

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060420\
 Data File : BN010772.D
 Acq On : 04 Jun 2020 21:01
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.413

Quant Time: Jun 04 21:40:38 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 : Thu Jun 04 12:51:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1550	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	4974	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	2579	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	5098	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	4169	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	3785	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	4963	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	9090	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	6189	0.422	ng/ul	99
5) 2-Methylnaphthalene	11.99	142	3903	0.424	ng/ul	98
7) Acenaphthylene	13.90	152	4809	0.418	ng/ul	100
8) Acenaphthene	14.25	153	4045	0.420	ng/ul	98
9) Fluorene	15.25	166	4469	0.422	ng/ul	100
11) Pentachlorophenol	16.60	266	502	0.421	ng/ul	99
12) Phenanthrene	16.98	178	6980	0.421	ng/ul	100
13) Anthracene	17.07	178	5632	0.411	ng/ul	100
15) Fluoranthene	19.00	202	7720	0.415	ng/ul	98
17) Pyrene	19.37	202	7733	0.427	ng/ul	99
18) Benzo(a)anthracene	21.13	228	6057	0.409	ng/ul	99
19) Chrysene	21.18	228	8021	0.426	ng/ul	98
21) Benzo(b)fluoranthene	22.66	252	6964	0.413	ng/ul	99
22) Benzo(k)fluoranthene	22.70	252	8149	0.457	ng/ul	99
23) Benzo(a)pyrene	23.21	252	6146	0.432	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.42	276	8467	0.434	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.43	278	6899	0.433	ng/ul	100
26) Benzo(g,h,i)perylene	26.06	276	7520	0.417	ng/ul	99

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

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