

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120821\
 Data File : BG051416.D
 Acq On : 8 Dec 2021 21:07
 Operator : CG/JU
 Sample : SSTD16033
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTD160433

Quant Time: Dec 09 03:08:58 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120821.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 09 02:58:27 2021
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 12/09/2021
 Supervised By :Yogesh Patel 12/16/2021

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.189	152	26055	20.000	ng/u1	0.00
20) Naphthalene-d8	11.021	136	119785	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.822	164	76334	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.578	188	171395	20.000	ng/u1	0.00
79) Chrysene-d12	21.885	240	144771	20.000	ng/u1	0.01
88) Perylene-d12	25.287	264	142057	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	3.959	84	353749	160.466	ng/u1	0.00
7) Phenol-d5	7.360	99	417562	162.261	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.513	67	258637	155.258	ng/u1	0.01
11) 2-Chlorophenol-d4	0.000	132	0d	0.000	ng/u1	
15) 4-Methylphenol-d8	8.923	113	323927	159.328	ng/u1	0.01
21) Nitrobenzene-d5	0.000	128	0d	0.000	ng/u1	
24) 2-Nitrophenol-d4	0.000	143	0d	0.000	ng/u1	
28) 2,4-Dichlorophenol-d3	0.000	165	0d	0.000	ng/u1	
31) 4-Chloroaniline-d4	11.168	131	419352	153.691	ng/u1	0.00
46) Dimethylphthalate-d6	0.000	166	0d	0.000	ng/u1	
49) Acenaphthylene-d8	0.000	160	0d	0.000	ng/u1	
54) 4-Nitrophenol-d4	15.075	143	152556	184.285	ng/u1	0.01
60) Fluorene-d10	0.000	176	0d	0.000	ng/u1	
65) 4,6-Dinitro-2-methylph...	15.968	200	174260	178.825	ng/u1	0.02
73) Anthracene-d10	0.000	188	0d	0.000	ng/u1	
81) Pyrene-d10	0.000	212	0d	0.000	ng/u1	
92) Benzo(a)pyrene-d12	0.000	264	0d	0.000	ng/u1	
Target Compounds						
					Qvalue	
5) Pyridine	3.982	79	362773	152.543	ng/u1	98
6) Benzaldehyde	7.325	77	112293	66.594	ng/u1	96
8) Phenol	7.396	94	416853	153.542	ng/u1	97
10) Bis(2-Chloroethyl)ether	7.607	93	313048	150.541	ng/u1	98
13) 2-Methylphenol	8.647	108	317770	157.211	ng/u1	100
14) 2,2'-oxybis(1-Chloropr...	8.712	45	458425	146.502	ng/u1	96
16) Acetophenone	9.035	105	465889	144.389	ng/u1	98
18) 4-Methylphenol	8.994	108	321176	151.255	ng/u1	96
32) 4-Chloroaniline	11.191	127	417099	148.204	ng/u1	100
34) Caprolactam	12.008	113	117574m	151.272	ng/u1	
40) Hexachlorocyclopentadiene	12.989	237	198925	158.048	ng/u1	98
51) 3-Nitroaniline	14.758	138	155927	128.538	ng/u1	96
53) 2,4-Dinitrophenol	14.975	184	117319	180.087	ng/u1	92
55) 4-Nitrophenol	15.093	109	140475	158.788	ng/u1	93
63) 4-Nitroaniline	15.933	138	122288	113.564	ng/u1	94
66) 4,6-Dinitro-2-methylph...	15.986	198	164805	166.432	ng/u1#	99
70) Atrazine	17.031	200	305831	148.346	ng/u1	98
71) Pentachlorophenol	17.237	266	141610	185.431	ng/u1	97
77) Carbazole	17.989	167	1188056	144.377	ng/u1	96
84) 3,3'-Dichlorobenzidine	21.767	252	418747	136.971	ng/u1	98
89) Di-n-octyl phthalate	22.977	149	1564487	149.752	ng/u1	100

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