

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072020\
 Data File : BN011210.D
 Acq On : 20 Jul 2020 15:55
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.411

Quant Time: Jul 20 16:44:30 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN072020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 20 12:53:55 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.49	152	1270	0.40	ng/ul	0.00
2) Naphthalene-d8	10.24	136	4539	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	2657	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.88	188	5672	0.40	ng/ul	0.00
16) Chrysene-d12	21.10	240	3902	0.40	ng/ul	0.00
20) Perylene-d12	23.23	264	3391	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.84	152	4722	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	10321	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.29	128	5598	0.412	ng/ul	99
5) 2-Methylnaphthalene	11.91	142	3721	0.419	ng/ul	100
7) Acenaphthylene	13.83	152	5318	0.413	ng/ul	100
8) Acenaphthene	14.19	153	4248	0.420	ng/ul	98
9) Fluorene	15.18	166	4827	0.420	ng/ul	97
11) Pentachlorophenol	16.54	266	516	0.426	ng/ul	97
12) Phenanthrene	16.92	178	7757	0.415	ng/ul	99
13) Anthracene	17.01	178	6414	0.416	ng/ul	99
15) Fluoranthene	18.95	202	8768	0.420	ng/ul	99
17) Pyrene	19.31	202	8743	0.420	ng/ul	99
18) Benzo(a)anthracene	21.08	228	5812	0.421	ng/ul	98
19) Chrysene	21.13	228	7566	0.412	ng/ul	100
21) Benzo(b)fluoranthene	22.61	252	5896	0.403	ng/ul	98
22) Benzo(k)fluoranthene	22.65	252	7088	0.422	ng/ul	100
23) Benzo(a)pyrene	23.14	252	5366	0.436	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.32	276	6243	0.430	ng/ul#	58
25) Dibenzo(a,h)anthracene	25.34	278	4962	0.453	ng/ul	96
26) Benzo(g,h,i)perylene	25.95	276	5683	0.394	ng/ul	98

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

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