

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
 Data File : BG050967.D
 Acq On : 11 Nov 2021 14:02
 Operator : CG/JU
 Sample : PB140678BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SLCS678

Quant Time: Nov 11 14:57:48 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 11 12:40:48 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.225	152	35866	20.000	ng/u1	0.00
20) Naphthalene-d8	11.051	136	159156	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.847	164	105261	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.596	188	229016	20.000	ng/u1	0.00
79) Chrysene-d12	21.891	240	202187	20.000	ng/u1	0.00
88) Perylene-d12	25.281	264	203443	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.589	96	6310	5.678	ng/uL	0.00
4) Pyridine-d5	4.012	84	88326	26.569	ng/u1	0.00
7) Phenol-d5	7.367	99	101070	26.415	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.543	67	67229	27.200	ng/u1	0.00
11) 2-Chlorophenol-d4	7.755	132	71028	26.786	ng/u1	0.00
15) 4-Methylphenol-d8	8.924	113	79479	26.385	ng/u1	0.00
21) Nitrobenzene-d5	9.400	128	37793	27.943	ng/u1	0.00
24) 2-Nitrophenol-d4	10.123	143	41448	27.561	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.663	165	70806	27.950	ng/u1	0.00
31) 4-Chloroaniline-d4	11.180	131	91622	23.882	ng/u1	0.00
46) Dimethylphthalate-d6	14.241	166	230884	28.670	ng/u1	0.00
49) Acenaphthylene-d8	14.547	160	287450	28.650	ng/u1	0.00
54) 4-Nitrophenol-d4	15.035	143	39160	26.819	ng/u1	0.00
60) Fluorene-d10	15.840	176	199543	27.971	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.951	200	38207	27.517	ng/u1	0.00
73) Anthracene-d10	17.690	188	310078	28.638	ng/u1	0.00
81) Pyrene-d10	19.970	212	364636	27.922	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.046	264	315096	28.017	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.624	88	13703	11.227	ng/uL#	93
5) Pyridine	4.030	79	92566	26.899	ng/u1	96
6) Benzaldehyde	7.361	77	69955	28.987	ng/u1	96
8) Phenol	7.396	94	109228	27.596	ng/u1	99
10) Bis(2-Chloroethyl)ether	7.637	93	83241	28.097	ng/u1	98
12) 2-Chlorophenol	7.790	128	75596	28.078	ng/u1	98
13) 2-Methylphenol	8.660	108	79860	27.297	ng/u1	98
14) 2,2'-oxybis(1-Chloropr...	8.742	45	126048	27.011	ng/u1	99
16) Acetophenone	9.053	105	126060	26.939	ng/u1	95
17) N-Nitroso-di-n-propyla...	9.030	70	76521	27.103	ng/u1	99
18) 4-Methylphenol	8.989	108	84862	27.243	ng/u1	100
19) Hexachloroethane	9.318	117	30650	27.228	ng/u1	98
22) Nitrobenzene	9.441	77	110216	29.218	ng/u1	97
23) Isophorone	9.958	82	211489	28.888	ng/u1	100
25) 2-Nitrophenol	10.158	139	43454	28.801	ng/u1	97
26) 2,4-Dimethylphenol	10.199	107	95002	28.611	ng/u1	100
27) Bis(2-Chloroethoxy)met...	10.440	93	115538	29.290	ng/u1	98
29) 2,4-Dichlorophenol	10.693	162	70502	28.531	ng/u1	93
30) Naphthalene	11.104	128	249884	28.712	ng/u1	98
32) 4-Chloroaniline	11.210	127	94411	24.784	ng/u1	99
33) Hexachlorobutadiene	11.374	225	45724	28.193	ng/u1	96
34) Caprolactam	11.968	113	28812	27.465	ng/u1	89
35) 4-Chloro-3-methylphenol	12.308	107	91171	28.882	ng/u1	98
36) 2-Methylnaphthalene	12.690	142	168083	28.349	ng/u1	100

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
 Data File : BG050967.D
 Acq On : 11 Nov 2021 14:02
 Operator : CG/JU
 Sample : PB140678BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SLCS678

Quant Time: Nov 11 14:57:48 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 11 12:40:48 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
37) 1-Methylnaphthalene	12.908	142	169959	28.290	ng/ul	96
39) 1,2,4,5-Tetrachloroben...	13.055	216	90323	29.449	ng/ul	99
40) Hexachlorocyclopentadiene	13.025	237	36925	25.071	ng/ul	95
41) 2,4,6-Trichlorophenol	13.290	196	58883	29.340	ng/ul	95
42) 2,4,5-Trichlorophenol	13.360	196	62544	29.028	ng/ul	99
43) 1,1'-Biphenyl	13.683	154	226304	29.414	ng/ul	99
44) 2-Chloronaphthalene	13.736	162	175708	29.143	ng/ul	98
45) 2-Nitroaniline	13.936	65	69891	29.177	ng/ul	96
47) Dimethylphthalate	14.288	163	239367	29.727	ng/ul	100
48) 2,6-Dinitrotoluene	14.424	165	50687	30.076	ng/ul	93
50) Acenaphthylene	14.576	152	295597	29.396	ng/ul	99
51) 3-Nitroaniline	14.753	138	47071	27.005	ng/ul	93
52) Acenaphthene	14.911	153	193960	29.334	ng/ul	98
53) 2,4-Dinitrophenol	14.964	184	16711	17.974	ng/ul	95
55) 4-Nitrophenol	15.052	109	36334	27.132	ng/ul	94
56) Dibenzofuran	15.246	168	272713	28.815	ng/ul	98
57) 2,4-Dinitrotoluene	15.205	165	71118	29.569	ng/ul	98
58) 2,3,4,6-Tetrachlorophenol	15.469	232	47668	28.151	ng/ul	99
59) Diethylphthalate	15.646	149	250951	29.115	ng/ul	99
61) Fluorene	15.892	166	216558	28.909	ng/ul	98
62) 4-Chlorophenyl-phenyle...	15.875	204	113645	29.134	ng/ul	98
63) 4-Nitroaniline	15.916	138	50972	29.477	ng/ul	98
66) 4,6-Dinitro-2-methylph...	15.969	198	38032	28.088	ng/ul	99
67) N-Nitrosodiphenylamine	16.092	169	191333	29.889	ng/ul	98
68) 4-Bromophenyl-phenylether	16.774	248	68583	30.106	ng/ul	95
69) Hexachlorobenzene	16.897	284	69470	29.663	ng/ul	97
70) Atrazine	17.032	200	78899	29.064	ng/ul	99
71) Pentachlorophenol	17.238	266	27513	25.588	ng/ul	90
72) Phenanthrene	17.637	178	365974	29.937	ng/ul	99
74) Anthracene	17.726	178	358710	29.245	ng/ul	99
75) 1,2,3,4-Tetrachloroben...	13.654	216	94450	30.289	ng/ul	97
76) Pentachlorobenzene	15.164	250	83983	29.067	ng/ul	98
77) Carbazole	17.996	167	338555	30.795	ng/ul	98
78) Di-n-butylphthalate	18.531	149	432694	29.952	ng/ul	100
80) Fluoranthene	19.635	202	450967	28.775	ng/ul	98
82) Pyrene	19.999	202	443679	28.973	ng/ul	99
83) Butylbenzylphthalate	20.863	149	194814	29.584	ng/ul	97
84) 3,3'-Dichlorobenzidine	21.774	252	126554	25.702	ng/ul	99
85) Benzo(a)anthracene	21.868	228	411891	29.427	ng/ul	99
86) Bis(2-ethylhexyl)phtha...	21.738	149	280512	29.676	ng/ul	99
87) Chrysene	21.938	228	392901	29.384	ng/ul	99
89) Di-n-octyl phthalate	23.002	149	474316	28.638	ng/ul	100
90) Benzo(b)fluoranthene	24.194	252	412390	28.454	ng/ul	100
91) Benzo(k)fluoranthene	24.271	252	389110	28.611	ng/ul	98
93) Benzo(a)pyrene	25.123	252	395009	28.616	ng/ul	99
94) Indeno(1,2,3-cd)pyrene	29.177	276	443381	28.855	ng/ul	96
95) Dibenzo(a,h)anthracene	29.247	278	375160	28.855	ng/ul	98
96) Benzo(g,h,i)perylene	30.405	276	369879	28.756	ng/ul	99

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
 Data File : BG050967.D
 Acq On : 11 Nov 2021 14:02
 Operator : CG/JU
 Sample : PB140678BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SLCS678

Quant Time: Nov 11 14:57:48 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 11 12:40:48 2021
 Response via : Initial Calibration

