

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG033018\
 Data File : BG033469.D
 Acq On : 31 Mar 2018 5:36
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 4/2/2018 5:07:21 PM

Quant Time: Mar 31 06:14:36 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 28 18:34:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.87	152	35957	20.00	ng	0.00
21) Naphthalene-d8	11.76	136	172787	20.00	ng	0.00
38) Acenaphthene-d10	15.49	164	137102	20.00	ng	0.00
63) Phenanthrene-d10	18.21	188	346488	20.00	ng	0.00
75) Chrysene-d12	22.56	240	353832	20.00	ng	0.00
86) Perylene-d12	26.33	264	391616	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.30	112	156517	81.62	ng	0.00
7) Phenol-d6	7.99	99	278783	84.61	ng	0.01
23) Nitrobenzene-d5	10.07	82	291485	77.51	ng	0.00
41) 2,4,6-Tribromophenol	16.96	330	155994	76.08	ng	0.00
44) 2-Fluorobiphenyl	14.11	172	717800	75.58	ng	0.00
78) Terphenyl-d14	20.73	244	1223905	73.97	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.93	88	37399	38.404	ng	89
3) Pyridine	4.39	79	103997	38.632	ng	97
4) n-Nitrosodimethylamine	4.30	42	46807m	42.369	ng	
6) Aniline	8.17	93	160369	41.389	ng	96
8) 2-Chlorophenol	8.42	128	89168	40.675	ng	98
9) Benzaldehyde	7.97	77	68783	32.850	ng	91
10) Phenol	8.02	94	136186	41.865	ng	93
11) bis(2-Chloroethyl)ether	8.25	93	109502	41.241	ng	97
12) 1,3-Dichlorobenzene	8.75	146	108610	37.895	ng	95
13) 1,4-Dichlorobenzene	8.91	146	108537	37.813	ng	92
14) 1,2-Dichlorobenzene	9.24	146	109175	38.971	ng	98
15) Benzyl Alcohol	9.11	79	111922	43.145	ng	92
16) 2,2'-oxybis(1-Chloropropan	9.39	45	175699	40.591	ng	88
17) 2-Methylphenol	9.31	107	94702m	42.735	ng	
18) Hexachloroethane	9.99	117	43726	39.470	ng	92
19) n-Nitroso-di-n-propylamine	9.68	70	106153	43.968	ng	96
20) 3+4-Methylphenols	9.65	107	141643	44.892	ng	91
22) Acetophenone	9.71	105	186252	39.345	ng	# 97
24) Nitrobenzene	10.12	77	152450	37.872	ng	93
25) Isophorone	10.63	82	289577	39.673	ng	98
26) 2-Nitrophenol	10.85	139	57106	37.471	ng	# 86
27) 2,4-Dimethylphenol	10.88	122	87888	39.698	ng	87
28) bis(2-Chloroethoxy)methane	11.11	93	136267	37.416	ng	97
29) 2,4-Dichlorophenol	11.39	162	109559	40.239	ng	92
30) 1,2,4-Trichlorobenzene	11.60	180	132335	37.410	ng	97
31) Naphthalene	11.80	128	341826	38.176	ng	98
32) Benzoic acid	11.01	122	67817m	32.951	ng	
33) 4-Chloroaniline	11.90	127	158948	39.623	ng	# 93
34) Hexachlorobutadiene	12.06	225	101224	37.839	ng	97
35) Caprolactam	12.66	113	45779	42.918	ng	97
36) 4-Chloro-3-methylphenol	12.97	107	149184	42.010	ng	98
37) 2-Methylnaphthalene	13.35	142	266169	38.299	ng	94
39) 1,2,4,5-Tetrachlorobenzene	13.71	216	182803	36.809	ng	97
40) Hexachlorocyclopentadiene	13.68	237	96628	33.191	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.94	196	115339	39.974	ng	93
43) 2,4,5-Trichlorophenol	14.02	196	143557	42.093	ng	97
45) 1,1'-Biphenyl	14.33	154	406043	38.549	ng	97
46) 2-Chloronaphthalene	14.38	162	307809	37.900	ng	94
47) 2-Nitroaniline	14.58	65	126426	41.560	ng	98
48) Acenaphthylene	15.22	152	526917	38.868	ng	97
49) Dimethylphthalate	14.91	163	464573	38.936	ng	99
50) 2,6-Dinitrotoluene	15.05	165	97392	39.920	ng	98
51) Acenaphthene	15.55	154	335548	38.396	ng	98
52) 3-Nitroaniline	15.39	138	97186	37.997	ng	# 96
53) 2,4-Dinitrophenol	15.59	184	22621	23.837	ng	85
54) Dibenzofuran	15.88	168	497385	38.531	ng	91
55) 4-Nitrophenol	15.68	139	63027	35.043	ng	91
56) 2,4-Dinitrotoluene	15.82	165	142303	39.615	ng	91
57) Fluorene	16.52	166	427111	38.838	ng	98
58) 2,3,4,6-Tetrachlorophenol	16.09	232	128477	40.015	ng	95
59) Diethylphthalate	16.25	149	452841	39.086	ng	98
60) 4-Chlorophenyl-phenylether	16.49	204	239009	37.807	ng	96
61) 4-Nitroaniline	16.54	138	103900	40.113	ng	88
62) Azobenzene	16.79	77	439985	39.203	ng	98
64) 4,6-Dinitro-2-methylphenol	16.58	198	71037	34.271	ng	96
65) n-Nitrosodiphenylamine	16.71	169	397700	40.205	ng	97
66) 4-Bromophenyl-phenylether	17.39	248	158177	38.872	ng	96
67) Hexachlorobenzene	17.51	284	166301	38.724	ng	96
68) Atrazine	17.63	200	155991	38.917	ng	98
69) Pentachlorophenol	17.86	266	121759	41.309	ng	97
70) Phenanthrene	18.26	178	699670	38.773	ng	99
71) Anthracene	18.35	178	713723	39.063	ng	99
72) Carbazole	18.61	167	637096	39.349	ng	97
73) Di-n-butylphthalate	19.10	149	755619	40.096	ng	# 98
74) Fluoranthene	20.21	202	853742	38.232	ng	97
76) Benzidine	20.37	184	385113	40.135	ng	98
77) Pyrene	20.57	202	876395	37.138	ng	98
79) Butylbenzylphthalate	21.42	149	343730	39.471	ng	98
80) Benzo(a)anthracene	22.54	228	861037	38.216	ng	99
81) 3,3'-Dichlorobenzidine	22.43	252	333710	41.623	ng	# 98
82) Chrysene	22.62	228	841863	38.601	ng	99
83) Bis(2-ethylhexyl)phthalate	22.36	149	488809	40.719	ng	98
84) Di-n-octyl phthalate	23.76	149	778759	42.369	ng	98
85) Indeno(1,2,3-cd)pyrene	30.71	276	967009	41.009	ng	# 77
87) Benzo(b)fluoranthene	25.12	252	846257	37.403	ng	# 97
88) Benzo(k)fluoranthene	25.20	252	856890	37.446	ng	# 96
89) Benzo(a)pyrene	26.15	252	825379	37.987	ng	# 97
90) Dibenzo(a,h)anthracene	30.79	278	796853	38.934	ng	# 97
91) Benzo(g,h,i)perylene	32.10	276	782909	38.630	ng	# 95

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

