

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
 Data File : BG057716.D
 Acq On : 14 Jun 2023 20:48
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Jun 15 01:24:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 14 03:52:08 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.935	152	33308	20.000	ng	0.00
21) Naphthalene-d8	10.738	136	131621	20.000	ng	0.00
39) Acenaphthene-d10	14.569	164	89261	20.000	ng	0.00
64) Phenanthrene-d10	17.313	188	211167	20.000	ng	0.00
76) Chrysene-d12	21.572	240	182383	20.000	ng	0.00
86) Perylene-d12	24.733	264	236849	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.509	112	170329	83.267	ng	0.00
7) Phenol-d6	7.095	99	242198	77.072	ng	0.00
23) Nitrobenzene-d5	9.099	82	196462	91.356	ng	0.00
42) 2,4,6-Tribromophenol	16.055	330	97543	97.950	ng	0.00
45) 2-Fluorobiphenyl	13.194	172	460999	76.984	ng	0.00
79) Terphenyl-d14	19.933	244	747158	75.741	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.423	88	38975	37.425	ng	91
3) Pyridine	3.817	79	112277	38.085	ng	94
4) n-Nitrosodimethylamine	3.729	42	50714	35.270	ng	87
6) Aniline	7.266	93	150575	36.520	ng	96
8) 2-Chlorophenol	7.501	128	80257	40.921	ng	99
9) Benzaldehyde	7.078	77	70631	36.464	ng	98
10) Phenol	7.119	94	117585	38.292	ng	99
11) bis(2-Chloroethyl)ether	7.360	93	88170	36.634	ng	98
12) 1,3-Dichlorobenzene	7.830	146	86218	38.003	ng	97
13) 1,4-Dichlorobenzene	7.971	146	87191	38.762	ng	98
14) 1,2-Dichlorobenzene	8.294	146	81856	37.453	ng	97
15) Benzyl Alcohol	8.170	79	87596	36.396	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.464	45	172994	33.010	ng	98
17) 2-Methylphenol	8.370	107	83862	37.113	ng	95
18) Hexachloroethane	9.022	117	30864	44.400	ng	96
19) n-Nitroso-di-n-propyla...	8.740	70	76352	33.602	ng	96
20) 3+4-Methylphenols	8.699	107	115757	37.467	ng	97
22) Acetophenone	8.758	105	140803	36.823	ng	97
24) Nitrobenzene	9.140	77	101866	43.466	ng	96
25) Isophorone	9.663	82	215487	35.322	ng	99
26) 2-Nitrophenol	9.851	139	36052	57.004	ng	# 85
27) 2,4-Dimethylphenol	9.909	122	83532	41.005	ng	91
28) bis(2-Chloroethoxy)met...	10.145	93	117503	37.051	ng	98
29) 2,4-Dichlorophenol	10.380	162	78566	45.465	ng	98
30) 1,2,4-Trichlorobenzene	10.603	180	89375	39.793	ng	96
31) Naphthalene	10.791	128	272697	38.550	ng	98
32) Benzoic acid	10.021	122	59936	75.290	ng	97
33) 4-Chloroaniline	10.897	127	122132	37.083	ng	96
34) Hexachlorobutadiene	11.073	225	54540	38.615	ng	99
35) Caprolactam	11.666	113	30688	43.049	ng	99
36) 4-Chloro-3-methylphenol	12.013	107	98908	40.401	ng	98
37) 2-Methylnaphthalene	12.401	142	193448	36.964	ng	99
38) 1-Methylnaphthalene	12.618	142	182889	37.251	ng	99
40) 1,2,4,5-Tetrachloroben...	12.765	216	98593	39.019	ng	99
41) Hexachlorocyclopentadiene	12.747	237	53144	55.144	ng	97
43) 2,4,6-Trichlorophenol	13.000	196	70024	43.179	ng	100

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44) 2,4,5-Trichlorophenol	13.070	196	83668	43.592	ng	95
46) 1,1'-Biphenyl	13.405	154	237279	38.474	ng	99
47) 2-Chloronaphthalene	13.447	162	184115	39.499	ng	98
48) 2-Nitroaniline	13.646	65	65397	50.957	ng	95
49) Acenaphthylene	14.293	152	314605	38.918	ng	98
50) Dimethylphthalate	14.028	163	253050	38.664	ng	99
51) 2,6-Dinitrotoluene	14.140	165	47788	49.495	ng	92
52) Acenaphthene	14.633	154	193368	39.906	ng	100
53) 3-Nitroaniline	14.469	138	55860	49.376	ng	# 93
54) 2,4-Dinitrophenol	14.675	184	18841	57.628	ng	98
55) Dibenzofuran	14.968	168	305959	37.617	ng	99
56) 4-Nitrophenol	14.769	139	46576	51.934	ng	91
57) 2,4-Dinitrotoluene	14.927	165	66291	51.884	ng	# 98
58) Fluorene	15.615	166	242828	37.478	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.192	232	68716	43.007	ng	97
60) Diethylphthalate	15.391	149	265310	39.271	ng	97
61) 4-Chlorophenyl-phenyle...	15.609	204	127744	36.802	ng	97
62) 4-Nitroaniline	15.638	138	57274	46.296	ng	94
63) Azobenzene	15.902	77	270213	37.198	ng	99
65) 4,6-Dinitro-2-methylph...	15.685	198	32161	59.775	ng	89
66) n-Nitrosodiphenylamine	15.826	169	216962	39.304	ng	97
67) 4-Bromophenyl-phenylether	16.508	248	84366	39.237	ng	97
68) Hexachlorobenzene	16.619	284	86421	38.459	ng	100
69) Atrazine	16.772	200	80198	43.667	ng	99
70) Pentachlorophenol	16.960	266	70054	51.928	ng	96
71) Phenanthrene	17.360	178	396908	38.518	ng	100
72) Anthracene	17.448	178	415316	39.786	ng	100
73) Carbazole	17.712	167	390561	40.399	ng	100
74) Di-n-butylphthalate	18.282	149	490849	45.807	ng	99
75) Fluoranthene	19.369	202	493663	39.769	ng	100
77) Benzidine	19.545	184	184778	42.499	ng	98
78) Pyrene	19.727	202	508868	39.214	ng	98
80) Butylbenzylphthalate	20.620	149	220616	46.822	ng	99
81) Benzo(a)anthracene	21.555	228	493080	39.148	ng	100
82) 3,3'-Dichlorobenzidine	21.472	252	172707	42.097	ng	99
83) Chrysene	21.619	228	469776	39.191	ng	99
84) Bis(2-ethylhexyl)phtha...	21.472	149	314014	49.118	ng	99
85) Di-n-octyl phthalate	22.671	149	570160	51.721	ng	97
87) Indeno(1,2,3-cd)pyrene	28.341	276	593919	38.755	ng	99
88) Benzo(b)fluoranthene	23.729	252	484601	37.715	ng	98
89) Benzo(k)fluoranthene	23.799	252	493956	37.067	ng	97
90) Benzo(a)pyrene	24.586	252	483312	38.257	ng	# 98
91) Dibenzo(a,h)anthracene	28.405	278	490142	38.703	ng	98
92) Benzo(g,h,i)perylene	29.463	276	491826	38.910	ng	98

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

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