

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062422\
 Data File : BN020403.D
 Acq On : 24 Jun 2022 18:28
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4378

Quant Time: Jun 25 00:55:25 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 00:17:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.753	152	2492	0.400	ng/u1	0.00
4) Naphthalene-d8	10.534	136	8211	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.381	164	5499	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.120	188	12738	0.400	ng/u1	0.00
17) Chrysene-d12	21.310	240	11587	0.400	ng/u1	0.00
23) Perylene-d12	23.593	264	9673	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	1223	0.424	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.129	152	5777	0.429	ng/u1	0.00
18) Fluoranthene-d10	19.150	212	15169	0.397	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.239	88	1266	0.429	ng/u1	95
5) Naphthalene	10.584	128	9667	0.381	ng/u1	99
7) 2-Methylnaphthalene	12.206	142	6449	0.380	ng/u1	99
8) 1-Methylnaphthalene	12.421	142	6965	0.433	ng/u1	100
10) Acenaphthylene	14.099	152	9764	0.375	ng/u1	99
11) Acenaphthene	14.441	153	7856	0.377	ng/u1	99
12) Fluorene	15.431	166	9267	0.375	ng/u1	100
14) Pentachlorophenol	16.786	266	733	0.285	ng/u1	99
15) Phenanthrene	17.162	178	16458	0.372	ng/u1	100
16) Anthracene	17.255	178	14097	0.369	ng/u1	100
19) Fluoranthene	19.182	202	18717	0.354	ng/u1	99
20) Pyrene	19.545	202	19452	0.354	ng/u1	100
21) Benzo(a)anthracene	21.296	228	15363	0.369	ng/u1	99
22) Chrysene	21.348	228	17606	0.380	ng/u1	100
24) Benzo(b)fluoranthene	22.903	252	16502	0.369	ng/u1	99
25) Benzo(k)fluoranthene	22.950	252	16939	0.375	ng/u1	99
26) Benzo(a)pyrene	23.496	252	15344	0.377	ng/u1	98
27) Indeno(1,2,3-cd)pyrene	25.927	276	15968	0.394	ng/u1	98
28) Dibenzo(a,h)anthracene	25.941	278	12920	0.395	ng/u1	98
29) Benzo(g,h,i)perylene	26.642	276	13602	0.391	ng/u1	99

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

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