

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110922\
 Data File : BG055519.D
 Acq On : 9 Nov 2022 11:47
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/10/2022
 Supervised By : Jagrut Upadhyay 11/10/2022

Quant Time: Nov 10 02:53:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 10 02:52:33 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.270	152	44149	20.000	ng	0.00
21) Naphthalene-d8	11.096	136	190054	20.000	ng	0.00
39) Acenaphthene-d10	14.886	164	140602	20.000	ng	0.00
64) Phenanthrene-d10	17.630	188	330879	20.000	ng	0.00
76) Chrysene-d12	21.925	240	324741	20.000	ng	# 0.00
86) Perylene-d12	25.344	264	339937	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.791	112	45721	19.725	ng	0.00
7) Phenol-d6	7.395	99	63296	19.769	ng	0.00
23) Nitrobenzene-d5	9.439	82	48284	19.355	ng	0.00
42) 2,4,6-Tribromophenol	16.367	330	23147	19.212	ng	0.00
45) 2-Fluorobiphenyl	13.517	172	192925	20.697	ng	0.00
79) Terphenyl-d14	20.215	244	343363	21.237	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.634	88	12512	10.628	ng	96
3) Pyridine	4.052	79	31918	9.905	ng	94
4) n-Nitrosodimethylamine	3.958	42	17638	9.899	ng	# 91
6) Aniline	7.583	93	43271	10.175	ng	94
8) 2-Chlorophenol	7.824	128	24593	9.477	ng	96
9) Benzaldehyde	7.395	77	27971	11.073	ng	96
10) Phenol	7.418	94	36746	9.969	ng	96
11) bis(2-Chloroethyl)ether	7.671	93	30670	10.335	ng	99
12) 1,3-Dichlorobenzene	8.159	146	33113	10.215	ng	91
13) 1,4-Dichlorobenzene	8.305	146	33252	10.136	ng	99
14) 1,2-Dichlorobenzene	8.629	146	32006	10.274	ng	95
15) Benzyl Alcohol	8.499	79	27255	9.762	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.793	45	59128	10.369	ng	98
17) 2-Methylphenol	8.693	107	26049	10.043	ng	99
18) Hexachloroethane	9.369	117	10489	9.631	ng	93
19) n-Nitroso-di-n-propyla...	9.069	70	27855	10.527	ng	97
20) 3+4-Methylphenols	9.016	107	35937	9.945	ng	99
22) Acetophenone	9.093	105	52184	10.322	ng	96
24) Nitrobenzene	9.481	77	27749	10.153	ng	99
25) Isophorone	10.003	82	70666	9.866	ng	96
26) 2-Nitrophenol	10.197	139	7321	9.931	ng	97
27) 2,4-Dimethylphenol	10.238	122	24040	7.931	ng	97
28) bis(2-Chloroethoxy)met...	10.485	93	39004	10.100	ng	98
29) 2,4-Dichlorophenol	10.732	162	24888	8.844	ng	99
30) 1,2,4-Trichlorobenzene	10.955	180	35754	10.142	ng	97
31) Naphthalene	11.149	128	104140	10.135	ng	99
32) Benzoic acid	10.291	122	8447m	11.421	ng	
33) 4-Chloroaniline	11.243	127	43688	10.172	ng	97
34) Hexachlorobutadiene	11.431	225	23338	9.927	ng	94
35) Caprolactam	11.983	113	8696	8.599	ng	# 83
36) 4-Chloro-3-methylphenol	12.336	107	31273	9.427	ng	99
37) 2-Methylnaphthalene	12.736	142	78890	10.219	ng	98
38) 1-Methylnaphthalene	12.953	142	78178	10.352	ng	98
40) 1,2,4,5-Tetrachloroben...	13.094	216	41498	9.881	ng	99
41) Hexachlorocyclopentadiene	13.076	237	15129	8.063	ng	100
43) 2,4,6-Trichlorophenol	13.323	196	20741	10.361	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.394	196	25139	10.253	ng	97
46) 1,1'-Biphenyl	13.729	154	104684	10.177	ng	100
47) 2-Chloronaphthalene	13.770	162	80179	10.092	ng	98
48) 2-Nitroaniline	13.958	65	13138	10.993	ng	93
49) Acenaphthylene	14.616	152	134714	10.248	ng	98
50) Dimethylphthalate	14.328	163	105750	9.898	ng	100
51) 2,6-Dinitrotoluene	14.451	165	12764	10.417	ng	94
52) Acenaphthene	14.951	154	82430	10.044	ng	99
53) 3-Nitroaniline	14.774	138	15812	10.221	ng	# 98
54) 2,4-Dinitrophenol	14.974	184	4769	8.855	ng	# 21
55) Dibenzofuran	15.286	168	136228	10.383	ng	97
56) 4-Nitrophenol	15.056	139	10773	11.302	ng	96
57) 2,4-Dinitrotoluene	15.227	165	15749	10.572	ng	97
58) Fluorene	15.932	166	107975	10.344	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.497	232	22617	10.198	ng	# 98
60) Diethylphthalate	15.685	149	108634	9.995	ng	99
61) 4-Chlorophenyl-phenyle...	15.914	204	56776	10.226	ng	99
62) 4-Nitroaniline	15.932	138	17822	9.902	ng	91
63) Azobenzene	16.208	77	109709	10.414	ng	99
65) 4,6-Dinitro-2-methylph...	15.985	198	6923	10.147	ng	99
66) n-Nitrosodiphenylamine	16.126	169	93299	10.065	ng	99
67) 4-Bromophenyl-phenylether	16.813	248	34626	10.258	ng	96
68) Hexachlorobenzene	16.931	284	40453	10.081	ng	98
69) Atrazine	17.060	200	32198	9.598	ng	98
70) Pentachlorophenol	17.266	266	15808	11.371	ng	93
71) Phenanthrene	17.671	178	186136	10.467	ng	98
72) Anthracene	17.759	178	181242	10.400	ng	98
73) Carbazole	18.023	167	164137	10.422	ng	98
74) Di-n-butylphthalate	18.570	149	168654	9.340	ng	99
75) Fluoranthene	19.663	202	220779	10.411	ng	99
77) Benzidine	19.833	184	82151	9.351	ng	96
78) Pyrene	20.021	202	229073	10.423	ng	98
80) Butylbenzylphthalate	20.902	149	58529	9.866	ng	98
81) Benzo(a)anthracene	21.907	228	225007	10.116	ng	100
82) 3,3'-Dichlorobenzidine	21.807	252	70161	9.665	ng	96
83) Chrysene	21.978	228	214698	10.151	ng	97
84) Bis(2-ethylhexyl)phtha...	21.801	149	94345	9.686	ng	99
85) Di-n-octyl phthalate	23.088	149	143530	7.488	ng	# 95
87) Indeno(1,2,3-cd)pyrene	29.269	276	254413	9.826	ng	98
88) Benzo(b)fluoranthene	24.251	252	216074	10.032	ng	99
89) Benzo(k)fluoranthene	24.322	252	224030	10.174	ng	99
90) Benzo(a)pyrene	25.186	252	185523	9.999	ng	100
91) Dibenzo(a,h)anthracene	29.357	278	214076	9.989	ng	96
92) Benzo(g,h,i)perylene	30.509	276	207723	9.915	ng	100

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

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