

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111122\
 Data File : BG055559.D
 Acq On : 11 Nov 2022 9:22
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Nov 11 16:29:37 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 03:39:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.268	152	53098	20.000	ng	0.00
21) Naphthalene-d8	11.100	136	224929	20.000	ng	0.00
39) Acenaphthene-d10	14.889	164	170746	20.000	ng	0.00
64) Phenanthrene-d10	17.633	188	419767	20.000	ng	0.00
76) Chrysene-d12	21.934	240	396288	20.000	ng	0.00
86) Perylene-d12	25.354	264	437415	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.794	112	233829	83.879	ng	0.00
7) Phenol-d6	7.404	99	322712	83.805	ng	0.00
23) Nitrobenzene-d5	9.443	82	339514	102.327	ng	0.00
42) 2,4,6-Tribromophenol	16.370	330	193064	115.886	ng	0.00
45) 2-Fluorobiphenyl	13.520	172	872100	77.043	ng	0.00
79) Terphenyl-d14	20.218	244	1519209	77.000	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.638	88	50862	35.921	ng	96
3) Pyridine	4.049	79	162681	41.321	ng	98
4) n-Nitrosodimethylamine	3.961	42	84550	39.456	ng	95
6) Aniline	7.586	93	203502	39.787	ng	97
8) 2-Chlorophenol	7.827	128	133830	42.882	ng	99
9) Benzaldehyde	7.398	77	113594	37.389	ng	98
10) Phenol	7.428	94	182570	41.185	ng	100
11) bis(2-Chloroethyl)ether	7.674	93	138560	38.823	ng	94
12) 1,3-Dichlorobenzene	8.162	146	154506	39.632	ng	99
13) 1,4-Dichlorobenzene	8.309	146	156414	39.644	ng	98
14) 1,2-Dichlorobenzene	8.632	146	147759	39.435	ng	98
15) Benzyl Alcohol	8.503	79	142480	42.430	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.791	45	258450	37.686	ng	99
17) 2-Methylphenol	8.697	107	130781	41.923	ng	98
18) Hexachloroethane	9.372	117	54577	41.667	ng	97
19) n-Nitroso-di-n-propyla...	9.079	70	127357	40.019	ng	98
20) 3+4-Methylphenols	9.026	107	184073	42.356	ng	97
22) Acetophenone	9.096	105	233558	39.034	ng	# 99
24) Nitrobenzene	9.484	77	177349	46.461	ng	99
25) Isophorone	10.007	82	335866	39.622	ng	99
26) 2-Nitrophenol	10.201	139	73956	58.171	ng	# 91
27) 2,4-Dimethylphenol	10.242	122	130645	43.107	ng	99
28) bis(2-Chloroethoxy)met...	10.489	93	179149	39.198	ng	99
29) 2,4-Dichlorophenol	10.730	162	151668	45.537	ng	96
30) 1,2,4-Trichlorobenzene	10.959	180	171934	41.211	ng	97
31) Naphthalene	11.153	128	485801	39.948	ng	99
32) Benzoic acid	10.365	122	109177	55.476	ng	95
33) 4-Chloroaniline	11.247	127	207623	40.845	ng	98
34) Hexachlorobutadiene	11.435	225	115545	41.526	ng	98
35) Caprolactam	12.022	113	55884	49.219	ng	94
36) 4-Chloro-3-methylphenol	12.345	107	174545	44.459	ng	98
37) 2-Methylnaphthalene	12.739	142	369589	40.450	ng	99
38) 1-Methylnaphthalene	12.956	142	358547	40.118	ng	97
40) 1,2,4,5-Tetrachloroben...	13.097	216	205366	40.265	ng	99
41) Hexachlorocyclopentadiene	13.080	237	107371	49.174	ng	97
43) 2,4,6-Trichlorophenol	13.327	196	146082	48.711	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111122\
 Data File : BG055559.D
 Acq On : 11 Nov 2022 9:22
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Nov 11 16:29:37 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 03:39:46 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.397	196	161104	46.925	ng	100
46) 1,1'-Biphenyl	13.732	154	481549	38.550	ng	98
47) 2-Chloronaphthalene	13.779	162	374186	38.783	ng	99
48) 2-Nitroaniline	13.967	65	126097	50.522	ng	97
49) Acenaphthylene	14.619	152	629593	39.437	ng	99
50) Dimethylphthalate	14.337	163	547246	42.179	ng	99
51) 2,6-Dinitrotoluene	14.455	165	108209	47.482	ng	96
52) Acenaphthene	14.954	154	408842	41.022	ng	99
53) 3-Nitroaniline	14.784	138	124670	45.604	ng	# 95
54) 2,4-Dinitrophenol	14.983	184	53157	65.894	ng	# 85
55) Dibenzofuran	15.289	168	630732	39.585	ng	99
56) 4-Nitrophenol	15.066	139	98632	51.339	ng	98
57) 2,4-Dinitrotoluene	15.236	165	152137	50.311	ng	97
58) Fluorene	15.935	166	506749	39.976	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.506	232	159448	50.107	ng	97
60) Diethylphthalate	15.694	149	553470	41.933	ng	100
61) 4-Chlorophenyl-phenylether	15.918	204	277693	41.185	ng	99
62) 4-Nitroaniline	15.947	138	129874	50.489	ng	99
63) Azobenzene	16.211	77	493044	38.538	ng	97
65) 4,6-Dinitro-2-methylph...	15.994	198	78472	55.163	ng	98
66) n-Nitrosodiphenylamine	16.129	169	455607	38.744	ng	99
67) 4-Bromophenyl-phenylether	16.817	248	187449	43.773	ng	95
68) Hexachlorobenzene	16.934	284	205052	40.277	ng	97
69) Atrazine	17.069	200	180007	42.298	ng	97
70) Pentachlorophenol	17.269	266	138368	49.609	ng	98
71) Phenanthrene	17.674	178	877647	38.903	ng	99
72) Anthracene	17.763	178	855155	38.680	ng	100
73) Carbazole	18.027	167	771956	38.635	ng	99
74) Di-n-butylphthalate	18.573	149	937292	40.914	ng	99
75) Fluoranthene	19.666	202	1039222	38.629	ng	99
77) Benzidine	19.837	184	408899	38.139	ng	99
78) Pyrene	20.030	202	1053086	39.264	ng	99
80) Butylbenzylphthalate	20.900	149	413487	41.005	ng	99
81) Benzo(a)anthracene	21.911	228	1070768	39.447	ng	99
82) 3,3'-Dichlorobenzidine	21.817	252	377018	42.558	ng	99
83) Chrysene	21.981	228	1020801	39.549	ng	99
84) Bis(2-ethylhexyl)phtha...	21.799	149	587028	43.368	ng	99
85) Di-n-octyl phthalate	23.086	149	1009931	43.175	ng	99
87) Indeno(1,2,3-cd)pyrene	29.296	276	1311809	39.376	ng	# 96
88) Benzo(b)fluoranthene	24.261	252	1083838	39.107	ng	98
89) Benzo(k)fluoranthene	24.337	252	1093253	38.584	ng	99
90) Benzo(a)pyrene	25.195	252	932459	39.058	ng	99
91) Dibenzo(a,h)anthracene	29.367	278	1077047	39.055	ng	98
92) Benzo(g,h,i)perylene	30.530	276	1069209	39.662	ng	98

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

Quant Time: Nov 11 16:29:37 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 03:39:46 2022
Response via : Initial Calibration

