

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080422\
 Data File : BN021011.D
 Acq On : 05 Aug 2022 05:23
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4268

Quant Time: Aug 05 06:21:09 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	2672	0.400	ng/ul	0.00
4) Naphthalene-d8	10.602	136	8495	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.457	164	4744	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.210	188	9116	0.400	ng/ul	0.00
17) Chrysene-d12	21.412	240	7550	0.400	ng/ul	0.00
23) Perylene-d12	23.754	264	6253	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	1288	0.377	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.202	152	4958	0.398	ng/ul	0.00
18) Fluoranthene-d10	19.245	212	9571	0.410	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.244	88	1319	0.375	ng/ul	95
5) Naphthalene	10.651	128	9039	0.346	ng/ul	99
7) 2-Methylnaphthalene	12.279	142	5480	0.346	ng/ul	100
8) 1-Methylnaphthalene	12.493	142	6097	0.385	ng/ul	99
10) Acenaphthylene	14.174	152	7693	0.347	ng/ul	99
11) Acenaphthene	14.517	153	6286	0.343	ng/ul	98
12) Fluorene	15.511	166	6920	0.345	ng/ul	99
14) Pentachlorophenol	16.880	266	456	0.324	ng/ul	98
15) Phenanthrene	17.248	178	10980	0.337	ng/ul	99
16) Anthracene	17.345	178	8818	0.349	ng/ul	100
19) Fluoranthene	19.277	202	11734	0.351	ng/ul	99
20) Pyrene	19.640	202	11974	0.347	ng/ul	98
21) Benzo(a)anthracene	21.398	228	8460	0.343	ng/ul	99
22) Chrysene	21.450	228	10700	0.330	ng/ul	99
24) Benzo(b)fluoranthene	23.043	252	9535	0.329	ng/ul	100
25) Benzo(k)fluoranthene	23.093	252	10178	0.333	ng/ul	99
26) Benzo(a)pyrene	23.651	252	8776	0.332	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.172	276	9529	0.302	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.196	278	7563	0.304	ng/ul	98
29) Benzo(g,h,i)perylene	26.913	276	8649	0.307	ng/ul	99

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

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