

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082318\
 Data File : BN002461.D
 Acq On : 23 Aug 2018 12:26
 Operator : JU/SJ
 Sample : SSTD0.811
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.811

Quant Time: Aug 23 13:13:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 23 13:07:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	1492	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	6086	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	3506	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	8525	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	9217	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	8292	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	7632	0.79	ng/ul	-0.01
14) Fluoranthene-d10	19.09	212	20249	0.78	ng/ul	0.00

Target Compounds						Qvalue
3) Naphthalene	10.50	128	12201	0.778	ng/ul	97
5) 2-Methylnaphthalene	12.12	142	8708	0.777	ng/ul	99
7) Acenaphthylene	14.03	152	13011	0.759	ng/ul	96
8) Acenaphthene	14.37	153	9899	0.785	ng/ul	94
9) Fluorene	15.36	166	11530	0.783	ng/ul#	93
11) Pentachlorophenol	16.73	266	926	0.574	ng/ul	97
12) Phenanthrene	17.10	178	19908	0.783	ng/ul#	89
13) Anthracene	17.20	178	17641	0.732	ng/ul#	92
15) Fluoranthene	19.12	202	24138	0.799	ng/ul	97
17) Pyrene	19.49	202	23198	0.692	ng/ul#	95
18) Benzo(a)anthracene	21.24	228	21714	0.681	ng/ul#	85
19) Chrysene	21.29	228	24070	0.759	ng/ul	97
21) Benzo(b)fluoranthene	22.81	252	23504	0.781	ng/ul	84
22) Benzo(k)fluoranthene	22.86	252	24615	0.820	ng/ul#	85
23) Benzo(a)pyrene	23.39	252	22359	0.800	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.70	276	24264	0.773	ng/ul#	85
25) Dibenzo(a,h)anthracene	25.71	278	19157	0.766	ng/ul#	89
26) Benzo(g,h,i)perylene	26.38	276	20439	0.766	ng/ul	94

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

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