

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101318\
 Data File : BN003163.D
 Acq On : 13 Oct 2018 05:28
 Operator : MJ/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.411

Quant Time: Oct 13 06:03:00 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 13 05:59:15 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	984	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	4936	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	2998	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	7089	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	5526	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	4619	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	3369	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	8762	0.43	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.47	128	5451	0.430	ng/ul	98
5) 2-Methylnaphthalene	12.11	142	3840	0.440	ng/ul	100
7) Acenaphthylene	14.00	152	5832	0.425	ng/ul	100
8) Acenaphthene	14.34	153	4604	0.442	ng/ul	99
9) Fluorene	15.35	166	5324	0.447	ng/ul	100
11) Pentachlorophenol	16.76	266	316	0.459	ng/ul	95
12) Phenanthrene	17.08	178	8697	0.440	ng/ul	100
13) Anthracene	17.18	178	7976	0.427	ng/ul	100
15) Fluoranthene	19.10	202	9998	0.431	ng/ul	99
17) Pyrene	19.47	202	10150	0.473	ng/ul	98
18) Benzo(a)anthracene	21.23	228	7680	0.432	ng/ul	99
19) Chrysene	21.28	228	8069	0.447	ng/ul	100
21) Benzo(b)fluoranthene	22.80	252	6938	0.417	ng/ul	98
22) Benzo(k)fluoranthene	22.84	252	7186	0.417	ng/ul	97
23) Benzo(a)pyrene	23.37	252	6578	0.441	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.69	276	7612	0.444	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.68	278	6130	0.447	ng/ul	98
26) Benzo(g,h,i)perylene	26.36	276	6420	0.444	ng/ul	99

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

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