

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN011020\
 Data File : BN009357.D
 Acq On : 11 Jan 2020 05:12
 Operator : JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.417

Quant Time: Jan 11 05:43:12 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN011020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 10 16:45:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	1235	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	4275	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	2527	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	5938	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	5573	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	5393	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	2801	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	7760	0.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.28	128	5181	0.403	ng/ul	99
5) 2-Methylnaphthalene	11.91	142	3563	0.403	ng/ul	100
7) Acenaphthylene	13.82	152	4806	0.373	ng/ul	98
8) Acenaphthene	14.17	153	3971	0.383	ng/ul	99
9) Fluorene	15.17	166	4509	0.378	ng/ul	99
11) Pentachlorophenol	16.54	266	631	0.388	ng/ul	99
12) Phenanthrene	16.90	178	8186	0.399	ng/ul	99
13) Anthracene	17.00	178	7125	0.392	ng/ul	99
15) Fluoranthene	18.93	202	10343	0.426	ng/ul	98
17) Pyrene	19.29	202	10604	0.396	ng/ul	99
18) Benzo(a)anthracene	21.05	228	9709	0.398	ng/ul	100
19) Chrysene	21.10	228	10565	0.400	ng/ul	98
21) Benzo(b)fluoranthene	22.56	252	9537	0.399	ng/ul	97
22) Benzo(k)fluoranthene	22.60	252	11113	0.408	ng/ul	97
23) Benzo(a)pyrene	23.09	252	9178	0.383	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.26	276	10695	0.400	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.27	278	8475	0.402	ng/ul	99
26) Benzo(g,h,i)perylene	25.89	276	9464	0.404	ng/ul	98

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

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