

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071221\
 Data File : BN015493.D
 Acq On : 12 Jul 2021 17:08
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4119

Quant Time: Jul 12 17:39:07 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 12 15:47:53 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.808	152	3052	0.400	ng/ul	0.00
4) Naphthalene-d8	10.601	136	10312	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.446	164	5258	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.191	188	9926	0.400	ng/ul	0.00
17) Chrysene-d12	21.389	240	8669	0.400	ng/ul	0.00
23) Perylene-d12	23.745	264	7881	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	1279	0.372	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.190	152	5712	0.353	ng/ul	0.00
18) Fluoranthene-d10	19.224	212	9984	0.350	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	1390	0.374	ng/ul	99
5) Naphthalene	10.645	128	11884	0.361	ng/ul	100
7) 2-Methylnaphthalene	12.267	142	7588	0.353	ng/ul	99
8) 1-Methylnaphthalene	12.481	142	7489	0.360	ng/ul	98
10) Acenaphthylene	14.164	152	9965	0.367	ng/ul	100
11) Acenaphthene	14.506	153	7471	0.359	ng/ul	99
12) Fluorene	15.496	166	7848	0.341	ng/ul	99
14) Pentachlorophenol	16.845	266	642	0.292	ng/ul	99
15) Phenanthrene	17.234	178	13027	0.362	ng/ul	99
16) Anthracene	17.327	178	11203	0.353	ng/ul	99
19) Fluoranthene	19.257	202	14325	0.358	ng/ul	98
20) Pyrene	19.619	202	14665	0.365	ng/ul	99
21) Benzo(a)anthracene	21.375	228	12733	0.397	ng/ul	98
22) Chrysene	21.424	228	13057	0.340	ng/ul	99
24) Benzo(b)fluoranthene	23.029	252	12647	0.344	ng/ul	98
25) Benzo(k)fluoranthene	23.075	252	12877	0.310	ng/ul	98
26) Benzo(a)pyrene	23.640	252	11055	0.357	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.169	276	13877	0.372	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.189	278	10924	0.373	ng/ul	99
29) Benzo(g,h,i)perylene	26.910	276	11767	0.340	ng/ul	100

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

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