

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN011720\
 Data File : BN009390.D
 Acq On : 17 Jan 2020 11:40
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.420

Quant Time: Jan 17 12:30:25 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 17 12:28:30 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	1291	0.40	ng/ul	0.00
2) Naphthalene-d8	10.22	136	4236	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	2420	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	5125	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	3856	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	3649	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	2747	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	5879	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.27	128	5156	0.405	ng/ul	99
5) 2-Methylnaphthalene	11.90	142	3453	0.395	ng/ul	99
7) Acenaphthylene	13.82	152	4404	0.357	ng/ul	98
8) Acenaphthene	14.17	153	3773	0.380	ng/ul	99
9) Fluorene	15.17	166	4252	0.373	ng/ul	100
11) Pentachlorophenol	16.54	266	382	0.272	ng/ul	97
12) Phenanthrene	16.90	178	7032	0.397	ng/ul	99
13) Anthracene	17.00	178	5840	0.372	ng/ul	98
15) Fluoranthene	18.93	202	7723	0.368	ng/ul	100
17) Pyrene	19.29	202	7841	0.423	ng/ul	98
18) Benzo(a)anthracene	21.05	228	6129	0.363	ng/ul	99
19) Chrysene	21.10	228	7703	0.421	ng/ul	100
21) Benzo(b)fluoranthene	22.56	252	6689	0.413	ng/ul	98
22) Benzo(k)fluoranthene	22.60	252	7748	0.420	ng/ul	96
23) Benzo(a)pyrene	23.10	252	6331	0.391	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.26	276	7205	0.399	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.27	278	5696	0.399	ng/ul	96
26) Benzo(g,h,i)perylene	25.88	276	6547	0.413	ng/ul	99

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

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