

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100918\
 Data File : BN002985.D
 Acq On : 08 Oct 2018 14:06
 Operator : MJ/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Oct 08 18:10:54 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 08 18:09:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	1156	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	5277	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	3058	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	7261	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	6569	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	4888	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	3539	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	9162	0.44	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.47	128	5777	0.427	ng/ul	99
5) 2-Methylnaphthalene	12.09	142	4053	0.435	ng/ul	99
7) Acenaphthylene	14.00	152	6126	0.438	ng/ul	100
8) Acenaphthene	14.34	153	4703	0.443	ng/ul	100
9) Fluorene	15.34	166	5346	0.440	ng/ul	99
11) Pentachlorophenol	16.75	266	246	0.349	ng/ul#	83
12) Phenanthrene	17.08	178	8804	0.435	ng/ul	100
13) Anthracene	17.17	178	8380	0.438	ng/ul	100
15) Fluoranthene	19.10	202	10491	0.442	ng/ul	99
17) Pyrene	19.46	202	10727	0.420	ng/ul	99
18) Benzo(a)anthracene	21.22	228	9207	0.436	ng/ul	100
19) Chrysene	21.27	228	9380	0.437	ng/ul	100
21) Benzo(b)fluoranthene	22.79	252	7715	0.438	ng/ul	99
22) Benzo(k)fluoranthene	22.84	252	7956	0.436	ng/ul	99
23) Benzo(a)pyrene	23.36	252	6933	0.439	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.67	276	7579	0.418	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.67	278	6080	0.419	ng/ul	99
26) Benzo(g,h,i)perylene	26.35	276	6324	0.414	ng/ul	100

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

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