quant Time: Nov 17 03:36:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 16:39:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.09	152	50218	20.00	ng	0.00
21) Naphthalene-d8	10.91	136	217731	20.00	ng	0.00
39) Acenaphthene-d10	14.72	164	132497	20.00	ng	0.00
64) Phenanthrene-d10	17.47	188	299324	20.00	ng	0.00
76) Chrysene-d12	21.75	240	272124	20.00	ng	0.00
87) Perylene-d12	25.07	264	289394	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.63	112	246714	84.82	ng	0.00
7) Phenol-d6	7.23	99	348584	83.96	ng	0.00
23) Nitrobenzene-d5	9.27	82	343678	84.50	ng	0.00
42) 2,4,6-Tribromophenol	16.21	330	119900	79.35	ng	0.00
45) 2-Fluorobiphenyl	13.34	172	744097	81.82	ng	0.00
79) Terphenyl-d14	20.07	244	963807	83.34	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	56720	41.286	ng	97
3) Pyridine	3.93	79	160981	41.078	ng	95
4) n-Nitrosodimethylamine	3.84	42	64312	45.234	ng	# 92
6) Aniline	7.42	93	219747	41.009	ng	97
8) 2-Chlorophenol	7.65	128	140834	41.730	ng	99
9) Benzaldehyde	7.23	77	122395	40.765	ng	99
10) Phenol	7.26	94	185743	42.237	ng	98
11) bis(2-Chloroethyl)ether	7.50	93	150875	42.024	ng	97
12) 1,3-Dichlorobenzene	7.97	146	161768	41.608	ng	98
13) 1,4-Dichlorobenzene	8.12	146	161623	41.152	ng	99
14) 1,2-Dichlorobenzene	8.44	146	154576	41.224	ng	99
15) Benzyl Alcohol	8.33	79	132274	44.030	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.60	45	288320	43.249	ng	96
17) 2-Methylphenol	8.52	107	124993	42.040	ng	98
18) Hexachloroethane	9.17	117	60593	42.392	ng	95
19) n-Nitroso-di-n-propylamine	8.89	70	121130	40.321	ng	98
20) 3+4-Methylphenols	8.85	107	175593	41.661	ng	99
22) Acetophenone	8.91	105	222779	41.124	ng	100
24) Nitrobenzene	9.31	77	178866	42.988	ng	96
25) Isophorone	9.82	82	299685	40.935	ng	96
26) 2-Nitrophenol	10.01	139	86511	41.648	ng	93
27) 2,4-Dimethylphenol	10.06	122	131707	42.083	ng	96
28) bis(2-Chloroethoxy)methane	10.30	93	196409	41.955	ng	97
29) 2,4-Dichlorophenol	10.55	162	147526	41.907	ng	99
30) 1,2,4-Trichlorobenzene	10.76	180	162615	41.223	ng	97
31) Naphthalene	10.96	128	442532	41.323	ng	99
32) Benzoic acid	10.19	122	88136m	44.832	ng	
33) 4-Chloroaniline	11.07	127	200383	40.226	ng	97
34) Hexachlorobutadiene	11.22	225	99782	41.099	ng	96
35) Caprolactam	11.86	113	55387	39.308	ng	93
36) 4-Chloro-3-methylphenol	12.18	107	159838	41.560	ng	97
37) 2-Methylnaphthalene	12.56	142	311084	40.440	ng	100
38) 1-Methylnaphthalene	12.77	142	299908	40.504	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.92	216	186190	42.053	ng	98

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111618\
 Data File : BG038048.D
 Acq On : 17 Nov 2018 00:31
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 11/19/2018 2:13:49 PM

Quant Time: Nov 17 03:36:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 16:39:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.88	237	97287	41.036	nq	100
43) 2,4,6-Trichlorophenol	13.16	196	122932	40.746	nq	98
44) 2,4,5-Trichlorophenol	13.23	196	127106	40.040	nq	96
46) 1,1'-Biphenyl	13.56	154	456402	41.002	nq	100
47) 2-Chloronaphthalene	13.60	162	346785	40.827	nq	99
48) 2-Nitroaniline	13.81	65	116755	43.674	nq	91
49) Acenaphthylene	14.45	152	528217	41.353	nq	100
50) Dimethylphthalate	14.17	163	421778	40.154	nq	99
51) 2,6-Dinitrotoluene	14.30	165	98381	41.383	nq	98
52) Acenaphthene	14.78	154	319370	41.532	nq	98
53) 3-Nitroaniline	14.64	138	108327	41.294	nq	# 95
54) 2,4-Dinitrophenol	14.84	184	54668	46.947	nq	# 86
55) Dibenzofuran	15.12	168	517421	40.693	nq	99
56) 4-Nitrophenol	14.93	139	79045	39.069	nq	99
57) 2,4-Dinitrotoluene	15.08	165	136404	42.170	nq	99
58) Fluorene	15.77	166	386501	40.739	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.34	232	108468	40.834	nq	98
60) Diethylphthalate	15.52	149	446531	40.023	nq	99
61) 4-Chlorophenyl-phenylether	15.75	204	222561	40.092	nq	98
62) 4-Nitroaniline	15.80	138	109287	41.937	nq	95
63) Azobenzene	16.05	77	415077	41.610	nq	98
65) 4,6-Dinitro-2-methylphenol	15.84	198	73113	38.618	nq	96
66) n-Nitrosodiphenylamine	15.97	169	375188	41.433	nq	99
67) 4-Bromophenyl-phenylether	16.65	248	135850	41.350	nq	98
68) Hexachlorobenzene	16.76	284	136900	40.370	nq	99
69) Atrazine	16.91	200	142588	42.715	nq	97
70) Pentachlorophenol	17.11	266	93907	40.774	nq	92
71) Phenanthrene	17.52	178	626504	40.881	nq	99
72) Anthracene	17.60	178	628156	41.582	nq	99
73) Carbazole	17.87	167	628745	41.484	nq	99
74) Di-n-butylphthalate	18.41	149	715312	42.044	nq	99
75) Fluoranthene	19.52	202	723215	41.363	nq	99
77) Benzidine	19.70	184	396717	41.547	nq	100
78) Pyrene	19.88	202	737586	41.457	nq	99
80) Butylbenzylphthalate	20.75	149	328331	42.199	nq	99
81) Benzo(a)anthracene	21.73	228	682450	41.500	nq	99
82) 3,3'-Dichlorobenzidine	21.65	252	266827	41.654	nq	99
83) Chrysene	21.80	228	654591	41.604	nq	99
84) Bis(2-ethylhexyl)phthalate	21.61	149	466399	42.034	nq	98
85) Di-n-octyl phthalate	22.84	149	772277	43.022	nq	100
86) Indeno(1,2,3-cd)pyrene	28.88	276	744656	42.613	nq	# 90
88) Benzo(b)fluoranthene	24.01	252	662852	41.622	nq	99
89) Benzo(k)fluoranthene	24.08	252	667800	42.496	nq	99
90) Benzo(a)pyrene	24.92	252	646740	42.494	nq	100
91) Dibenzo(a,h)anthracene	28.95	278	617519	42.169	nq	95
92) Benzo(g,h,i)perylene	30.08	276	612080	42.036	nq	98

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

