

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110922\
 Data File : BG055522.D
 Acq On : 9 Nov 2022 13:50
 Operator : CG/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC050

Quant Time: Nov 10 03:01:11 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 02:56:38 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.274	152	50294	20.000	ng	0.00
21) Naphthalene-d8	11.100	136	208575	20.000	ng	0.00
39) Acenaphthene-d10	14.890	164	151167	20.000	ng	0.00
64) Phenanthrene-d10	17.628	188	343777	20.000	ng	0.00
76) Chrysene-d12	21.934	240	321379	20.000	ng	0.00
86) Perylene-d12	25.354	264	348722	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.794	112	262378	99.367	ng	0.00
7) Phenol-d6	7.398	99	365315	100.158	ng	0.00
23) Nitrobenzene-d5	9.443	82	344801	95.994	ng	0.00
42) 2,4,6-Tribromophenol	16.370	330	158905	95.692	ng	0.00
45) 2-Fluorobiphenyl	13.521	172	933850	93.183	ng	0.00
79) Terphenyl-d14	20.219	244	1498020	93.623	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.638	88	62165	46.351	ng	98
3) Pyridine	4.049	79	185573	50.551	ng	99
4) n-Nitrosodimethylamine	3.955	42	98575	48.565	ng	98
6) Aniline	7.586	93	234483	48.400	ng	100
8) 2-Chlorophenol	7.827	128	147501	49.897	ng	100
9) Benzaldehyde	7.398	77	132330	45.985	ng	98
10) Phenol	7.428	94	204626	48.734	ng	100
11) bis(2-Chloroethyl)ether	7.680	93	158958	47.021	ng	97
12) 1,3-Dichlorobenzene	8.162	146	175039	47.402	ng	98
13) 1,4-Dichlorobenzene	8.309	146	174997	46.827	ng	99
14) 1,2-Dichlorobenzene	8.632	146	167022	47.062	ng	98
15) Benzyl Alcohol	8.503	79	161127	50.658	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.797	45	305732	47.066	ng	99
17) 2-Methylphenol	8.697	107	144993	49.070	ng	98
18) Hexachloroethane	9.373	117	59575	48.018	ng	92
19) n-Nitroso-di-n-propyla...	9.079	70	146747	48.683	ng	98
20) 3+4-Methylphenols	9.026	107	206634	50.198	ng	96
22) Acetophenone	9.096	105	266045	47.950	ng	# 99
24) Nitrobenzene	9.484	77	189764	48.456	ng	98
25) Isophorone	10.007	82	387044	49.240	ng	99
26) 2-Nitrophenol	10.201	139	60445	49.755	ng	93
27) 2,4-Dimethylphenol	10.242	122	146426	48.155	ng	100
28) bis(2-Chloroethoxy)met...	10.489	93	208121	49.107	ng	99
29) 2,4-Dichlorophenol	10.736	162	163728	53.013	ng	98
30) 1,2,4-Trichlorobenzene	10.959	180	187694	48.516	ng	100
31) Naphthalene	11.153	128	544256	48.264	ng	99
32) Benzoic acid	10.365	122	84933	49.220	ng	98
33) 4-Chloroaniline	11.247	127	232859	49.402	ng	98
34) Hexachlorobutadiene	11.435	225	125624	48.688	ng	97
35) Caprolactam	12.022	113	56383	50.803	ng	95
36) 4-Chloro-3-methylphenol	12.340	107	188735	51.843	ng	98
37) 2-Methylnaphthalene	12.739	142	412133	48.643	ng	100
38) 1-Methylnaphthalene	12.957	142	399227	48.172	ng	98
40) 1,2,4,5-Tetrachloroben...	13.098	216	215370	47.695	ng	98
41) Hexachlorocyclopentadiene	13.080	237	101902	50.515	ng	99
43) 2,4,6-Trichlorophenol	13.327	196	144808	48.345	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.397	196	162921	48.686	ng	98
46) 1,1'-Biphenyl	13.732	154	528930	47.827	ng	99
47) 2-Chloronaphthalene	13.779	162	411779	48.207	ng	99
48) 2-Nitroaniline	13.967	65	112865	46.897	ng	95
49) Acenaphthylene	14.619	152	679018	48.042	ng	99
50) Dimethylphthalate	14.337	163	572158	49.810	ng	99
51) 2,6-Dinitrotoluene	14.455	165	101236	46.945	ng	100
52) Acenaphthene	14.954	154	427309	48.428	ng	99
53) 3-Nitroaniline	14.784	138	116335	47.863	ng	99
54) 2,4-Dinitrophenol	14.978	184	36011	48.834	ng	# 58
55) Dibenzofuran	15.289	168	671757	47.620	ng	99
56) 4-Nitrophenol	15.066	139	89689	47.900	ng	97
57) 2,4-Dinitrotoluene	15.236	165	135869	47.230	ng	98
58) Fluorene	15.935	166	535784	47.741	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.501	232	151449	48.465	ng	98
60) Diethylphthalate	15.689	149	578844	49.535	ng	99
61) 4-Chlorophenyl-phenyle...	15.918	204	282102	47.258	ng	99
62) 4-Nitroaniline	15.941	138	120018	47.822	ng	99
63) Azobenzene	16.212	77	540523	47.721	ng	99
65) 4,6-Dinitro-2-methylph...	15.994	198	55159	49.327	ng	97
66) n-Nitrosodiphenylamine	16.129	169	473815	49.199	ng	99
67) 4-Bromophenyl-phenylether	16.811	248	171186	48.812	ng	98
68) Hexachlorobenzene	16.934	284	205522	49.293	ng	96
69) Atrazine	17.069	200	178158	51.117	ng	98
70) Pentachlorophenol	17.269	266	121496	48.455	ng	95
71) Phenanthrene	17.675	178	893546	48.363	ng	99
72) Anthracene	17.763	178	872270	48.175	ng	99
73) Carbazole	18.027	167	784817	47.961	ng	99
74) Di-n-butylphthalate	18.574	149	962887	51.322	ng	99
75) Fluoranthene	19.666	202	1035345	46.992	ng	99
77) Benzidine	19.837	184	443385	50.995	ng	100
78) Pyrene	20.031	202	1056383	48.568	ng	99
80) Butylbenzylphthalate	20.906	149	408614	49.232	ng	95
81) Benzo(a)anthracene	21.911	228	1062536	48.268	ng	99
82) 3,3'-Dichlorobenzidine	21.811	252	363335	50.573	ng	99
83) Chrysene	21.981	228	1008700	48.189	ng	100
84) Bis(2-ethylhexyl)phtha...	21.805	149	593892	49.137	ng	99
85) Di-n-octyl phthalate	23.092	149	995379	52.472	ng	99
87) Indeno(1,2,3-cd)pyrene	29.302	276	1298006	48.871	ng	99
88) Benzo(b)fluoranthene	24.261	252	1070263	48.439	ng	100
89) Benzo(k)fluoranthene	24.337	252	1087715	48.153	ng	99
90) Benzo(a)pyrene	25.195	252	927720	48.743	ng	99
91) Dibenzo(a,h)anthracene	29.373	278	1065777	48.476	ng	99
92) Benzo(g,h,i)perylene	30.536	276	1051494	48.925	ng	100

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

