

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101218\
 Data File : BN003122.D
 Acq On : 12 Oct 2018 02:28
 Operator : MJ/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.406

Quant Time: Oct 12 05:38:50 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 : Fri Oct 12 05:29:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	1428	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	5807	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	2642	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	5085	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	4813	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	5152	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	3522	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	6232	0.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	6356	0.427	ng/ul	100
5) 2-Methylnaphthalene	12.09	142	4014	0.391	ng/ul	100
7) Acenaphthylene	14.00	152	5469	0.452	ng/ul	100
8) Acenaphthene	14.33	153	3995	0.436	ng/ul	99
9) Fluorene	15.34	166	4178	0.398	ng/ul	99
11) Pentachlorophenol	16.74	266	288	0.584	ng/ul	87
12) Phenanthrene	17.07	178	6093	0.430	ng/ul	99
13) Anthracene	17.17	178	5783	0.432	ng/ul	99
15) Fluoranthene	19.09	202	7059	0.424	ng/ul	97
17) Pyrene	19.46	202	7327	0.392	ng/ul	100
18) Benzo(a)anthracene	21.22	228	6816	0.440	ng/ul	98
19) Chrysene	21.27	228	6879	0.437	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	7278	0.392	ng/ul	99
22) Benzo(k)fluoranthene	22.83	252	7573	0.394	ng/ul	100
23) Benzo(a)pyrene	23.35	252	7138	0.429	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.66	276	8714	0.456	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.66	278	6976	0.456	ng/ul	97
26) Benzo(g,h,i)perylene	26.34	276	7366	0.457	ng/ul	99

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

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