

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN062719\
 Data File : BN006596.D
 Acq On : 27 Jun 2019 11:11
 Operator : HP/JU
 Sample : SSTD0.113
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.113

Quant Time: Jun 27 14:38:46 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 27 14:04:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2110	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	8704	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	5280	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	11942	0.40	ng/ul	0.00
16) Chrysene-d12	21.28	240	11119	0.40	ng/ul	0.01
20) Perylene-d12	23.54	264	12302	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	1299	0.10	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	3452	0.11	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.46	128	2509	0.104	ng/ul#	90
5) 2-Methylnaphthalene	12.09	142	1637	0.098	ng/ul	99
7) Acenaphthylene	14.00	152	2608	0.089	ng/ul	94
8) Acenaphthene	14.34	153	2038	0.100	ng/ul	98
9) Fluorene	15.34	166	2175	0.091	ng/ul#	98
12) Phenanthrene	17.08	178	3720	0.097	ng/ul	96
13) Anthracene	17.18	178	3038	0.083	ng/ul#	94
15) Fluoranthene	19.11	202	4751	0.112	ng/ul	98
17) Pyrene	19.48	202	5142	0.088	ng/ul	99
18) Benzo(a)anthracene	21.27	228	4511	0.092	ng/ul	96
19) Chrysene	21.32	228	5273	0.114	ng/ul	96
21) Benzo(b)fluoranthene	22.88	252	4832	0.094	ng/ul	82
22) Benzo(k)fluoranthene	22.92	252	5189	0.104	ng/ul#	82
23) Benzo(a)pyrene	23.45	252	4952	0.104	ng/ul#	74
24) Indeno(1,2,3-cd)pyrene	25.79	276	4902	0.095	ng/ul	98
25) Dibenzo(a,h)anthracene	25.80	278	3890	0.094	ng/ul#	85
26) Benzo(a,h,i)perylene	26.48	276	4141	0.096	ng/ul#	89

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

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