

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110518\
 Data File : BN003422.D
 Acq On : 05 Nov 2018 10:38
 Operator : MJ/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.437

Quant Time: Nov 05 13:40:20 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	8788	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	36193	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	20173	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	43750	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	32908	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	29733	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	23493	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	47427	0.39	ng/ul	0.00

Target Compounds					Qvalue
3) Naphthalene	10.70	128	41265	0.406	ng/ul 100
5) 2-Methylnaphthalene	12.31	142	30111	0.412	ng/ul 100
7) Acenaphthylene	14.21	152	41758	0.476	ng/ul 99
8) Acenaphthene	14.55	153	32187	0.411	ng/ul 99
9) Fluorene	15.54	166	36405	0.395	ng/ul 99
11) Pentachlorophenol	16.88	266	3827	0.437	ng/ul 99
12) Phenanthrene	17.27	178	57414	0.400	ng/ul 99
13) Anthracene	17.36	178	51839	0.443	ng/ul 100
15) Fluoranthene	19.29	202	60924	0.378	ng/ul 100
17) Pyrene	19.65	202	59400	0.496	ng/ul 97
18) Benzo(a)anthracene	21.39	228	48918	0.474	ng/ul 99
19) Chrysene	21.44	228	46685	0.398	ng/ul 99
21) Benzo(b)fluoranthene	23.02	252	45394	0.354	ng/ul 96
22) Benzo(k)fluoranthene	23.06	252	44590	0.359	ng/ul 96
23) Benzo(a)pyrene	23.62	252	43114	0.403	ng/ul 96
24) Indeno(1,2,3-cd)pyrene	26.07	276	51063	0.446	ng/ul 100
25) Dibenzo(a,h)anthracene	26.09	278	40500	0.434	ng/ul 97
26) Benzo(g,h,i)perylene	26.80	276	43897	0.447	ng/ul 98

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

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