

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101419\
 Data File : BN008165.D
 Acq On : 15 Oct 2019 10:07
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.417

Quant Time: Oct 15 11:03:48 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 : Tue Oct 15 10:50:49 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1988	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	10351	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	6863	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	16474	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	15313	0.40	ng/ul	0.00
20) Perylene-d12	23.38	264	13192	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	6584	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	19333	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.43	128	10840	0.368	ng/ul	99
5) 2-Methylnaphthalene	12.06	142	7659	0.379	ng/ul	99
7) Acenaphthylene	13.97	152	11796	0.361	ng/ul	99
8) Acenaphthene	14.32	153	10319	0.406	ng/ul	99
9) Fluorene	15.31	166	10635	0.368	ng/ul	99
11) Pentachlorophenol	16.67	266	1240	0.311	ng/ul	100
12) Phenanthrene	17.04	178	17368	0.360	ng/ul	99
13) Anthracene	17.14	178	15396	0.363	ng/ul	100
15) Fluoranthene	19.06	202	21797	0.342	ng/ul	98
17) Pyrene	19.43	202	25202	0.413	ng/ul	99
18) Benzo(a)anthracene	21.18	228	20232	0.363	ng/ul	100
19) Chrysene	21.24	228	23952	0.351	ng/ul	99
21) Benzo(b)fluoranthene	22.73	252	18493	0.431	ng/ul	96
22) Benzo(k)fluoranthene	22.77	252	20080	0.413	ng/ul	99
23) Benzo(a)pyrene	23.29	252	17547	0.388	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.55	276	17745	0.333	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.56	278	14138	0.336	ng/ul	99
26) Benzo(g,h,i)perylene	26.21	276	15770	0.323	ng/ul	98

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

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