

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061019\
 Data File : BN006239.D
 Acq On : 11 Jun 2019 00:38
 Operator : JU/SJ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.408

Quant Time: Jun 11 01:40:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jun 09 05:10:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	5251	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	21705	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	10787	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	19542	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	12055	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	13666	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	13170	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	19179	0.37	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.47	128	24936	0.416	ng/ul	100
5) 2-Methylnaphthalene	12.10	142	16463	0.394	ng/ul	98
7) Acenaphthylene	14.01	152	23724	0.396	ng/ul	100
8) Acenaphthene	14.36	153	17276	0.415	ng/ul	99
9) Fluorene	15.35	166	17845	0.367	ng/ul	97
11) Pentachlorophenol	16.71	266	1535	0.309	ng/ul	99
12) Phenanthrene	17.09	178	24805	0.395	ng/ul	99
13) Anthracene	17.18	178	21746	0.363	ng/ul	99
15) Fluoranthene	19.12	202	25188	0.364	ng/ul	97
17) Pyrene	19.49	202	25915	0.408	ng/ul	99
18) Benzo(a)anthracene	21.25	228	19775	0.371	ng/ul	99
19) Chrysene	21.30	228	23481	0.469	ng/ul	99
21) Benzo(b)fluoranthene	22.82	252	21330	0.374	ng/ul	96
22) Benzo(k)fluoranthene	22.87	252	23115	0.416	ng/ul	97
23) Benzo(a)pyrene	23.39	252	21891	0.412	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.73	276	25662	0.447	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.73	278	20066	0.437	ng/ul	96
26) Benzo(g,h,i)perylene	26.41	276	22055	0.458	ng/ul	97

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

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