

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN011320\
 Data File : BN009361.D
 Acq On : 13 Jan 2020 13:02
 Operator : JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.419

Quant Time: Jan 13 13:44:23 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	7.47	152	1079	0.40	ng/ul	0.00
2) Naphthalene-d8	10.22	136	3717	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.11	164	2173	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	4598	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	3735	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	3560	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	2395	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	5475	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.27	128	4488	0.402	ng/ul	99
5) 2-Methylnaphthalene	11.91	142	3043	0.396	ng/ul	99
7) Acenaphthylene	13.82	152	4060	0.366	ng/ul	98
8) Acenaphthene	14.17	153	3365	0.377	ng/ul	99
9) Fluorene	15.17	166	3817	0.372	ng/ul	99
11) Pentachlorophenol	16.54	266	414	0.329	ng/ul	98
12) Phenanthrene	16.90	178	6380	0.402	ng/ul	99
13) Anthracene	17.00	178	5325	0.378	ng/ul	100
15) Fluoranthene	18.93	202	7148	0.380	ng/ul	99
17) Pyrene	19.29	202	7278	0.406	ng/ul	97
18) Benzo(a)anthracene	21.05	228	5997	0.366	ng/ul	99
19) Chrysene	21.10	228	7437	0.420	ng/ul	99
21) Benzo(b)fluoranthene	22.57	252	6635	0.420	ng/ul	98
22) Benzo(k)fluoranthene	22.61	252	7503	0.417	ng/ul	97
23) Benzo(a)pyrene	23.10	252	5986	0.379	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.27	276	6999	0.397	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.28	278	5526	0.397	ng/ul	96
26) Benzo(g,h,i)perylene	25.89	276	6300	0.408	ng/ul	99

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

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