

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101118\
 Data File : BN003097.D
 Acq On : 11 Oct 2018 10:59
 Operator : MJ/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.403

Quant Time: Oct 11 13:47:54 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 : Thu Oct 11 08:10:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	1128	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	5163	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	3141	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	7413	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	6150	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	4730	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	3475	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	9233	0.44	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	5667	0.428	ng/ul	100
5) 2-Methylnaphthalene	12.09	142	3958	0.434	ng/ul	100
7) Acenaphthylene	14.00	152	6180	0.430	ng/ul	100
8) Acenaphthene	14.33	153	4798	0.440	ng/ul	98
9) Fluorene	15.34	166	5521	0.442	ng/ul	99
11) Pentachlorophenol	16.75	266	298	0.414	ng/ul	92
12) Phenanthrene	17.07	178	8929	0.432	ng/ul	100
13) Anthracene	17.17	178	8425	0.432	ng/ul	100
15) Fluoranthene	19.09	202	10492	0.433	ng/ul	98
17) Pyrene	19.46	202	10691	0.448	ng/ul	99
18) Benzo(a)anthracene	21.22	228	8613	0.435	ng/ul	99
19) Chrysene	21.27	228	8785	0.437	ng/ul	100
21) Benzo(b)fluoranthene	22.79	252	7048	0.414	ng/ul	99
22) Benzo(k)fluoranthene	22.83	252	7585	0.430	ng/ul	99
23) Benzo(a)pyrene	23.36	252	6570	0.430	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.68	276	7556	0.431	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.67	278	6057	0.431	ng/ul	98
26) Benzo(g,h,i)perylene	26.35	276	6448	0.436	ng/ul	99

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

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