

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090523\
 Data File : BN027368.D
 Acq On : 05 Sep 2023 10:33
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4255

Quant Time: Sep 06 00:41:05 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 02 06:16:48 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.059	152	3520	0.400	ng/ul	0.00
4) Naphthalene-d8	10.883	136	11275	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.689	164	6383	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.443	188	15405	0.400	ng/ul	0.00
17) Chrysene-d12	21.618	240	13065	0.400	ng/ul	0.00
23) Perylene-d12	24.143	264	10897	0.400	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.418	96	1973	0.421	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.456	152	7270	0.425	ng/ul	0.00
18) Fluoranthene-d10	19.459	212	18932	0.423	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.456	88	2115	0.446	ng/ul#	81
5) Naphthalene	10.933	128	14122	0.432	ng/ul	100
7) 2-Methylnaphthalene	12.527	142	8727	0.427	ng/ul	99
8) 1-Methylnaphthalene	12.747	142	9156	0.420	ng/ul	99
10) Acenaphthylene	14.416	152	12298	0.421	ng/ul	99
11) Acenaphthene	14.749	153	10780	0.429	ng/ul	100
12) Fluorene	15.739	166	12541	0.445	ng/ul	100
14) Pentachlorophenol	17.075	266	1207	0.402	ng/ul	98
15) Phenanthrene	17.485	178	21472	0.428	ng/ul	99
16) Anthracene	17.578	178	17038	0.427	ng/ul	100
19) Fluoranthene	19.492	202	25157	0.419	ng/ul	99
20) Pyrene	19.854	202	26075	0.426	ng/ul	99
21) Benzo(a)anthracene	21.603	228	18831	0.400	ng/ul	99
22) Chrysene	21.659	228	23270	0.424	ng/ul	100
24) Benzo(b)fluoranthene	23.360	252	21233	0.413	ng/ul	99
25) Benzo(k)fluoranthene	23.410	252	21636	0.430	ng/ul	99
26) Benzo(a)pyrene	24.029	252	16552	0.412	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.820	276	17896	0.401	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.850	278	13050	0.382	ng/ul	98
29) Benzo(g,h,i)perylene	27.639	276	17875	0.444	ng/ul	99

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

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