

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110218\
 Data File : BN003420.D
 Acq On : 03 Nov 2018 01:08
 Operator : MJ/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.436

Quant Time: Nov 03 00:41:45 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 02 23:49:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	11335	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	46765	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	25009	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	53139	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	38306	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	35411	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	29435	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	56123	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.70	128	54014	0.412	ng/ul	100
5) 2-Methylnaphthalene	12.31	142	38239	0.405	ng/ul	100
7) Acenaphthylene	14.21	152	52748	0.485	ng/ul	99
8) Acenaphthene	14.55	153	39403	0.406	ng/ul	99
9) Fluorene	15.54	166	44239	0.387	ng/ul	100
11) Pentachlorophenol	16.88	266	5477	0.515	ng/ul	100
12) Phenanthrene	17.27	178	68587	0.394	ng/ul	100
13) Anthracene	17.36	178	64246	0.452	ng/ul	99
15) Fluoranthene	19.29	202	71767	0.366	ng/ul	99
17) Pyrene	19.65	202	70347	0.505	ng/ul	100
18) Benzo(a)anthracene	21.39	228	57589	0.480	ng/ul	99
19) Chrysene	21.45	228	53551	0.392	ng/ul	99
21) Benzo(b)fluoranthene	23.02	252	53266	0.349	ng/ul	94
22) Benzo(k)fluoranthene	23.07	252	51996	0.352	ng/ul#	94
23) Benzo(a)pyrene	23.62	252	50194	0.394	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	26.08	276	61354	0.450	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.10	278	49652	0.447	ng/ul	95
26) Benzo(g,h,i)perylene	26.81	276	51782	0.443	ng/ul	96

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

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