

Data Path : Z:\HPCHEM1\BNA G\DATA\BG031418\
 Data File : BG033143.D
 Acq On : 14 Mar 2018 11:57
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 3/15/2018 6:59:57 PM

Quant Time: Mar 14 14:02:00 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 14 13:59:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.90	152	50732	20.00	ng	0.00
21) Naphthalene-d8	11.79	136	243586	20.00	ng	0.00
38) Acenaphthene-d10	15.52	164	146772	20.00	ng	0.00
63) Phenanthrene-d10	18.25	188	338958	20.00	ng	0.00
75) Chrysene-d12	22.61	240	295934	20.00	ng	0.00
86) Perylene-d12	26.41	264	317210	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.32	112	234801	74.05	ng	0.00
7) Phenol-d6	8.01	99	330367	74.74	ng	0.00
23) Nitrobenzene-d5	10.10	82	324139	91.49	ng	0.00
41) 2,4,6-Tribromophenol	16.99	330	152515	98.83	ng	0.00
44) 2-Fluorobiphenyl	14.14	172	739127	72.63	ng	0.00
78) Terphenyl-d14	20.76	244	990349	75.19	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.95	88	45640	33.163	ng	94
3) Pyridine	4.42	79	137353	36.211	ng	90
4) n-Nitrosodimethylamine	4.32	42	68548	45.075	ng	# 79
6) Aniline	8.19	93	219255	36.647	ng	99
8) 2-Chlorophenol	8.45	128	129752	37.383	ng	98
9) Benzaldehyde	8.00	77	100879	39.600	ng	91
10) Phenol	8.04	94	169741	38.096	ng	98
11) bis(2-Chloroethyl)ether	8.28	93	128666	35.315	ng	96
12) 1,3-Dichlorobenzene	8.78	146	149477	36.769	ng	99
13) 1,4-Dichlorobenzene	8.93	146	149412	36.512	ng	99
14) 1,2-Dichlorobenzene	9.27	146	145670	37.148	ng	94
15) Benzyl Alcohol	9.13	79	132732	40.848	ng	92
16) 2,2'-oxybis(1-Chloropropan	9.42	45	266206	42.503	ng	99
17) 2-Methylphenol	9.34	107	124685	37.949	ng	100
18) Hexachloroethane	10.03	117	54890	39.396	ng	# 87
19) n-Nitroso-di-n-propylamine	9.71	70	122332	39.204	ng	# 90
20) 3+4-Methylphenols	9.67	107	178982	39.179	ng	95
22) Acetophenone	9.74	105	226012	36.740	ng	# 96
24) Nitrobenzene	10.14	77	167271	43.717	ng	95
25) Isophorone	10.67	82	326688	38.210	ng	96
26) 2-Nitrophenol	10.87	139	81032	55.798	ng	95
27) 2,4-Dimethylphenol	10.91	122	131983	35.536	ng	92
28) bis(2-Chloroethoxy)methane	11.14	93	175734	35.283	ng	97
29) 2,4-Dichlorophenol	11.42	162	130768	38.793	ng	95
30) 1,2,4-Trichlorobenzene	11.63	180	142845	36.151	ng	96
31) Naphthalene	11.84	128	464262	36.475	ng	99
32) Benzoic acid	10.97	122	37939m	23.118	ng	
33) 4-Chloroaniline	11.93	127	207173	36.403	ng	98
34) Hexachlorobutadiene	12.10	225	85458	38.557	ng	96
35) Caprolactam	12.69	113	57219	42.392	ng	99
36) 4-Chloro-3-methylphenol	13.00	107	163595	41.704	ng	97
37) 2-Methylnaphthalene	13.39	142	341272	37.060	ng	93
39) 1,2,4,5-Tetrachlorobenzene	13.74	216	163621	37.554	ng	# 98
40) Hexachlorocyclopentadiene	13.71	237	90552	40.781	ng	88

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.97	196	114094	41.524	ng	100
43) 2,4,5-Trichlorophenol	14.04	196	129561	40.741	ng	95
45) 1,1'-Biphenyl	14.36	154	455997	35.554	ng	96
46) 2-Chloronaphthalene	14.41	162	339625	35.449	ng	94
47) 2-Nitroaniline	14.60	65	126509	51.466	ng	95
48) Acenaphthylene	15.25	152	573945	36.591	ng	98
49) Dimethylphthalate	14.94	163	460109	36.920	ng	99
50) 2,6-Dinitrotoluene	15.08	165	97391	47.361	ng	96
51) Acenaphthene	15.58	154	362870	37.146	ng	99
52) 3-Nitroaniline	15.41	138	113490	45.028	ng	# 90
53) 2,4-Dinitrophenol	15.61	184	30942	54.522	ng	# 64
54) Dibenzofuran	15.91	168	507625	36.495	ng	93
55) 4-Nitrophenol	15.69	139	85209	45.138	ng	95
56) 2,4-Dinitrotoluene	15.85	165	137265	54.849	ng	89
57) Fluorene	16.55	166	423392	36.879	ng	98
58) 2,3,4,6-Tetrachlorophenol	16.12	232	113530	45.982	ng	91
59) Diethylphthalate	16.28	149	502110	38.199	ng	98
60) 4-Chlorophenyl-phenylether	16.53	204	213561	38.026	ng	91
61) 4-Nitroaniline	16.56	138	119460	44.690	ng	96
62) Azobenzene	16.82	77	437212	37.180	ng	95
64) 4,6-Dinitro-2-methylphenol	16.61	198	64936	59.812	ng	93
65) n-Nitrosodiphenylamine	16.74	169	393199	35.659	ng	99
66) 4-Bromophenyl-phenylether	17.42	248	135079	37.688	ng	# 90
67) Hexachlorobenzene	17.55	284	147536	38.091	ng	95
68) Atrazine	17.66	200	128870	35.505	ng	99
69) Pentachlorophenol	17.89	266	95302	39.064	ng	100
70) Phenanthrene	18.29	178	678647	35.887	ng	98
71) Anthracene	18.38	178	685340	36.099	ng	100
72) Carbazole	18.64	167	660931	35.319	ng	98
73) Di-n-butylphthalate	19.13	149	830241	36.165	ng	# 98
74) Fluoranthene	20.24	202	737083	35.286	ng	96
76) Benzidine	20.40	184	422865	41.920	ng	99
77) Pyrene	20.60	202	757416	36.293	ng	99
79) Butylbenzylphthalate	21.46	149	353923	38.250	ng	93
80) Benzo(a)anthracene	22.58	228	686222	36.161	ng	99
81) 3,3'-Dichlorobenzidine	22.47	252	264928	37.694	ng	# 99
82) Chrysene	22.66	228	661161	36.473	ng	97
83) Bis(2-ethylhexyl)phthalate	22.40	149	503363	36.752	ng	94
84) Di-n-octyl phthalate	23.82	149	779095	37.319	ng	99
85) Indeno(1,2,3-cd)pyrene	30.83	276	784129	37.090	ng	95
87) Benzo(b)fluoranthene	25.19	252	673338	35.901	ng	99
88) Benzo(k)fluoranthene	25.27	252	673924	36.155	ng	# 97
89) Benzo(a)pyrene	26.23	252	645625	36.191	ng	# 97
90) Dibenzo(a,h)anthracene	30.91	278	656572	36.565	ng	# 96
91) Benzo(g,h,i)perylene	32.24	276	640969	36.211	ng	97

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

