

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081722\
 Data File : BG054643.D
 Acq On : 17 Aug 2022 16:50
 Operator : CG/JU
 Sample : SSTD02037
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTD020437

Manual Integrations

APPROVED

Reviewed By :Jagrut Upadhyay 08/18/2022

Supervised By :Sohil Jodhani 08/18/2022

Quant Time: Aug 18 02:23:32 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081722.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Aug 18 02:20:33 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.165	152	44527	20.000	ng/u1	0.00
20) Naphthalene-d8	10.991	136	187878	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.798	164	124027	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.542	188	277314	20.000	ng/u1	0.00
79) Chrysene-d12	21.843	240	259287	20.000	ng/u1	0.00
88) Perylene-d12	25.221	264	267616	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.517	96	9000	9.617	ng/uL	0.00
4) Pyridine-d5	3.940	84	63290	24.865	ng/u1	0.00
7) Phenol-d5	7.301	99	72852	22.838	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.483	67	47585	24.729	ng/u1	0.00
11) 2-Chlorophenol-d4	7.689	132	52678	20.708	ng/u1	0.00
15) 4-Methylphenol-d8	8.858	113	55286	20.326	ng/u1	0.00
21) Nitrobenzene-d5	9.334	128	22363	16.036	ng/u1	0.00
24) 2-Nitrophenol-d4	10.062	143	20144	13.595	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.597	165	54547	16.639	ng/u1	0.00
31) 4-Chloroaniline-d4	11.120	131	77751	18.774	ng/u1	0.00
46) Dimethylphthalate-d6	14.193	166	164620	16.872	ng/u1	0.00
49) Acenaphthylene-d8	14.492	160	196262	18.498	ng/u1	0.00
54) 4-Nitrophenol-d4	14.963	143	24488	18.329	ng/u1	0.00
60) Fluorene-d10	15.785	176	150040	17.120	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.885	200	15242	10.076	ng/u1	0.00
73) Anthracene-d10	17.642	188	233001	18.604	ng/u1	0.00
81) Pyrene-d10	19.921	212	266101	19.683	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.986	264	254365	18.830	ng/u1	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.552	88	9928	9.722	ng/uL	100
5) Pyridine	3.958	79	65394	24.773	ng/u1	100
6) Benzaldehyde	7.295	77	37553	23.118	ng/u1	100
8) Phenol	7.330	94	71953	21.617	ng/u1	100
10) Bis(2-Chloroethyl)ether	7.577	93	60801	24.115	ng/u1	100
12) 2-Chlorophenol	7.724	128	56705	21.833	ng/u1	100
13) 2-Methylphenol	8.593	108	55208	21.367	ng/u1	100
14) 2,2'-oxybis(1-Chloropr...	8.687	45	97870	26.133	ng/u1	100
16) Acetophenone	8.993	105	88848	21.122	ng/u1	100
17) N-Nitroso-di-n-propyla...	8.970	70	47883	22.330	ng/u1	100
18) 4-Methylphenol	8.923	108	58776	21.265	ng/u1	100
19) Hexachloroethane	9.263	117	22872	20.782	ng/u1	100
22) Nitrobenzene	9.381	77	59853	18.019	ng/u1	100
23) Isophorone	9.904	82	131487	20.417	ng/u1	100
25) 2-Nitrophenol	10.092	139	23978	15.433	ng/u1	100
26) 2,4-Dimethylphenol	10.139	107	64818	19.105	ng/u1	100
27) Bis(2-Chloroethoxy)met...	10.386	93	79301	21.873	ng/u1	100
29) 2,4-Dichlorophenol	10.626	162	51947	15.846	ng/u1	100
30) Naphthalene	11.044	128	187574	20.126	ng/u1	100
32) 4-Chloroaniline	11.143	127	79889	19.453	ng/u1	100
33) Hexachlorobutadiene	11.320	225	38623	12.062	ng/u1	100
34) Caprolactam	11.896	113	17196m	17.578	ng/u1	
35) 4-Chloro-3-methylphenol	12.242	107	57217	18.149	ng/u1	100
36) 2-Methylnaphthalene	12.636	142	128772	18.678	ng/u1	100

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081722\
 Data File : BG054643.D
 Acq On : 17 Aug 2022 16:50
 Operator : CG/JU
 Sample : SSTD02037
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTD020437

Quant Time: Aug 18 02:23:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 18 02:20:33 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 08/18/2022
 Supervised By :Sohil Jodhani 08/18/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
37) 1-Methylnaphthalene	12.853	142	127584	18.556	ng/u1	100
39) 1,2,4,5-Tetrachloroben...	12.994	216	72163	14.314	ng/u1	100
40) Hexachlorocyclopentadiene	12.977	237	43327	15.324	ng/u1	100
41) 2,4,6-Trichlorophenol	13.229	196	42473	16.233	ng/u1	100
42) 2,4,5-Trichlorophenol	13.300	196	47415	16.029	ng/u1	100
43) 1,1'-Biphenyl	13.635	154	170962	19.888	ng/u1	100
44) 2-Chloronaphthalene	13.676	162	128905	18.057	ng/u1	100
45) 2-Nitroaniline	13.870	65	30676	17.081	ng/u1	100
47) Dimethylphthalate	14.240	163	168613	17.980	ng/u1	100
48) 2,6-Dinitrotoluene	14.363	165	25713	13.633	ng/u1	100
50) Acenaphthylene	14.522	152	217049	19.924	ng/u1	100
51) 3-Nitroaniline	14.692	138	25389	16.582	ng/u1	100
52) Acenaphthene	14.857	153	144604	19.877	ng/u1	100
53) 2,4-Dinitrophenol	14.892	184	10684	12.228	ng/u1	100
55) 4-Nitrophenol	14.974	109	20720	15.821	ng/u1	100
56) Dibenzofuran	15.192	168	198525	17.835	ng/u1	100
57) 2,4-Dinitrotoluene	15.145	165	35554	12.995	ng/u1	100
58) 2,3,4,6-Tetrachlorophenol	15.409	232	38119	14.830	ng/u1	100
59) Diethylphthalate	15.603	149	167407	18.206	ng/u1	100
61) Fluorene	15.844	166	162468	17.921	ng/u1	100
62) 4-Chlorophenyl-phenyle...	15.832	204	82964	14.393	ng/u1	100
63) 4-Nitroaniline	15.850	138	24489m	16.634	ng/u1	
66) 4,6-Dinitro-2-methylph...	15.903	198	16723	11.516	ng/u1	100
67) N-Nitrosodiphenylamine	16.044	169	144301	21.284	ng/u1	100
68) 4-Bromophenyl-phenylether	16.725	248	54977	15.960	ng/u1	100
69) Hexachlorobenzene	16.837	284	63557	15.743	ng/u1	100
70) Atrazine	16.984	200	54779	16.450	ng/u1	100
71) Pentachlorophenol	17.183	266	32524	19.006	ng/u1	100
72) Phenanthrene	17.583	178	266261	19.380	ng/u1	100
74) Anthracene	17.677	178	270403	19.529	ng/u1	100
75) 1,2,3,4-Tetrachloroben...	13.599	216	76262	16.146	ng/uL	100
76) Pentachlorobenzene	15.109	250	72000	16.574	ng/uL	100
77) Carbazole	17.941	167	244776	21.442	ng/u1	100
78) Di-n-butylphthalate	18.494	149	287259	22.102	ng/u1	100
80) Fluoranthene	19.586	202	331640	21.975	ng/u1	100
82) Pyrene	19.951	202	328611	21.901	ng/u1	100
83) Butylbenzylphthalate	20.832	149	113267	25.049	ng/u1	100
84) 3,3'-Dichlorobenzidine	21.725	252	110447	20.583	ng/u1	100
85) Benzo(a)anthracene	21.819	228	312031	19.524	ng/u1	100
86) Bis(2-ethylhexyl)phtha...	21.725	149	168684	25.529	ng/u1	100
87) Chrysene	21.890	228	299016	19.502	ng/u1	100
89) Di-n-octyl phthalate	23.000	149	280101	24.371	ng/u1	100
90) Benzo(b)fluoranthene	24.140	252	323930	19.603	ng/u1	100
91) Benzo(k)fluoranthene	24.210	252	308354	19.488	ng/u1	100
93) Benzo(a)pyrene	25.057	252	315878	19.844	ng/u1	100
94) Indeno(1,2,3-cd)pyrene	29.105	276	362467	18.926	ng/u1	100
95) Dibenzo(a,h)anthracene	29.187	278	310480	19.264	ng/u1	100
96) Benzo(g,h,i)perylene	30.321	276	303349	18.760	ng/u1	100

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081722\
 Data File : BG054643.D
 Acq On : 17 Aug 2022 16:50
 Operator : CG/JU
 Sample : SST02037
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTD020437

Manual Integrations

APPROVED

Reviewed By :Jagrut Upadhyay 08/18/2022
 Supervised By :Sohil Jodhani 08/18/2022

Quant Time: Aug 18 02:23:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 18 02:20:33 2022
 Response via : Initial Calibration

