

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111918\
 Data File : BN003622.D
 Acq On : 20 Nov 2018 07:31
 Operator : MJ/SJ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.470

Quant Time: Nov 20 08:29:06 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 : Tue Nov 20 03:27:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	5728	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	23069	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	12661	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	30380	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	33467	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	31230	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	14321	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	36279	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.69	128	25968	0.401	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	18457	0.408	ng/ul	99
7) Acenaphthylene	14.20	152	23705	0.416	ng/ul	100
8) Acenaphthene	14.54	153	20042	0.401	ng/ul	99
9) Fluorene	15.53	166	23129	0.397	ng/ul	100
11) Pentachlorophenol	16.87	266	2554	0.467	ng/ul	99
12) Phenanthrene	17.26	178	39825	0.399	ng/ul	100
13) Anthracene	17.35	178	34102	0.416	ng/ul	100
15) Fluoranthene	19.28	202	47585	0.404	ng/ul	99
17) Pyrene	19.64	202	47328	0.394	ng/ul	99
18) Benzo(a)anthracene	21.38	228	45100	0.427	ng/ul	99
19) Chrysene	21.43	228	47663	0.394	ng/ul	99
21) Benzo(b)fluoranthene	23.01	252	50950	0.394	ng/ul	99
22) Benzo(k)fluoranthene	23.05	252	49706	0.395	ng/ul	99
23) Benzo(a)pyrene	23.61	252	45217	0.404	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.06	276	48229	0.376	ng/ul	99
25) Dibenzo(a,h)anthracene	26.07	278	38992	0.378	ng/ul	99
26) Benzo(g,h,i)perylene	26.79	276	39949	0.358	ng/ul	100

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

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