

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101520\
 Data File : BG046917.D
 Acq On : 15 Oct 2020 17:47
 Operator : CG/JU
 Sample : SSTD01029
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTD01029

Manual Integrations
 APPROVED

mohammad
 10/16/2020 12:46:58 PM

Quant Time: Oct 15 19:09:33 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG101520MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 15 19:04:31 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	46281	20.00	ng/ul	0.00
18) Naphthalene-d8	10.89	136	176679	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.70	164	128887	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.45	188	322794	20.00	ng/ul	0.00
78) Chrysene-d12	21.72	240	402346	20.00	ng/ul	0.00
86) Perylene-d12	24.97	264	404280	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.49	96	3764	3.99	ng/uL	0.00
5) Phenol-d5	7.23	99	37616	9.50	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.40	67	24455	10.12	ng/ul	0.00
9) 2-Chlorophenol-d4	7.61	132	30648	10.60	ng/ul	0.00
13) 4-Methylphenol-d8	8.78	113	30381	9.81	ng/ul	0.00
19) Nitrobenzene-d5	9.24	128	14931	11.29	ng/ul	0.00
22) 2-Nitrophenol-d4	9.96	143	16231	12.08	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.50	165	32513	10.28	ng/ul	0.00
29) 4-Chloroaniline-d4	11.02	131	40902	10.89	ng/ul	0.00
44) Dimethylphthalate-d6	14.10	166	105856	10.75	ng/ul	0.00
47) Acenaphthylene-d8	14.40	160	122296	10.46	ng/ul	0.00
52) 4-Nitrophenol-d4	14.89	143	17038	10.15	ng/ul	0.00
58) Fluorene-d10	15.70	176	99640	10.97	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.80	200	19739	12.24	ng/ul	0.00
71) Anthracene-d10	17.55	188	161073	10.88	ng/ul	0.00
79) Pyrene-d10	19.83	212	216030	10.35	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.75	264	225402	10.34	ng/ul	0.00

Target Compounds

					Ovalue
2) 1,4-Dioxane	3.53	88	5382	4.453 ng/uL#	88
4) Benzaldehyde	7.21	77	29108	10.419 ng/ul	93
6) Phenol	7.25	94	40881	9.852 ng/ul	96
8) Bis(2-Chloroethyl)ether	7.49	93	33014	10.332 ng/ul	91
10) 2-Chlorophenol	7.64	128	30423	10.320 ng/ul	94
11) 2-Methylphenol	8.51	108	28935	9.459 ng/ul	93
12) 2,2'-oxybis(1-Chloropropan	8.61	45	46623m	9.273 ng/ul	
14) Acetophenone	8.90	105	51917	10.338 ng/ul	94
15) N-Nitroso-di-n-propylamine	8.88	70	27301	10.065 ng/ul	97
16) 4-Methylphenol	8.84	108	31272	9.584 ng/ul	96
17) Hexachloroethane	9.17	117	14690	10.728 ng/ul	93
20) Nitrobenzene	9.28	77	39991	10.601 ng/ul	96
21) Isophorone	9.80	82	73090	9.714 ng/ul	98
23) 2-Nitrophenol	10.00	139	17901	12.227 ng/ul#	91
24) 2,4-Dimethylphenol	10.05	107	38714	10.556 ng/ul	94
25) Bis(2-Chloroethoxy)methane	10.29	93	42257	9.641 ng/ul	99
27) 2,4-Dichlorophenol	10.53	162	32532	10.570 ng/ul	99
28) Naphthalene	10.94	128	101797	10.407 ng/ul	98
30) 4-Chloroaniline	11.04	127	39904	10.908 ng/ul	94
31) Hexachlorobutadiene	11.23	225	30268	11.444 ng/ul	89
32) Caprolactam	11.80	113	10981m	10.340 ng/ul	
33) 4-Chloro-3-methylphenol	12.15	107	33603	9.990 ng/ul	95
34) 2-Methylnaphthalene	12.54	142	74671	10.442 ng/ul	100

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101520\
 Data File : BG046917.D
 Acq On : 15 Oct 2020 17:47
 Operator : CG/JU
 Sample : SSTD01029
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTD01029

Manual Integrations
 APPROVED

mohammad
 10/16/2020 12:46:58 PM

Quant Time: Oct 15 19:09:33 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG101520MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 15 19:04:31 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
35) 1-Methylnaphthalene	12.76	142	73375	10.588	ng/uL#	97
37) 1,2,4,5-Tetrachlorobenzene	12.91	216	51934	11.006	ng/uL#	89
38) Hexachlorocyclopentadiene	12.89	237	30211	9.901	ng/uL#	95
39) 2,4,6-Trichlorophenol	13.14	196	29075	10.412	ng/uL	91
40) 2,4,5-Trichlorophenol	13.21	196	31330	10.572	ng/uL	91
41) 1,1'-Biphenyl	13.54	154	105187	10.610	ng/uL	99
42) 2-Chloronaphthalene	13.59	162	85725	10.795	ng/uL	95
43) 2-Nitroaniline	13.78	65	24003	11.411	ng/uL	100
45) Dimethylphthalate	14.15	163	111490	11.186	ng/uL	99
46) 2,6-Dinitrotoluene	14.27	165	22054	12.504	ng/uL#	87
48) Acenaphthylene	14.43	152	127029	11.026	ng/uL	97
49) 3-Nitroaniline	14.60	138	19318	10.997	ng/uL#	86
50) Acenaphthene	14.77	153	86661	10.570	ng/uL	97
51) 2,4-Dinitrophenol	14.81	184	11032	10.414	ng/uL	93
53) 4-Nitrophenol	14.90	109	18067	10.862	ng/uL	91
54) Dibenzofuran	15.10	168	131211	10.892	ng/uL	97
55) 2,4-Dinitrotoluene	15.06	165	31299	12.826	ng/uL	92
56) 2,3,4,6-Tetrachlorophenol	15.33	232	30108	10.904	ng/uL	91
57) Diethylphthalate	15.51	149	112081	11.254	ng/uL	95
59) Fluorene	15.75	166	106938	10.932	ng/uL	97
60) 4-Chlorophenyl-phenylether	15.74	204	61508	10.923	ng/uL	95
61) 4-Nitroaniline	15.76	138	22118	11.451	ng/uL	97
64) 4,6-Dinitro-2-methylphenol	15.82	198	20300	11.463	ng/uL#	94
65) N-Nitrosodiphenylamine	15.96	169	94794	10.202	ng/uL	95
66) 4-Bromophenyl-phenylether	16.64	248	41654	10.334	ng/uL	97
67) Hexachlorobenzene	16.75	284	44806	12.895	ng/uL#	92
68) Atrazine	16.90	200	42876	11.139	ng/uL	98
69) Pentachlorophenol	17.10	266	23589	9.286	ng/uL	98
70) Phenanthrene	17.49	178	192027	10.882	ng/uL	99
72) Anthracene	17.58	178	191434	10.707	ng/uL	98
73) 1,2,3,4-Tetrachlorobenzene	13.51	216	51353	10.310	ng/uL	96
74) Pentachlorobenzene	15.02	250	52627	10.398	ng/uL	95
75) Carbazole	17.85	167	163203	11.452	ng/uL	98
76) Di-n-butylphthalate	18.41	149	209988	11.881	ng/uL	99
77) Fluoranthene	19.50	202	263932	12.595	ng/uL	99
80) Pvrene	19.86	202	275213	10.642	ng/uL	98
81) Butylbenzylphthalate	20.74	149	96907	10.752	ng/uL	97
82) 3,3'-Dichlorobenzidine	21.61	252	99516	10.576	ng/uL	100
83) Benzo(a)anthracene	21.70	228	287072	10.905	ng/uL	99
84) Bis(2-ethylhexyl)phthalate	21.61	149	142237	10.612	ng/uL	96
85) Chrysene	21.77	228	274397	10.745	ng/uL	99
87) Di-n-octyl phthalate	22.82	149	239418	10.843	ng/uL	100
88) Benzo(b)fluoranthene	23.93	252	281105	10.540	ng/uL	98
89) Benzo(k)fluoranthene	24.00	252	282434	10.864	ng/uL	99
91) Benzo(a)pyrene	24.82	252	252921	10.382	ng/uL	98
92) Indeno(1,2,3-cd)pyrene	28.66	276	299669	10.026	ng/uL	97
93) Dibenzo(a,h)anthracene	28.73	278	258778	10.449	ng/uL	100
94) Benzo(a,h,i)perylene	29.83	276	259868	10.378	ng/uL	98

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101520\
Data File : BG046917.D
Acq On : 15 Oct 2020 17:47
Operator : CG/JU
Sample : SSTD01029
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SSTD01029

Manual Integrations
APPROVED

mohammad
10/16/2020 12:46:58 PM

Quant Time: Oct 15 19:09:33 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG101520MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Oct 15 19:04:31 2020
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

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

