

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071922\
 Data File : BN020794.D
 Acq On : 20 Jul 2022 03:41
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4396

Quant Time: Jul 20 04:19:26 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 : Wed Jul 20 01:02:30 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.821	152	2022	0.400	ng/ul	0.00
4) Naphthalene-d8	10.617	136	8894	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.469	164	5402	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.221	188	11619	0.400	ng/ul	0.00
17) Chrysene-d12	21.422	240	11513	0.400	ng/ul	0.00
23) Perylene-d12	23.769	264	9787	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	1022	0.436	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.218	152	5052	0.346	ng/ul	0.00
18) Fluoranthene-d10	19.258	212	12539	0.330	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.256	88	991	0.414	ng/ul	88
5) Naphthalene	10.667	128	10987	0.400	ng/ul	100
7) 2-Methylnaphthalene	12.289	142	6978	0.380	ng/ul	99
8) 1-Methylnaphthalene	12.509	142	6279	0.360	ng/ul	100
10) Acenaphthylene	14.192	152	10290	0.402	ng/ul	99
11) Acenaphthene	14.534	153	8391	0.409	ng/ul	99
12) Fluorene	15.524	166	9529	0.392	ng/ul	99
14) Pentachlorophenol	16.888	266	791	0.338	ng/ul	99
15) Phenanthrene	17.263	178	16575	0.410	ng/ul	100
16) Anthracene	17.356	178	13537	0.388	ng/ul	99
19) Fluoranthene	19.290	202	19162	0.364	ng/ul	100
20) Pyrene	19.653	202	20008	0.367	ng/ul	99
21) Benzo(a)anthracene	21.407	228	14920	0.361	ng/ul	100
22) Chrysene	21.460	228	19403	0.422	ng/ul	100
24) Benzo(b)fluoranthene	23.056	252	18627	0.411	ng/ul	98
25) Benzo(k)fluoranthene	23.102	252	18852	0.412	ng/ul	98
26) Benzo(a)pyrene	23.667	252	16398	0.398	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	26.203	276	16536	0.404	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.216	278	13040	0.394	ng/ul	98
29) Benzo(g,h,i)perylene	26.938	276	16859	0.480	ng/ul	99

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

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