

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110118\
 Data File : BN003395.D
 Acq On : 02 Nov 2018 00:51
 Operator : MJ/SJ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.433

Quant Time: Nov 02 04:53:55 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 04:49:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	9104	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	37541	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	22458	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	53611	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	52911	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	43957	0.40	ng/ul	-0.02

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	24927	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	63217	0.42	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.70	128	42619	0.405	ng/ul	100
5) 2-Methylnaphthalene	12.31	142	31577	0.417	ng/ul	99
7) Acenaphthylene	14.21	152	45549	0.467	ng/ul	100
8) Acenaphthene	14.55	153	35409	0.406	ng/ul	99
9) Fluorene	15.54	166	41402	0.403	ng/ul	99
11) Pentachlorophenol	16.87	266	5720	0.533	ng/ul	100
12) Phenanthrene	17.27	178	69171	0.393	ng/ul	99
13) Anthracene	17.36	178	63047	0.440	ng/ul	100
15) Fluoranthene	19.29	202	80739	0.409	ng/ul	99
17) Pyrene	19.65	202	81569	0.424	ng/ul	98
18) Benzo(a)anthracene	21.39	228	77104	0.465	ng/ul	99
19) Chrysene	21.45	228	73947	0.392	ng/ul	99
21) Benzo(b)fluoranthene	23.02	252	72019	0.380	ng/ul	98
22) Benzo(k)fluoranthene	23.07	252	68044	0.371	ng/ul	98
23) Benzo(a)pyrene	23.62	252	64454	0.408	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.08	276	65578	0.387	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.09	278	53297	0.386	ng/ul	98
26) Benzo(g,h,i)perylene	26.80	276	53922	0.371	ng/ul	99

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

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