

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG102622\
 Data File : BG055299.D
 Acq On : 26 Oct 2022 14:49
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

Reviewed By : Christian
 Giraldo

Quant Time : Oct 27 10:35:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG102422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 24 16:54:05 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.276	152	46142	20.000	ng	0.00
21) Naphthalene-d8	11.108	136	199936	20.000	ng	0.00
39) Acenaphthene-d10	14.886	164	136071	20.000	ng	0.00
64) Phenanthrene-d10	17.618	188	284710	20.000	ng	0.00
76) Chrysene-d12	21.901	240	244294	20.000	ng	0.00
86) Perylene-d12	25.297	264	261404	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.797	112	201658	76.524	ng	0.00
7) Phenol-d6	7.418	99	292201	76.536	ng	0.00
23) Nitrobenzene-d5	9.451	82	313048	79.952	ng	0.00
42) 2,4,6-Tribromophenol	16.367	330	125535	84.866	ng	0.00
45) 2-Fluorobiphenyl	13.517	172	698373	84.005	ng	0.00
79) Terphenyl-d14	20.192	244	966949	82.238	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.605	88	43556	34.231	ng	97
3) Pyridine	4.028	79	139958	36.526	ng	96
4) n-Nitrosodimethylamine	3.934	42	59474	34.936	ng	# 97
6) Aniline	7.589	93	188882	37.792	ng	99
8) 2-Chlorophenol	7.836	128	120474	39.323	ng	99
9) Benzaldehyde	7.401	77	88119	33.805	ng	97
10) Phenol	7.448	94	162569	38.015	ng	99
11) bis(2-Chloroethyl)ether	7.677	93	124874	38.683	ng	99
12) 1,3-Dichlorobenzene	8.165	146	133637	39.978	ng	98
13) 1,4-Dichlorobenzene	8.311	146	136573	40.000	ng	100
14) 1,2-Dichlorobenzene	8.635	146	128973	39.233	ng	98
15) Benzyl Alcohol	8.511	79	124157	38.861	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.793	45	206651	39.117	ng	99
17) 2-Methylphenol	8.711	107	115183	38.086	ng	98
18) Hexachloroethane	9.369	117	50433	41.481	ng	97
19) n-Nitroso-di-n-propyla...	9.081	70	122242	38.558	ng	99
20) 3+4-Methylphenols	9.040	107	163934	37.846	ng	98
22) Acetophenone	9.105	105	204735	40.054	ng	# 97
24) Nitrobenzene	9.492	77	163318	40.059	ng	97
25) Isophorone	10.015	82	311518	39.390	ng	97
26) 2-Nitrophenol	10.203	139	78059	43.722	ng	98
27) 2,4-Dimethylphenol	10.256	122	112290	40.287	ng	97
28) bis(2-Chloroethoxy)met...	10.491	93	166965	41.814	ng	99
29) 2,4-Dichlorophenol	10.744	162	129955	43.462	ng	98
30) 1,2,4-Trichlorobenzene	10.961	180	134107	43.945	ng	99
31) Naphthalene	11.155	128	434209	40.710	ng	99
32) Benzoic acid	10.403	122	80421	39.039	ng	99
33) 4-Chloroaniline	11.255	127	176527	38.713	ng	98
34) Hexachlorobutadiene	11.431	225	81385	46.676	ng	98
35) Caprolactam	12.037	113	48046	34.622	ng	96
36) 4-Chloro-3-methylphenol	12.360	107	150485	39.997	ng	98
37) 2-Methylnaphthalene	12.736	142	317439	41.279	ng	98
38) 1-Methylnaphthalene	12.953	142	305694	40.294	ng	99
40) 1,2,4,5-Tetrachloroben...	13.100	216	148581	45.273	ng	99
41) Hexachlorocyclopentadiene	13.076	237	73199	48.083	ng	98
43) 2,4,6-Trichlorophenol	13.329	196	108671	44.832	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.406	196	120886	45.089	ng	99
46) 1,1'-Biphenyl	13.729	154	396510	42.486	ng	99
47) 2-Chloronaphthalene	13.776	162	307033	41.720	ng	99
48) 2-Nitroaniline	13.970	65	113780	39.533	ng	97
49) Acenaphthylene	14.616	152	516786	40.679	ng	99
50) Dimethylphthalate	14.328	163	422235	39.522	ng	100
51) 2,6-Dinitrotoluene	14.457	165	89894	41.350	ng	93
52) Acenaphthene	14.951	154	335637m	41.447	ng	10/27/2022
53) 3-Nitroaniline	14.786	138	104499	38.245	ng	99
54) 2,4-Dinitrophenol	14.992	184	49702	41.130	ng	99
55) Dibenzofuran	15.280	168	485863	39.992	ng	100
56) 4-Nitrophenol	15.086	139	76162	38.971	ng	96
57) 2,4-Dinitrotoluene	15.239	165	125529	39.790	ng	92
58) Fluorene	15.926	166	396788	39.252	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.503	232	104067	43.540	ng	95
60) Diethylphthalate	15.673	149	436167	38.380	ng	98
61) 4-Chlorophenyl-phenyle...	15.908	204	196411	42.281	ng	99
62) 4-Nitroaniline	15.944	138	105522	36.750	ng	99
63) Azobenzene	16.202	77	425568	38.439	ng	98
65) 4,6-Dinitro-2-methylph...	15.997	198	73143	46.732	ng	99
66) n-Nitrosodiphenylamine	16.120	169	347346	43.831	ng	100
67) 4-Bromophenyl-phenylether	16.802	248	129179	46.701	ng	97
68) Hexachlorobenzene	16.925	284	140589	44.952	ng	97
69) Atrazine	17.054	200	113437	42.097	ng	100
70) Pentachlorophenol	17.272	266	73321	42.722	ng	99
71) Phenanthrene	17.665	178	621711	40.946	ng	99
72) Anthracene	17.753	178	614729	41.413	ng	99
73) Carbazole	18.018	167	562621	39.407	ng	99
74) Di-n-butylphthalate	18.547	149	717145	40.183	ng	100
75) Fluoranthene	19.651	202	696194	39.349	ng	98
77) Benzidine	19.816	184	222466	37.610	ng	99
78) Pyrene	20.010	202	701773	41.910	ng	99
80) Butylbenzylphthalate	20.867	149	313779	39.655	ng	95
81) Benzo(a)anthracene	21.878	228	646772	39.426	ng	99
82) 3,3'-Dichlorobenzidine	21.778	252	216822	40.582	ng	100
83) Chrysene	21.948	228	616557	39.567	ng	99
84) Bis(2-ethylhexyl)phtha...	21.743	149	450292	39.836	ng	99
85) Di-n-octyl phthalate	23.012	149	771152	40.951	ng	99
87) Indeno(1,2,3-cd)pyrene	29.211	276	801557	41.517	ng	# 93
88) Benzo(b)fluoranthene	24.211	252	640500	40.734	ng	98
89) Benzo(k)fluoranthene	24.281	252	654197	41.203	ng	98
90) Benzo(a)pyrene	25.139	252	551860	40.927	ng	98
91) Dibenzo(a,h)anthracene	29.281	278	663605	41.884	ng	98
92) Benzo(g,h,i)perylene	30.450	276	647163	41.181	ng	98

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

Manual Integrations

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