

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101819\
 Data File : BN008259.D
 Acq On : 18 Oct 2019 22:10
 Operator : JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.402

Quant Time: Oct 18 23:51:24 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 18 23:47:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	2188	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	10790	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	6797	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	14209	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	13520	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	12461	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	7320	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	18732	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.41	128	12716	0.414	ng/ul	99
5) 2-Methylnaphthalene	12.03	142	8633	0.410	ng/ul	100
7) Acenaphthylene	13.95	152	12663	0.391	ng/ul	100
8) Acenaphthene	14.29	153	10345	0.411	ng/ul	99
9) Fluorene	15.29	166	11247	0.393	ng/ul	98
11) Pentachlorophenol	16.66	266	1155	0.336	ng/ul	99
12) Phenanthrene	17.02	178	17871	0.430	ng/ul	100
13) Anthracene	17.11	178	16165	0.442	ng/ul	100
15) Fluoranthene	19.04	202	21542	0.392	ng/ul	95
17) Pyrene	19.41	202	23066	0.428	ng/ul	99
18) Benzo(a)anthracene	21.17	228	20887	0.425	ng/ul	100
19) Chrysene	21.22	228	23315	0.387	ng/ul	100
21) Benzo(b)fluoranthene	22.71	252	20089	0.495	ng/ul	97
22) Benzo(k)fluoranthene	22.75	252	22179	0.483	ng/ul	97
23) Benzo(a)pyrene	23.26	252	19653	0.460	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.52	276	19059	0.379	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.53	278	15396	0.387	ng/ul	98
26) Benzo(g,h,i)perylene	26.17	276	16276	0.353	ng/ul	99

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

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