

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042924\
 Data File : BN031309.D
 Acq On : 29 Apr 2024 15:41
 Operator : MA/JU
 Sample : SSTD16049
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD160229

Quant Time: Apr 29 16:15:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 29 14:23:49 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 04/30/2024
 Supervised By :mohammad ahmed 05/02/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	249914	20.000	ng/u1	0.00
20) Naphthalene-d8	10.404	136	1123104	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.269	164	702643	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.028	188	1425930	20.000	ng/u1	0.01
79) Chrysene-d12	21.269	240	923491	20.000	ng/u1	0.01
88) Perylene-d12	24.157	264	918737	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	3.540	84	2327487	148.269	ng/u1	0.00
7) Phenol-d5	6.816	99	3082568	155.308	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	6.975	67	1744939	144.516	ng/u1	0.01
11) 2-Chlorophenol-d4	0.000	132	0d	0.000	ng/u1	
15) 4-Methylphenol-d8	8.352	113	2560433	153.992	ng/u1	0.02
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	10.557	131	3311657	134.273	ng/u1	0.02
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	14.528	143	1368407	159.160	ng/u1	0.04
60) Fluorene-d10	0.000	176	0d	0.000	ng/u1	
65) 4,6-Dinitro-2-methylph...	15.422	200	1234951	191.452	ng/u1	0.03
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	3.558	79	2259878	143.261	ng/u1	99
6) Benzaldehyde	6.775	77	1120098	121.663	ng/u1	97
8) Phenol	6.846	94	3037850	147.647	ng/u1	96
10) Bis(2-Chloroethyl)ether	7.069	93	2409070	144.133	ng/u1	99
13) 2-Methylphenol	8.075	108	2412214	151.675	ng/u1	98
14) 2,2'-oxybis(1-Chloropr...	8.157	45	3166917	136.019	ng/u1	99
16) Acetophenone	8.457	105	3386436	130.783	ng/u1	98
18) 4-Methylphenol	8.428	108	2551037	143.924	ng/u1	98
32) 4-Chloroaniline	10.587	127	3106278	130.822	ng/u1	98
34) Caprolactam	11.440	113	931236m	162.000	ng/u1	
40) Hexachlorocyclopentadiene	12.422	237	1745117	174.274	ng/u1	92
51) 3-Nitroaniline	14.210	138	1478845	159.551	ng/u1	99
53) 2,4-Dinitrophenol	14.416	184	1046405	245.003	ng/u1	90
55) 4-Nitrophenol	14.545	109	1102972	162.234	ng/u1	93
63) 4-Nitroaniline	15.398	138	1267231	142.529	ng/u1	94
66) 4,6-Dinitro-2-methylph...	15.439	198	1273315	184.069	ng/u1#	97
70) Atrazine	16.516	200	1715848	132.711	ng/u1	94
71) Pentachlorophenol	16.681	266	1522687	165.710	ng/u1	95
77) Carbazole	17.445	167	8471565	135.696	ng/u1	98
84) 3,3'-Dichlorobenzidine	21.186	252	2367226	144.255	ng/u1	97
89) Di-n-octyl phthalate	22.269	149	10305257	177.365	ng/u1	100

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