

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN011020\
 Data File : BN009351.D
 Acq On : 11 Jan 2020 01:02
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.416

Quant Time: Jan 11 01:40:48 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	1077	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	3800	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	2238	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.87	188	4808	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	4094	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	3897	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.84	152	2447	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	5729	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.28	128	4583	0.401	ng/ul	99
5) 2-Methylnaphthalene	11.91	142	3122	0.398	ng/ul	100
7) Acenaphthylene	13.83	152	4243	0.372	ng/ul	98
8) Acenaphthene	14.18	153	3475	0.378	ng/ul	97
9) Fluorene	15.17	166	3976	0.377	ng/ul	100
11) Pentachlorophenol	16.54	266	452	0.343	ng/ul	99
12) Phenanthrene	16.90	178	6526	0.393	ng/ul	99
13) Anthracene	17.00	178	5643	0.383	ng/ul	100
15) Fluoranthene	18.93	202	7511	0.382	ng/ul	99
17) Pyrene	19.29	202	7744	0.394	ng/ul	98
18) Benzo(a)anthracene	21.05	228	6763	0.377	ng/ul	99
19) Chrysene	21.10	228	7809	0.402	ng/ul	97
21) Benzo(b)fluoranthene	22.56	252	7110	0.411	ng/ul	97
22) Benzo(k)fluoranthene	22.60	252	7836	0.398	ng/ul	97
23) Benzo(a)pyrene	23.10	252	6633	0.383	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.26	276	7549	0.391	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.27	278	5943	0.390	ng/ul	96
26) Benzo(g,h,i)perylene	25.89	276	6766	0.400	ng/ul	99

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

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