

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110518\
 Data File : BN003435.D
 Acq On : 05 Nov 2018 18:21
 Operator : MJ/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.438

Quant Time: Nov 06 01:48:38 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 : Tue Nov 06 01:43:43 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	8727	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	36296	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	19816	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	43143	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	31823	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	28781	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	23375	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	46099	0.38	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.70	128	41454	0.407	ng/ul	100
5) 2-Methylnaphthalene	12.31	142	29684	0.405	ng/ul	100
7) Acenaphthylene	14.21	152	42583	0.495	ng/ul	99
8) Acenaphthene	14.55	153	31484	0.409	ng/ul	99
9) Fluorene	15.54	166	35355	0.390	ng/ul	99
11) Pentachlorophenol	16.88	266	4295	0.497	ng/ul	99
12) Phenanthrene	17.27	178	55813	0.394	ng/ul	99
13) Anthracene	17.36	178	52016	0.451	ng/ul	99
15) Fluoranthene	19.29	202	59363	0.373	ng/ul	100
17) Pyrene	19.65	202	57789	0.499	ng/ul	97
18) Benzo(a)anthracene	21.39	228	48064	0.482	ng/ul	99
19) Chrysene	21.44	228	45015	0.397	ng/ul	99
21) Benzo(b)fluoranthene	23.02	252	44126	0.356	ng/ul	96
22) Benzo(k)fluoranthene	23.06	252	42835	0.357	ng/ul#	96
23) Benzo(a)pyrene	23.62	252	41748	0.403	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	26.08	276	50176	0.453	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.09	278	40539	0.449	ng/ul	96
26) Benzo(g,h,i)perylene	26.80	276	42433	0.446	ng/ul	97

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

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