

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080222\
 Data File : BN020972.D
 Acq On : 03 Aug 2022 10:45
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4265

Quant Time: Aug 03 13:43:00 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 04:19:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	2891	0.400	ng/ul	0.00
4) Naphthalene-d8	10.602	136	9263	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.457	164	5337	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.210	188	10446	0.400	ng/ul	0.00
17) Chrysene-d12	21.418	240	8990	0.400	ng/ul	0.00
23) Perylene-d12	23.759	264	8081	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	1374	0.390	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.202	152	5546	0.406	ng/ul	0.00
18) Fluoranthene-d10	19.245	212	11022	0.401	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.244	88	1415	0.373	ng/ul	93
5) Naphthalene	10.651	128	9980	0.394	ng/ul	98
7) 2-Methylnaphthalene	12.279	142	6200	0.402	ng/ul	99
8) 1-Methylnaphthalene	12.493	142	6805	0.400	ng/ul	100
10) Acenaphthylene	14.174	152	8658	0.397	ng/ul	99
11) Acenaphthene	14.517	153	7056	0.387	ng/ul	99
12) Fluorene	15.511	166	7719	0.368	ng/ul	98
14) Pentachlorophenol	16.876	266	554	0.332	ng/ul	100
15) Phenanthrene	17.248	178	12463	0.352	ng/ul	100
16) Anthracene	17.345	178	10107	0.332	ng/ul	99
19) Fluoranthene	19.277	202	13358	0.331	ng/ul	99
20) Pyrene	19.640	202	13618	0.322	ng/ul	98
21) Benzo(a)anthracene	21.404	228	10112	0.310	ng/ul	100
22) Chrysene	21.453	228	12590	0.312	ng/ul	100
24) Benzo(b)fluoranthene	23.049	252	11897	0.308	ng/ul	98
25) Benzo(k)fluoranthene	23.096	252	12859	0.308	ng/ul	98
26) Benzo(a)pyrene	23.657	252	11295	0.310	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.182	276	12999	0.304	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.196	278	10421	0.297	ng/ul	96
29) Benzo(g,h,i)perylene	26.920	276	11138	0.286	ng/ul	99

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

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