

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN010423\
 Data File : BN023546.D
 Acq On : 04 Jan 2023 09:54
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4366

Quant Time: Jan 05 04:35:07 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN010323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 04 00:28:51 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	5192	0.400	ng/u1	0.00
4) Naphthalene-d8	10.766	136	15748	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.604	164	9023	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.352	188	19949	0.400	ng/u1	0.00
17) Chrysene-d12	21.543	240	18683	0.400	ng/u1	0.00
23) Perylene-d12	23.969	264	14967	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.302	96	2197	0.382	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.355	152	8626	0.356	ng/u1	0.00
18) Fluoranthene-d10	19.383	212	19756	0.361	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.340	88	2192	0.364	ng/u1#	87
5) Naphthalene	10.815	128	16522	0.365	ng/u1	99
7) 2-Methylnaphthalene	12.427	142	10338	0.355	ng/u1	100
8) 1-Methylnaphthalene	12.647	142	10666	0.358	ng/u1	100
10) Acenaphthylene	14.322	152	12659	0.344	ng/u1	99
11) Acenaphthene	14.664	153	11690	0.360	ng/u1	98
12) Fluorene	15.654	166	13280	0.356	ng/u1	100
14) Pentachlorophenol	17.002	266	2024	0.315	ng/u1	99
15) Phenanthrene	17.394	178	22524	0.361	ng/u1	100
16) Anthracene	17.483	178	18360	0.343	ng/u1	100
19) Fluoranthene	19.411	202	26154	0.370	ng/u1	98
20) Pyrene	19.773	202	26549	0.357	ng/u1	97
21) Benzo(a)anthracene	21.525	228	23819	0.345	ng/u1	99
22) Chrysene	21.581	228	25549	0.367	ng/u1	99
24) Benzo(b)fluoranthene	23.227	252	26596	0.384	ng/u1	99
25) Benzo(k)fluoranthene	23.273	252	24606	0.377	ng/u1	99
26) Benzo(a)pyrene	23.858	252	21011	0.365	ng/u1	98
27) Indeno(1,2,3-cd)pyrene	26.493	276	27979	0.391	ng/u1#	99
28) Dibenzo(a,h)anthracene	26.510	278	22247	0.392	ng/u1	100
29) Benzo(g,h,i)perylene	27.271	276	22769	0.398	ng/u1	100

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

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