

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101618\
 Data File : BN003171.D
 Acq On : 16 Oct 2018 16:51
 Operator : MJ/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.413

Quant Time: Oct 17 01:28:54 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 : Wed Oct 17 01:25:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	814	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	4340	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.28	164	2705	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	6308	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	4597	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	4045	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	2980	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	7458	0.41	ng/ul	0.00

Target Compounds					Qvalue
3) Naphthalene	10.48	128	4808	0.432	ng/ul 98
5) 2-Methylnaphthalene	12.12	142	3340	0.436	ng/ul 97
7) Acenaphthylene	14.01	152	5193	0.420	ng/ul 99
8) Acenaphthene	14.34	153	4121	0.439	ng/ul 100
9) Fluorene	15.36	166	4777	0.444	ng/ul 97
11) Pentachlorophenol	16.77	266	229	0.374	ng/ul 97
12) Phenanthrene	17.09	178	7727	0.439	ng/ul 99
13) Anthracene	17.19	178	6877	0.414	ng/ul 99
15) Fluoranthene	19.10	202	8592	0.416	ng/ul 97
17) Pyrene	19.47	202	8786	0.492	ng/ul 99
18) Benzo(a)anthracene	21.23	228	6159	0.416	ng/ul 99
19) Chrysene	21.28	228	6798	0.453	ng/ul 99
21) Benzo(b)fluoranthene	22.80	252	5780	0.397	ng/ul 97
22) Benzo(k)fluoranthene	22.84	252	6513	0.431	ng/ul 97
23) Benzo(a)pyrene	23.38	252	5764	0.441	ng/ul 98
24) Indeno(1,2,3-cd)pyrene	25.70	276	6196	0.413	ng/ul# 98
25) Dibenzo(a,h)anthracene	25.70	278	4828	0.402	ng/ul 98
26) Benzo(g,h,i)perylene	26.38	276	5498	0.435	ng/ul 98

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

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