

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102022\
 Data File : BN022200.D
 Acq On : 19 Oct 2022 17:32
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4316

Quant Time: Oct 20 01:07:15 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 19 23:32:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.953	152	3856	0.400	ng/ul	0.00
4) Naphthalene-d8	10.783	136	11487	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.628	164	6530	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.383	188	14416	0.400	ng/ul	0.00
17) Chrysene-d12	21.585	240	11790	0.400	ng/ul	0.00
23) Perylene-d12	24.049	264	9311	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.261	96	2061	0.411	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.378	152	7045	0.405	ng/ul	0.00
18) Fluoranthene-d10	19.417	212	16070	0.406	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.295	88	2046	0.398	ng/ul	99
5) Naphthalene	10.833	128	13934	0.411	ng/ul	100
7) 2-Methylnaphthalene	12.450	142	8669	0.406	ng/ul	100
8) 1-Methylnaphthalene	12.670	142	8958	0.407	ng/ul	100
10) Acenaphthylene	14.346	152	11625	0.397	ng/ul	99
11) Acenaphthene	14.688	153	10114	0.409	ng/ul	99
12) Fluorene	15.678	166	11718	0.408	ng/ul	99
14) Pentachlorophenol	17.032	266	1235	0.360	ng/ul	99
15) Phenanthrene	17.425	178	19847	0.410	ng/ul	99
16) Anthracene	17.518	178	17274	0.424	ng/ul	100
19) Fluoranthene	19.445	202	21752	0.403	ng/ul	100
20) Pyrene	19.812	202	21891	0.404	ng/ul	100
21) Benzo(a)anthracene	21.567	228	18386	0.416	ng/ul	99
22) Chrysene	21.620	228	20787	0.444	ng/ul	99
24) Benzo(b)fluoranthene	23.292	252	19467	0.416	ng/ul	98
25) Benzo(k)fluoranthene	23.339	252	19720	0.434	ng/ul	99
26) Benzo(a)pyrene	23.938	252	16028	0.430	ng/ul#	95
27) Indeno(1,2,3-cd)pyrene	26.632	276	19545	0.422	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.655	278	16215	0.438	ng/ul	99
29) Benzo(g,h,i)perylene	27.430	276	14990	0.397	ng/ul	99

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

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