

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN060320\
 Data File : BN010752.D
 Acq On : 03 Jun 2020 14:57
 Operator : CG/JU
 Sample : SST0.803
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.803

Quant Time: Jun 03 15:41:16 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN060320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 03 14:56:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1597	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	5188	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	2663	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	5328	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	4671	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	4265	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	9768	1.16	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	18495	1.19	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.37	128	12048	0.753	ng/ul	98
5) 2-Methylnaphthalene	11.99	142	7700	0.749	ng/ul	99
7) Acenaphthylene	13.90	152	9429	0.738	ng/ul	99
8) Acenaphthene	14.25	153	7892	0.753	ng/ul	99
9) Fluorene	15.25	166	8760	0.757	ng/ul	99
11) Pentachlorophenol	16.60	266	1021	0.719	ng/ul	96
12) Phenanthrene	16.98	178	13780	0.753	ng/ul	99
13) Anthracene	17.07	178	11580	0.760	ng/ul	98
15) Fluoranthene	19.00	202	15914	0.787	ng/ul	98
17) Pyrene	19.37	202	15632	0.767	ng/ul	99
18) Benzo(a)anthracene	21.13	228	13081	0.739	ng/ul	98
19) Chrysene	21.18	228	16101	0.752	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	14960	0.774	ng/ul	93
22) Benzo(k)fluoranthene	22.71	252	16000	0.774	ng/ul	93
23) Benzo(a)pyrene	23.21	252	12357	0.744	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.42	276	17413	0.763	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.44	278	14374	0.773	ng/ul	95
26) Benzo(g,h,i)perylene	26.06	276	15898	0.788	ng/ul	96

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

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