

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN101819\
 Data File : BN008245.D
 Acq On : 18 Oct 2019 13:49
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.401

Quant Time: Oct 18 17:28:12 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 17:11:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	1932	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	8997	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	5803	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	12568	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	12727	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	14513	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	6126	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	17656	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.41	128	10563	0.412	ng/ul	99
5) 2-Methylnaphthalene	12.03	142	7198	0.410	ng/ul	99
7) Acenaphthylene	13.95	152	10796	0.391	ng/ul	100
8) Acenaphthene	14.29	153	8725	0.406	ng/ul	99
9) Fluorene	15.29	166	9694	0.397	ng/ul	99
11) Pentachlorophenol	16.66	266	1281	0.421	ng/ul	100
12) Phenanthrene	17.02	178	15617	0.424	ng/ul	100
13) Anthracene	17.11	178	14319	0.442	ng/ul	99
15) Fluoranthene	19.05	202	20060	0.413	ng/ul	97
17) Pyrene	19.41	202	21732	0.429	ng/ul	97
18) Benzo(a)anthracene	21.17	228	19451	0.420	ng/ul	99
19) Chrysene	21.22	228	22303	0.393	ng/ul	100
21) Benzo(b)fluoranthene	22.72	252	21132	0.447	ng/ul	96
22) Benzo(k)fluoranthene	22.76	252	22840	0.427	ng/ul	98
23) Benzo(a)pyrene	23.27	252	21601	0.434	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.53	276	25804	0.441	ng/ul	99
25) Dibenzo(a,h)anthracene	25.54	278	20310	0.439	ng/ul	97
26) Benzo(g,h,i)perylene	26.18	276	23230	0.433	ng/ul	98

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

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