

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080422\
 Data File : BN021020.D
 Acq On : 05 Aug 2022 10:56
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4269

Quant Time: Aug 06 00:27:38 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 03 16:31:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.809	152	2957	0.400	ng/ul	0.00
4) Naphthalene-d8	10.602	136	9357	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.452	164	5169	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.209	188	9911	0.400	ng/ul	0.00
17) Chrysene-d12	21.412	240	8397	0.400	ng/ul	0.00
23) Perylene-d12	23.756	264	7124	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	1400	0.370	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.202	152	5515	0.402	ng/ul	0.00
18) Fluoranthene-d10	19.244	212	10599	0.409	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.244	88	1443	0.370	ng/ul	96
5) Naphthalene	10.651	128	9957	0.346	ng/ul	98
7) 2-Methylnaphthalene	12.273	142	6064	0.348	ng/ul	100
8) 1-Methylnaphthalene	12.493	142	6704	0.384	ng/ul	100
10) Acenaphthylene	14.174	152	8500	0.352	ng/ul	99
11) Acenaphthene	14.517	153	6863	0.343	ng/ul	99
12) Fluorene	15.511	166	7490	0.343	ng/ul	98
14) Pentachlorophenol	16.876	266	542	0.355	ng/ul	100
15) Phenanthrene	17.247	178	11892	0.335	ng/ul	100
16) Anthracene	17.345	178	9827	0.358	ng/ul	100
19) Fluoranthene	19.277	202	12839	0.345	ng/ul	99
20) Pyrene	19.640	202	13207	0.344	ng/ul	99
21) Benzo(a)anthracene	21.398	228	9577	0.350	ng/ul	99
22) Chrysene	21.450	228	11847	0.328	ng/ul	99
24) Benzo(b)fluoranthene	23.043	252	10662	0.323	ng/ul	99
25) Benzo(k)fluoranthene	23.090	252	11320	0.325	ng/ul	98
26) Benzo(a)pyrene	23.648	252	9987	0.332	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.172	276	10960	0.305	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.195	278	8733	0.308	ng/ul	98
29) Benzo(g,h,i)perylene	26.910	276	9643	0.301	ng/ul	99

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

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