

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101823\
 Data File : BN028325.D
 Acq On : 18 Oct 2023 16:06
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4300

Quant Time: Oct 18 16:48:28 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN100323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 16:16:04 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.882	152	2304	0.400	ng/ul	0.00
4) Naphthalene-d8	10.702	136	7403	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.536	164	4245	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.299	188	9484	0.400	ng/ul	0.00
17) Chrysene-d12	21.495	240	9672	0.400	ng/ul	0.00
23) Perylene-d12	23.945	264	9950	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	1252	0.433	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.297	152	4388	0.368	ng/ul	0.00
18) Fluoranthene-d10	19.329	212	10892	0.355	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.296	88	1237	0.404	ng/ul	95
5) Naphthalene	10.757	128	8029	0.382	ng/ul	97
7) 2-Methylnaphthalene	12.374	142	4902	0.370	ng/ul	98
8) 1-Methylnaphthalene	12.583	142	5097	0.371	ng/ul	100
10) Acenaphthylene	14.268	152	6888	0.358	ng/ul	98
11) Acenaphthene	14.601	153	5969	0.389	ng/ul	99
12) Fluorene	15.596	166	6782	0.398	ng/ul	99
14) Pentachlorophenol	16.953	266	541	0.266	ng/ul	93
15) Phenanthrene	17.341	178	11126	0.394	ng/ul	99
16) Anthracene	17.439	178	9302	0.368	ng/ul	97
19) Fluoranthene	19.357	202	13271	0.359	ng/ul	99
20) Pyrene	19.729	202	13740	0.359	ng/ul	100
21) Benzo(a)anthracene	21.481	228	12845	0.364	ng/ul	99
22) Chrysene	21.533	228	13951	0.391	ng/ul	99
24) Benzo(b)fluoranthene	23.185	252	14657	0.397	ng/ul	99
25) Benzo(k)fluoranthene	23.234	252	15132	0.412	ng/ul	99
26) Benzo(a)pyrene	23.837	252	12345	0.392	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.512	276	17262	0.454	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.529	278	13929	0.450	ng/ul	99
29) Benzo(g,h,i)perylene	27.320	276	14943	0.480	ng/ul	99

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

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