

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121118\
 Data File : BN003843.D
 Acq On : 12 Dec 2018 09:07
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.475

Quant Time: Dec 12 10:46:41 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 03:46:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	4532	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	18225	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	8797	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	17234	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	12560	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	12821	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	10075	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	16080	0.32	ng/ul	0.00

Target Compounds						Qvalue
3) Naphthalene	10.68	128	19603	0.383	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	13268	0.371	ng/ul	99
7) Acenaphthylene	14.19	152	17260	0.436	ng/ul	100
8) Acenaphthene	14.53	153	13342	0.384	ng/ul	100
9) Fluorene	15.52	166	14324	0.354	ng/ul	99
11) Pentachlorophenol	16.86	266	1189	0.383	ng/ul	96
12) Phenanthrene	17.25	178	21353	0.378	ng/ul	100
13) Anthracene	17.35	178	19454	0.419	ng/ul	99
15) Fluoranthene	19.27	202	21634	0.324	ng/ul	96
17) Pyrene	19.63	202	21604	0.480	ng/ul	98
18) Benzo(a)anthracene	21.38	228	17919	0.452	ng/ul	98
19) Chrysene	21.43	228	17361	0.382	ng/ul	99
21) Benzo(b)fluoranthene	23.01	252	18004	0.339	ng/ul	93
22) Benzo(k)fluoranthene	23.06	252	18370	0.356	ng/ul#	93
23) Benzo(a)pyrene	23.61	252	17585	0.383	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	26.06	276	21680	0.412	ng/ul	98
25) Dibenzo(a,h)anthracene	26.07	278	17186	0.405	ng/ul#	93
26) Benzo(g,h,i)perylene	26.79	276	18392	0.402	ng/ul	96

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

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