

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090523\
 Data File : BN027394.D
 Acq On : 06 Sep 2023 03:22
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4257

Quant Time: Sep 06 03:49:29 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	8.055	152	3477	0.400	ng/ul	0.00
4) Naphthalene-d8	10.878	136	10690	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.689	164	6043	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.439	188	14102	0.400	ng/ul	0.00
17) Chrysene-d12	21.615	240	12572	0.400	ng/ul	0.00
23) Perylene-d12	24.138	264	11126	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.422	96	35	0.008	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.456	152	6232	0.384	ng/ul	0.00
18) Fluoranthene-d10	19.459	212	15942	0.370	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.456	88	1758	0.376	ng/ul	98
5) Naphthalene	10.933	128	12543	0.404	ng/ul	99
7) 2-Methylnaphthalene	12.528	142	8165	0.421	ng/ul	99
8) 1-Methylnaphthalene	12.748	142	8035	0.389	ng/ul	100
10) Acenaphthylene	14.416	152	11837	0.428	ng/ul	100
11) Acenaphthene	14.749	153	9566	0.402	ng/ul	100
12) Fluorene	15.734	166	11052	0.414	ng/ul	99
14) Pentachlorophenol	17.071	266	1334	0.485	ng/ul	99
15) Phenanthrene	17.481	178	18434	0.401	ng/ul	99
16) Anthracene	17.574	178	14964	0.410	ng/ul	100
19) Fluoranthene	19.487	202	21933	0.379	ng/ul	98
20) Pyrene	19.854	202	22939	0.390	ng/ul	97
21) Benzo(a)anthracene	21.597	228	17354	0.383	ng/ul	99
22) Chrysene	21.653	228	20959	0.396	ng/ul	99
24) Benzo(b)fluoranthene	23.351	252	19188	0.365	ng/ul	97
25) Benzo(k)fluoranthene	23.407	252	19654	0.383	ng/ul	97
26) Benzo(a)pyrene	24.024	252	16626	0.405	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.807	276	17475	0.383	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.841	278	13323	0.382	ng/ul	98
29) Benzo(g,h,i)perylene	27.632	276	16701	0.406	ng/ul	100

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

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