

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN031620\
 Data File : BN010228.D
 Acq On : 16 Mar 2020 15:02
 Operator : CG/JU
 Sample : SSTD3.203
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.203

Quant Time: Mar 16 15:32:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN031620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 16 13:58:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1858	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	6101	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	3631	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	8271	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	8309	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	7627	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	32759	3.30	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	85953	3.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	59968	3.269	ng/ul	98
5) 2-Methylnaphthalene	12.06	142	41363	3.282	ng/ul	100
7) Acenaphthylene	13.97	152	64450	3.479	ng/ul	99
8) Acenaphthene	14.32	153	47334	3.177	ng/ul	99
9) Fluorene	15.31	166	57637	3.366	ng/ul	99
11) Pentachlorophenol	16.65	266	6098	2.693	ng/ul	97
12) Phenanthrene	17.04	178	96320	3.371	ng/ul	98
13) Anthracene	17.13	178	90518	3.573	ng/ul	99
15) Fluoranthene	19.06	202	121386	3.587	ng/ul	99
17) Pyrene	19.42	202	119549	2.996	ng/ul	98
18) Benzo(a)anthracene	21.18	228	115870	3.182	ng/ul	100
19) Chrysene	21.23	228	118796	3.014	ng/ul	100
21) Benzo(b)fluoranthene	22.72	252	113361	3.352	ng/ul	96
22) Benzo(k)fluoranthene	22.77	252	118911	3.085	ng/ul	97
23) Benzo(a)pyrene	23.28	252	99628	2.942	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.52	276	127587	3.378	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.54	278	104792	3.512	ng/ul	97
26) Benzo(g,h,i)perylene	26.17	276	108368	3.274	ng/ul	98

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

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