

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060520\
 Data File : BN010783.D
 Acq On : 05 Jun 2020 16:14
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.415

Quant Time: Jun 05 16:48:54 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 16:40:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1456	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	4715	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	2429	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	4842	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	3867	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	3528	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	4683	0.42	ng/ul	-0.01
14) Fluoranthene-d10	18.97	212	8639	0.41	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.36	128	5841	0.420	ng/ul	100
5) 2-Methylnaphthalene	11.98	142	3721	0.427	ng/ul	98
7) Acenaphthylene	13.90	152	4545	0.420	ng/ul	98
8) Acenaphthene	14.25	153	3809	0.420	ng/ul	99
9) Fluorene	15.24	166	4216	0.423	ng/ul	98
11) Pentachlorophenol	16.60	266	490	0.432	ng/ul	98
12) Phenanthrene	16.97	178	6637	0.421	ng/ul	100
13) Anthracene	17.07	178	5326	0.409	ng/ul	99
15) Fluoranthene	19.00	202	7265	0.411	ng/ul	99
17) Pyrene	19.37	202	7174	0.427	ng/ul	99
18) Benzo(a)anthracene	21.13	228	5674	0.413	ng/ul	98
19) Chrysene	21.18	228	7503	0.430	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	6628	0.422	ng/ul	99
22) Benzo(k)fluoranthene	22.70	252	6948	0.418	ng/ul	99
23) Benzo(a)pyrene	23.20	252	5420	0.408	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.42	276	7577	0.417	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.43	278	6254	0.421	ng/ul	99
26) Benzo(g,h,i)perylene	26.05	276	6911	0.411	ng/ul	98

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

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