

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101019\
 Data File : BN008059.D
 Acq On : 10 Oct 2019 12:53
 Operator : JU
 Sample : SSTD0.202
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.202

Quant Time: Oct 10 16:14:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 10 16:00:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	2514	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	11630	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	7533	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	17802	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	18972	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	23055	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	3811	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	12124	0.21	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	6634	0.202	ng/ul	98
5) 2-Methylnaphthalene	12.08	142	4390	0.199	ng/ul	100
7) Acenaphthylene	13.99	152	7068	0.169	ng/ul	99
8) Acenaphthene	14.33	153	5568	0.185	ng/ul	99
9) Fluorene	15.32	166	6182	0.190	ng/ul	99
11) Pentachlorophenol	16.68	266	796	0.198	ng/ul	98
12) Phenanthrene	17.06	178	10090	0.182	ng/ul	100
13) Anthracene	17.15	178	8487	0.162	ng/ul	98
15) Fluoranthene	19.08	202	13543	0.182	ng/ul	99
17) Pyrene	19.44	202	14779	0.159	ng/ul	100
18) Benzo(a)anthracene	21.20	228	13005	0.153	ng/ul	99
19) Chrysene	21.25	228	16984	0.186	ng/ul	100
21) Benzo(b)fluoranthene	22.75	252	14097	0.140	ng/ul	91
22) Benzo(k)fluoranthene	22.80	252	16743	0.159	ng/ul#	93
23) Benzo(a)pyrene	23.31	252	15291	0.152	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	25.59	276	17845	0.159	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.60	278	13965	0.157	ng/ul	93
26) Benzo(g,h,i)perylene	26.25	276	16547	0.172	ng/ul	98

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

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