

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120619\
 Data File : BN008854.D
 Acq On : 06 Dec 2019 11:53
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.401

Quant Time: Dec 06 12:30:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 06 12:03:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	2072	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	8440	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	5475	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	12326	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	11530	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	12611	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	5544	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	14869	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	9928	0.390	ng/ul	99
5) 2-Methylnaphthalene	12.00	142	7176	0.404	ng/ul	100
7) Acenaphthylene	13.91	152	10660	0.352	ng/ul	100
8) Acenaphthene	14.26	153	8165	0.363	ng/ul	99
9) Fluorene	15.25	166	9737	0.369	ng/ul	100
11) Pentachlorophenol	16.60	266	1343	0.288	ng/ul	98
12) Phenanthrene	16.98	178	16934	0.407	ng/ul	99
13) Anthracene	17.07	178	15390	0.392	ng/ul	99
15) Fluoranthene	19.00	202	20318	0.398	ng/ul	99
17) Pyrene	19.37	202	20509	0.443	ng/ul	99
18) Benzo(a)anthracene	21.13	228	20193	0.435	ng/ul	99
19) Chrysene	21.18	228	20073	0.441	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	20621	0.397	ng/ul	99
22) Benzo(k)fluoranthene	22.70	252	20893	0.405	ng/ul	98
23) Benzo(a)pyrene	23.20	252	20196	0.414	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.41	276	24491	0.429	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.42	278	19720	0.428	ng/ul	99
26) Benzo(g,h,i)perylene	26.05	276	21026	0.440	ng/ul	98

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

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