

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101018\
 Data File : BN003062.D
 Acq On : 10 Oct 2018 11:55
 Operator : MJ/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.499

Quant Time: Oct 10 12:41:33 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 : Wed Oct 10 08:36:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	1311	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	6183	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	3824	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	9489	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	9569	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	7891	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	4236	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	12520	0.46	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.46	128	6801	0.429	ng/ul	99
5) 2-Methylnaphthalene	12.08	142	4883	0.447	ng/ul	100
7) Acenaphthylene	13.99	152	7678	0.439	ng/ul	99
8) Acenaphthene	14.33	153	5806	0.437	ng/ul	98
9) Fluorene	15.33	166	6808	0.448	ng/ul	98
11) Pentachlorophenol	16.73	266	503	0.546	ng/ul	94
12) Phenanthrene	17.07	178	11437	0.432	ng/ul	100
13) Anthracene	17.17	178	11044	0.442	ng/ul	99
15) Fluoranthene	19.09	202	14261	0.460	ng/ul	99
17) Pyrene	19.46	202	14748	0.397	ng/ul	98
18) Benzo(a)anthracene	21.22	228	13734	0.446	ng/ul	99
19) Chrysene	21.27	228	13669	0.437	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	12422	0.437	ng/ul	95
22) Benzo(k)fluoranthene	22.83	252	12548	0.426	ng/ul	95
23) Benzo(a)pyrene	23.36	252	11046	0.434	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.67	276	11215	0.383	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.67	278	9038	0.386	ng/ul	94
26) Benzo(g,h,i)perylene	26.34	276	9261	0.375	ng/ul	97

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

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