

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101618\
 Data File : BN003165.D
 Acq On : 16 Oct 2018 13:17
 Operator : MJ/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.412

Quant Time: Oct 17 01:25:21 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 : Sat Oct 13 05:59:15 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	790	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	3986	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	2532	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	6080	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	4521	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	4162	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	2747	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	7399	0.43	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.48	128	4483	0.438	ng/ul	96
5) 2-Methylnaphthalene	12.11	142	3123	0.444	ng/ul	98
7) Acenaphthylene	14.00	152	4905	0.423	ng/ul	99
8) Acenaphthene	14.34	153	3920	0.446	ng/ul	99
9) Fluorene	15.35	166	4569	0.454	ng/ul	99
11) Pentachlorophenol	16.77	266	213	0.361	ng/ul	87
12) Phenanthrene	17.08	178	7498	0.442	ng/ul	99
13) Anthracene	17.18	178	6725	0.420	ng/ul	99
15) Fluoranthene	19.10	202	8518	0.428	ng/ul	99
17) Pyrene	19.47	202	8804	0.501	ng/ul	100
18) Benzo(a)anthracene	21.23	228	6216	0.427	ng/ul	99
19) Chrysene	21.28	228	6830	0.462	ng/ul	99
21) Benzo(b)fluoranthene	22.80	252	5902	0.394	ng/ul	98
22) Benzo(k)fluoranthene	22.84	252	6617	0.426	ng/ul	98
23) Benzo(a)pyrene	23.37	252	5903	0.439	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.69	276	6123	0.397	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.69	278	4608	0.373	ng/ul	99
26) Benzo(g,h,i)perylene	26.37	276	5567	0.428	ng/ul	100

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

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