

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072120\
 Data File : BN011212.D
 Acq On : 21 Jul 2020 09:19
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.412

Quant Time: Jul 21 10:09:10 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 : Tue Jul 21 10:07:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.49	152	1034	0.40	ng/ul	0.00
2) Naphthalene-d8	10.24	136	3585	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	2155	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.87	188	4806	0.40	ng/ul	0.00
16) Chrysene-d12	21.09	240	3669	0.40	ng/ul	0.00
20) Perylene-d12	23.23	264	3301	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.84	152	3758	0.43	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	9227	0.45	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.28	128	4489	0.419	ng/ul	99
5) 2-Methylnaphthalene	11.91	142	2957	0.422	ng/ul	99
7) Acenaphthylene	13.83	152	4303	0.412	ng/ul	99
8) Acenaphthene	14.18	153	3428	0.418	ng/ul	97
9) Fluorene	15.18	166	3931	0.422	ng/ul	99
11) Pentachlorophenol	16.54	266	429	0.418	ng/ul	97
12) Phenanthrene	16.92	178	6586	0.416	ng/ul	99
13) Anthracene	17.01	178	5472	0.419	ng/ul	100
15) Fluoranthene	18.94	202	7870	0.445	ng/ul	98
17) Pyrene	19.31	202	7884	0.402	ng/ul	99
18) Benzo(a)anthracene	21.08	228	5429	0.419	ng/ul	98
19) Chrysene	21.13	228	7151	0.414	ng/ul	100
21) Benzo(b)fluoranthene	22.60	252	5717	0.401	ng/ul	98
22) Benzo(k)fluoranthene	22.64	252	6946	0.425	ng/ul	99
23) Benzo(a)pyrene	23.14	252	5012	0.418	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.32	276	5930	0.420	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.34	278	4514	0.424	ng/ul	96
26) Benzo(g,h,i)perylene	25.94	276	5588	0.398	ng/ul	99

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

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