

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101920\
 Data File : BN012343.D
 Acq On : 19 Oct 2020 10:02
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.423

Quant Time: Oct 19 11:52:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 19 11:52:04 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	1178	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	5167	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	2967	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	5992	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	5262	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	4967	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	3159	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	5810	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	5833	0.391	ng/ul	100
5) 2-Methylnaphthalene	12.04	142	3910	0.413	ng/ul	100
7) Acenaphthylene	13.96	152	5220	0.406	ng/ul	100
8) Acenaphthene	14.30	153	4380	0.403	ng/ul	100
9) Fluorene	15.30	166	4842	0.400	ng/ul	100
11) Pentachlorophenol	16.65	266	654	0.378	ng/ul	100
12) Phenanthrene	17.03	178	7453	0.390	ng/ul	100
13) Anthracene	17.12	178	6562	0.401	ng/ul	100
15) Fluoranthene	19.05	202	8174	0.366	ng/ul	100
17) Pyrene	19.42	202	8347	0.352	ng/ul	100
18) Benzo(a)anthracene	21.17	228	7000	0.359	ng/ul	100
19) Chrysene	21.22	228	8584	0.361	ng/ul	100
21) Benzo(b)fluoranthene	22.71	252	8320	0.415	ng/ul	100
22) Benzo(k)fluoranthene	22.76	252	8774	0.395	ng/ul	100
23) Benzo(a)pyrene	23.27	252	7741	0.397	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.53	276	9770	0.390	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.54	278	7921	0.395	ng/ul	100
26) Benzo(g,h,i)perylene	26.20	276	8677	0.378	ng/ul	100

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

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