

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101018\
 Data File : BN003029.D
 Acq On : 09 Oct 2018 16:35
 Operator : MJ/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.497

Quant Time: Oct 09 18:37:31 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 : Tue Oct 09 08:56:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	989	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	4738	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	2811	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	6709	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	6744	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	5578	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	3186	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	8900	0.46	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.47	128	5208	0.428	ng/ul	100
5) 2-Methylnaphthalene	12.09	142	3646	0.436	ng/ul	99
7) Acenaphthylene	14.00	152	5627	0.438	ng/ul	100
8) Acenaphthene	14.33	153	4278	0.438	ng/ul	99
9) Fluorene	15.34	166	4871	0.436	ng/ul	100
11) Pentachlorophenol	16.74	266	329	0.505	ng/ul	94
12) Phenanthrene	17.08	178	8101	0.433	ng/ul	100
13) Anthracene	17.17	178	7810	0.442	ng/ul	100
15) Fluoranthene	19.10	202	10106	0.461	ng/ul	99
17) Pyrene	19.46	202	10368	0.396	ng/ul	99
18) Benzo(a)anthracene	21.22	228	9544	0.440	ng/ul	99
19) Chrysene	21.27	228	9549	0.433	ng/ul	100
21) Benzo(b)fluoranthene	22.79	252	8672	0.432	ng/ul	88
22) Benzo(k)fluoranthene	22.83	252	8683	0.417	ng/ul#	87
23) Benzo(a)pyrene	23.36	252	7737	0.430	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.67	276	8108	0.392	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.67	278	6570	0.397	ng/ul#	88
26) Benzo(g,h,i)perylene	26.35	276	6802	0.390	ng/ul#	93

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

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