

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN112018\
 Data File : BN003647.D
 Acq On : 21 Nov 2018 10:05
 Operator : MJ/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.475

Quant Time: Nov 21 15:53:02 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 Nov 21 00:09:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	4244	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	16619	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	8936	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	22010	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	23331	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	20640	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	10144	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	25670	0.40	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.69	128	18850	0.404	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	13131	0.403	ng/ul	99
7) Acenaphthylene	14.20	152	16090	0.400	ng/ul	100
8) Acenaphthene	14.54	153	14206	0.403	ng/ul	100
9) Fluorene	15.53	166	16404	0.399	ng/ul	99
11) Pentachlorophenol	16.87	266	1489	0.376	ng/ul	98
12) Phenanthrene	17.26	178	28839	0.399	ng/ul	99
13) Anthracene	17.35	178	23567	0.397	ng/ul	100
15) Fluoranthene	19.27	202	34115	0.400	ng/ul	98
17) Pyrene	19.64	202	33542	0.401	ng/ul	99
18) Benzo(a)anthracene	21.38	228	28992	0.393	ng/ul	99
19) Chrysene	21.43	228	33550	0.398	ng/ul	99
21) Benzo(b)fluoranthene	23.01	252	34207	0.401	ng/ul	100
22) Benzo(k)fluoranthene	23.05	252	33331	0.401	ng/ul	99
23) Benzo(a)pyrene	23.61	252	29524	0.399	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.06	276	32398	0.382	ng/ul	99
25) Dibenzo(a,h)anthracene	26.08	278	25819	0.378	ng/ul	99
26) Benzo(g,h,i)perylene	26.79	276	27819	0.378	ng/ul	99

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

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