

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062922\
 Data File : BN020565.D
 Acq On : 29 Jun 2022 12:23
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4383

Quant Time: Jun 29 12:52:32 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	7.741	152	3380	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.518	136	11101	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.368	164	6443	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.112	188	12713	0.400	ng/ul	0.00
17) Chrysene-d12	21.303	240	9349	0.400	ng/ul	0.00
23) Perylene-d12	23.586	264	6797	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.197	96	1568	0.400	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.119	152	6735	0.370	ng/ul	0.00
18) Fluoranthene-d10	19.141	212	13473	0.437	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.227	88	1607	0.402	ng/ul	94
5) Naphthalene	10.568	128	12138	0.354	ng/ul	99
7) 2-Methylnaphthalene	12.190	142	7574	0.330	ng/ul	99
8) 1-Methylnaphthalene	12.410	142	8236	0.379	ng/ul	100
10) Acenaphthylene	14.090	152	10718	0.351	ng/ul	99
11) Acenaphthene	14.432	153	8612	0.352	ng/ul	98
12) Fluorene	15.418	166	9639	0.333	ng/ul	100
14) Pentachlorophenol	16.778	266	522	0.204	ng/ul	98
15) Phenanthrene	17.150	178	15586	0.353	ng/ul	100
16) Anthracene	17.247	178	12900	0.338	ng/ul	100
19) Fluoranthene	19.174	202	16804	0.394	ng/ul	99
20) Pyrene	19.536	202	16994	0.384	ng/ul	99
21) Benzo(a)anthracene	21.289	228	10963	0.326	ng/ul	100
22) Chrysene	21.341	228	13627	0.365	ng/ul	99
24) Benzo(b)fluoranthene	22.896	252	11004	0.350	ng/ul	99
25) Benzo(k)fluoranthene	22.940	252	11362	0.358	ng/ul	99
26) Benzo(a)pyrene	23.484	252	9996	0.350	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.916	276	10354	0.364	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.933	278	8311	0.362	ng/ul	99
29) Benzo(g,h,i)perylene	26.624	276	9351	0.383	ng/ul	99

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

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