

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101119\
 Data File : BN008122.D
 Acq On : 12 Oct 2019 20:43
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.413

Quant Time: Oct 12 21:41:30 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 : Fri Oct 11 19:05:54 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2303	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	12029	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	7914	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	19358	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	19224	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	16569	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	7751	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	24095	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	12483	0.365	ng/ul	100
5) 2-Methylnaphthalene	12.07	142	8884	0.378	ng/ul	99
7) Acenaphthylene	13.98	152	14351	0.381	ng/ul	100
8) Acenaphthene	14.32	153	11800	0.403	ng/ul	99
9) Fluorene	15.31	166	12135	0.365	ng/ul	98
11) Pentachlorophenol	16.67	266	1629	0.348	ng/ul	99
12) Phenanthrene	17.05	178	20030	0.353	ng/ul	99
13) Anthracene	17.14	178	18579	0.373	ng/ul	99
15) Fluoranthene	19.07	202	26175	0.350	ng/ul	100
17) Pyrene	19.44	202	30399	0.397	ng/ul	99
18) Benzo(a)anthracene	21.19	228	25861	0.370	ng/ul	98
19) Chrysene	21.24	228	29143	0.340	ng/ul	100
21) Benzo(b)fluoranthene	22.74	252	23050	0.427	ng/ul	97
22) Benzo(k)fluoranthene	22.78	252	25769	0.422	ng/ul	99
23) Benzo(a)pyrene	23.30	252	22722	0.400	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.57	276	22162	0.332	ng/ul	100
25) Dibenzo(a,h)anthracene	25.58	278	17949	0.340	ng/ul	95
26) Benzo(g,h,i)perylene	26.23	276	19002	0.310	ng/ul	96

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

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