

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101118\
 Data File : BN003064.D
 Acq On : 10 Oct 2018 14:45
 Operator : MJ/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.401

Quant Time: Oct 11 01:20:10 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 10 08:36:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	1308	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	6069	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	3572	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	8073	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	5662	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	4991	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	3724	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	8433	0.37	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.46	128	6092	0.391	ng/ul	99
5) 2-Methylnaphthalene	12.09	142	4293	0.400	ng/ul	100
7) Acenaphthylene	13.99	152	6443	0.394	ng/ul	100
8) Acenaphthene	14.33	153	4971	0.401	ng/ul	100
9) Fluorene	15.34	166	5621	0.396	ng/ul	99
11) Pentachlorophenol	16.74	266	222	0.283	ng/ul	93
12) Phenanthrene	17.07	178	8834	0.393	ng/ul	100
13) Anthracene	17.17	178	8216	0.387	ng/ul	100
15) Fluoranthene	19.09	202	9555	0.362	ng/ul	99
17) Pyrene	19.46	202	9513	0.433	ng/ul	99
18) Benzo(a)anthracene	21.22	228	7198	0.395	ng/ul	99
19) Chrysene	21.28	228	7355	0.398	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	6639	0.369	ng/ul	98
22) Benzo(k)fluoranthene	22.84	252	6843	0.367	ng/ul	98
23) Benzo(a)pyrene	23.36	252	6203	0.385	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.67	276	7596	0.410	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.67	278	6112	0.412	ng/ul	97
26) Benzo(g,h,i)perylene	26.35	276	6489	0.416	ng/ul	99

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

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