

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101818\
 Data File : BG037412.D
 Acq On : 18 Oct 2018 9:53
 Operator : JU/SJ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Sohil
 10/19/2018 3:57:41 PM

Quant Time: Oct 18 12:33:39 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 18 12:15:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	53095	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	247750	20.00	ng	0.00
39) Acenaphthene-d10	14.74	164	169944	20.00	ng	0.00
64) Phenanthrene-d10	17.49	188	406851	20.00	ng	0.00
76) Chrysene-d12	21.77	240	381994	20.00	ng	0.00
87) Perylene-d12	25.11	264	390137	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.64	112	63947	21.20	ng	0.00
7) Phenol-d6	7.24	99	97821	21.36	ng	0.00
23) Nitrobenzene-d5	9.28	82	88054	23.37	ng	0.00
42) 2,4,6-Tribromophenol	16.23	330	36286	21.77	ng	0.00
45) 2-Fluorobiphenyl	13.36	172	237781	21.01	ng	0.00
79) Terphenyl-d14	20.09	244	393946	21.65	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	16996	10.031	ng	92
3) Pyridine	3.94	79	39320	9.774	ng	93
4) n-Nitrosodimethylamine	3.86	42	14118	9.037	ng	# 82
6) Aniline	7.43	93	58651	9.784	ng	96
8) 2-Chlorophenol	7.67	128	38520	10.909	ng	98
9) Benzaldehyde	7.25	77	34290	10.876	ng	97
10) Phenol	7.27	94	52049	10.792	ng	93
11) bis(2-Chloroethyl)ether	7.52	93	39175	9.943	ng	91
12) 1,3-Dichlorobenzene	7.99	146	42580	10.258	ng	96
13) 1,4-Dichlorobenzene	8.15	146	43956	10.459	ng	97
14) 1,2-Dichlorobenzene	8.46	146	42393	10.410	ng	98
15) Benzyl Alcohol	8.34	79	34373	9.596	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.63	45	70404	9.039	ng	97
17) 2-Methylphenol	8.53	107	32995	10.215	ng	97
18) Hexachloroethane	9.19	117	14337	9.995	ng	98
19) n-Nitroso-di-n-propylamine	8.90	70	31799	9.172	ng	94
20) 3+4-Methylphenols	8.87	107	49884	11.045	ng	91
22) Acetophenone	8.93	105	63791	10.214	ng	# 97
24) Nitrobenzene	9.33	77	46833	11.603	ng	99
25) Isophorone	9.84	82	78628	9.140	ng	97
26) 2-Nitrophenol	10.04	139	17820	12.408	ng	97
27) 2,4-Dimethylphenol	10.08	122	37786	10.509	ng	97
28) bis(2-Chloroethoxy)methane	10.33	93	53556	9.951	ng	95
29) 2,4-Dichlorophenol	10.57	162	41179	11.282	ng	98
30) 1,2,4-Trichlorobenzene	10.78	180	44538	10.027	ng	95
31) Naphthalene	10.98	128	125404	10.415	ng	98
32) Benzoic acid	10.14	122	19869m	10.007	ng	
33) 4-Chloroaniline	11.09	127	58959	10.439	ng	93
34) Hexachlorobutadiene	11.25	225	25815	9.359	ng	95
35) Caprolactam	11.85	113	16706	11.227	ng	88
36) 4-Chloro-3-methylphenol	12.18	107	47045	10.913	ng	97
37) 2-Methylnaphthalene	12.58	142	89855	10.226	ng	98
38) 1-Methylnaphthalene	12.79	142	89490	10.428	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.94	216	54371	9.879	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.91	237	21707	9.568	nq	98
43) 2,4,6-Trichlorophenol	13.17	196	36059	11.294	nq	94
44) 2,4,5-Trichlorophenol	13.24	196	39049	11.368	nq	95
46) 1,1'-Biphenyl	13.58	154	144077	10.306	nq	98
47) 2-Chloronaphthalene	13.62	162	107844	10.213	nq	99
48) 2-Nitroaniline	13.83	65	31083	12.181	nq	95
49) Acenaphthylene	14.47	152	166694	10.319	nq	98
50) Dimethylphthalate	14.19	163	133827	10.020	nq	100
51) 2,6-Dinitrotoluene	14.32	165	28951	12.011	nq	95
52) Acenaphthene	14.80	154	101224	10.145	nq	98
53) 3-Nitroaniline	14.65	138	30775	11.013	nq #	97
54) 2,4-Dinitrophenol	14.85	184	2719	4.069	nq #	81
55) Dibenzofuran	15.14	168	171710	10.606	nq	97
56) 4-Nitrophenol	14.93	139	18490	8.489	nq	98
57) 2,4-Dinitrotoluene	15.10	165	34837	11.504	nq #	96
58) Fluorene	15.78	166	129641	10.587	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.36	232	33793	11.111	nq #	98
60) Diethylphthalate	15.54	149	138157	9.982	nq	97
61) 4-Chlorophenyl-phenylether	15.77	204	73405	10.256	nq	95
62) 4-Nitroaniline	15.81	138	31208	10.641	nq	91
63) Azobenzene	16.07	77	136580	10.292	nq	97
65) 4,6-Dinitro-2-methylphenol	15.86	198	11005	9.707	nq	91
66) n-Nitrosodiphenylamine	15.99	169	125830	10.532	nq	98
67) 4-Bromophenyl-phenylether	16.67	248	44320	10.094	nq	94
68) Hexachlorobenzene	16.78	284	46713	10.261	nq	99
69) Atrazine	16.93	200	45711	10.294	nq	99
70) Pentachlorophenol	17.12	266	23758	8.089	nq	99
71) Phenanthrene	17.53	178	219739	10.610	nq	98
72) Anthracene	17.62	178	217656	10.737	nq	97
73) Carbazole	17.89	167	214850	10.287	nq	98
74) Di-n-butylphthalate	18.43	149	237699	10.618	nq	98
75) Fluoranthene	19.53	202	251177	10.377	nq	98
77) Benzidine	19.72	184	132118	9.041	nq	98
78) Pyrene	19.90	202	261276	10.495	nq	97
80) Butylbenzylphthalate	20.77	149	97982	10.116	nq	95
81) Benzo(a)anthracene	21.75	228	232486	10.162	nq	98
82) 3,3'-Dichlorobenzidine	21.67	252	87188	9.879	nq	97
83) Chrysene	21.82	228	223031	10.220	nq	96
84) Bis(2-ethylhexyl)phthalate	21.63	149	137184	9.863	nq	100
85) Di-n-octyl phthalate	22.88	149	222064	9.836	nq	97
86) Indeno(1,2,3-cd)pyrene	28.92	276	224329m	9.274	nq	
88) Benzo(b)fluoranthene	24.04	252	213172	10.050	nq	98
89) Benzo(k)fluoranthene	24.10	252	212000	10.129	nq	98
90) Benzo(a)pyrene	24.95	252	202538	9.958	nq	98
91) Dibenzo(a,h)anthracene	29.00	278	179357	9.216	nq	97
92) Benzo(g,h,i)perylene	30.13	276	182391	9.384	nq	96

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

