

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082223\
 Data File : BG058801.D
 Acq On : 22 Aug 2023 9:01
 Operator : MA/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Aug 22 17:16:59 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG081823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 21 01:40:57 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.809	152	30662	20.000	ng	0.00
21) Naphthalene-d8	10.611	136	125418	20.000	ng	0.00
39) Acenaphthene-d10	14.460	164	89825	20.000	ng	0.00
64) Phenanthrene-d10	17.203	188	208467	20.000	ng	0.00
76) Chrysene-d12	21.451	240	172227	20.000	ng	0.00
86) Perylene-d12	24.524	264	216837	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.382	112	152778	80.184	ng	0.00
7) Phenol-d6	6.980	99	222486	83.227	ng	0.00
23) Nitrobenzene-d5	8.978	82	206055	78.989	ng	0.00
42) 2,4,6-Tribromophenol	15.952	330	111143	78.912	ng	0.00
45) 2-Fluorobiphenyl	13.079	172	479374	77.621	ng	0.00
79) Terphenyl-d14	19.824	244	764861	79.742	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.267	88	33980	37.933	ng	94
3) Pyridine	3.666	79	93945	39.686	ng	97
4) n-Nitrosodimethylamine	3.584	42	48665	37.815	ng	90
6) Aniline	7.133	93	132986	40.696	ng	97
8) 2-Chlorophenol	7.374	128	78923	40.967	ng	96
9) Benzaldehyde	6.951	77	60425	40.297	ng	99
10) Phenol	7.010	94	107391	41.661	ng	97
11) bis(2-Chloroethyl)ether	7.233	93	81502	40.520	ng	98
12) 1,3-Dichlorobenzene	7.697	146	82429	40.281	ng	99
13) 1,4-Dichlorobenzene	7.844	146	84826	40.992	ng	96
14) 1,2-Dichlorobenzene	8.161	146	78846	39.422	ng	98
15) Benzyl Alcohol	8.049	79	81067	42.415	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.337	45	175784	41.008	ng	99
17) 2-Methylphenol	8.255	107	78689	41.562	ng	98
18) Hexachloroethane	8.884	117	31211	41.397	ng	95
19) n-Nitroso-di-n-propyla...	8.614	70	71930	41.626	ng	98
20) 3+4-Methylphenols	8.590	107	108302	42.548	ng	96
22) Acetophenone	8.637	105	137029	40.362	ng	# 96
24) Nitrobenzene	9.019	77	105623	40.710	ng	97
25) Isophorone	9.536	82	207142	39.929	ng	99
26) 2-Nitrophenol	9.724	139	46196	42.572	ng	97
27) 2,4-Dimethylphenol	9.789	122	78182	42.512	ng	97
28) bis(2-Chloroethoxy)met...	10.018	93	113037	41.135	ng	99
29) 2,4-Dichlorophenol	10.265	162	80404	42.547	ng	98
30) 1,2,4-Trichlorobenzene	10.476	180	92475	40.524	ng	97
31) Naphthalene	10.658	128	265114	40.259	ng	99
32) Benzoic acid	9.959	122	41213	37.932	ng	97
33) 4-Chloroaniline	10.776	127	118104	42.525	ng	99
34) Hexachlorobutadiene	10.940	225	63228	41.053	ng	99
35) Caprolactam	11.563	113	26454	39.066	ng	94
36) 4-Chloro-3-methylphenol	11.910	107	98334	41.843	ng	99
37) 2-Methylnaphthalene	12.274	142	195950	41.430	ng	97
38) 1-Methylnaphthalene	12.497	142	182109	40.562	ng	99
40) 1,2,4,5-Tetrachloroben...	12.650	216	111822	40.381	ng	98
41) Hexachlorocyclopentadiene	12.621	237	36870	38.924	ng	98
43) 2,4,6-Trichlorophenol	12.891	196	74452	42.253	ng	99

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44) 2,4,5-Trichlorophenol	12.973	196	87236	41.343	ng	95
46) 1,1'-Biphenyl	13.290	154	246773	39.880	ng	97
47) 2-Chloronaphthalene	13.332	162	192148	39.847	ng	99
48) 2-Nitroaniline	13.543	65	74471	40.663	ng	96
49) Acenaphthylene	14.183	152	315736	39.482	ng	98
50) Dimethylphthalate	13.919	163	266195	38.274	ng	99
51) 2,6-Dinitrotoluene	14.042	165	58964	40.355	ng	94
52) Acenaphthene	14.524	154	184761	39.537	ng	98
53) 3-Nitroaniline	14.377	138	57886	41.076	ng	99
54) 2,4-Dinitrophenol	14.595	184	22966	32.587	ng	93
55) Dibenzofuran	14.859	168	317447	39.380	ng	98
56) 4-Nitrophenol	14.695	139	33841	35.802	ng	97
57) 2,4-Dinitrotoluene	14.836	165	82559	40.510	ng	# 96
58) Fluorene	15.505	166	250116	38.987	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.088	232	71672	41.625	ng	95
60) Diethylphthalate	15.276	149	278071	38.778	ng	98
61) 4-Chlorophenyl-phenyle...	15.500	204	141496	39.792	ng	98
62) 4-Nitroaniline	15.541	138	60158	42.064	ng	95
63) Azobenzene	15.793	77	257433	39.097	ng	99
65) 4,6-Dinitro-2-methylph...	15.599	198	46948	39.842	ng	96
66) n-Nitrosodiphenylamine	15.717	169	222685	41.445	ng	97
67) 4-Bromophenyl-phenylether	16.393	248	95163	41.599	ng	99
68) Hexachlorobenzene	16.510	284	98849	40.149	ng	98
69) Atrazine	16.669	200	67489	35.493	ng	97
70) Pentachlorophenol	16.863	266	50876	39.407	ng	96
71) Phenanthrene	17.250	178	399368	39.905	ng	98
72) Anthracene	17.339	178	409335	40.781	ng	99
73) Carbazole	17.615	167	366496	39.873	ng	99
74) Di-n-butylphthalate	18.161	149	458700	39.560	ng	99
75) Fluoranthene	19.266	202	486489	39.199	ng	99
77) Benzidine	19.454	184	124670	40.390	ng	99
78) Pyrene	19.624	202	489588	40.798	ng	99
80) Butylbenzylphthalate	20.511	149	200716	41.241	ng	98
81) Benzo(a)anthracene	21.434	228	487898	40.697	ng	97
82) 3,3'-Dichlorobenzidine	21.352	252	169599	40.976	ng	99
83) Chrysene	21.493	228	466531	40.373	ng	99
84) Bis(2-ethylhexyl)phtha...	21.328	149	293356	41.097	ng	99
85) Di-n-octyl phthalate	22.480	149	507601	41.344	ng	100
87) Indeno(1,2,3-cd)pyrene	28.026	276	605776	41.292	ng	# 93
88) Benzo(b)fluoranthene	23.549	252	491761	40.874	ng	100
89) Benzo(k)fluoranthene	23.614	252	506635	40.832	ng	99
90) Benzo(a)pyrene	24.383	252	492598	41.582	ng	99
91) Dibenzo(a,h)anthracene	28.079	278	498065	41.316	ng	99
92) Benzo(g,h,i)perylene	29.119	276	493828	40.972	ng	97

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

