

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070122\
 Data File : BN020600.D
 Acq On : 01 Jul 2022 17:44
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 70 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4389

Quant Time: Jul 02 00:49:00 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 : Sat Jul 02 00:47:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.830	152	3279	0.400	ng/ul	0.00
4) Naphthalene-d8	10.628	136	9933	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.479	164	5393	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.234	188	10819	0.400	ng/ul	0.00
17) Chrysene-d12	21.441	240	9348	0.400	ng/ul	0.00
23) Perylene-d12	23.794	264	7968	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.227	96	1546	0.407	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.229	152	5844	0.359	ng/ul	0.00
18) Fluoranthene-d10	19.267	212	11610	0.377	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.260	88	1572	0.405	ng/ul	95
5) Naphthalene	10.678	128	10803	0.352	ng/ul	100
7) 2-Methylnaphthalene	12.306	142	6572	0.320	ng/ul	99
8) 1-Methylnaphthalene	12.520	142	7111	0.365	ng/ul	100
10) Acenaphthylene	14.201	152	9011	0.353	ng/ul	99
11) Acenaphthene	14.544	153	7308	0.357	ng/ul	99
12) Fluorene	15.534	166	7985	0.329	ng/ul	99
14) Pentachlorophenol	16.901	266	489	0.224	ng/ul	99
15) Phenanthrene	17.272	178	13172	0.350	ng/ul	100
16) Anthracene	17.369	178	10824	0.333	ng/ul	100
19) Fluoranthene	19.299	202	14487	0.339	ng/ul	100
20) Pyrene	19.662	202	14828	0.335	ng/ul	97
21) Benzo(a)anthracene	21.426	228	10758	0.320	ng/ul	100
22) Chrysene	21.476	228	13280	0.355	ng/ul	100
24) Benzo(b)fluoranthene	23.077	252	12435	0.337	ng/ul	99
25) Benzo(k)fluoranthene	23.127	252	13218	0.355	ng/ul	99
26) Benzo(a)pyrene	23.691	252	11412	0.341	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.228	276	12888	0.386	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.255	278	10280	0.382	ng/ul	99
29) Benzo(g,h,i)perylene	26.976	276	11009	0.385	ng/ul	99

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

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