

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120620\
 Data File : BN012930.D
 Acq On : 06 Dec 2020 23:19
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.417

Quant Time: Dec 07 04:10:30 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 07 03:06:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1204	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	4940	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.06	164	2763	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	5585	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	4545	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	4582	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	2977	0.41	ng/ul	-0.02
14) Fluoranthene-d10	18.86	212	5910	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.20	128	5664	0.396	ng/ul	93
5) 2-Methylnaphthalene	11.84	142	3778	0.402	ng/ul	98
7) Acenaphthylene	13.77	152	5392	0.412	ng/ul	95
8) Acenaphthene	14.12	153	4273	0.398	ng/ul	100
9) Fluorene	15.12	166	4660	0.379	ng/ul	97
11) Pentachlorophenol	16.48	266	679	0.420	ng/ul	95
12) Phenanthrene	16.86	178	7478	0.399	ng/ul	98
13) Anthracene	16.95	178	6751	0.424	ng/ul	96
15) Fluoranthene	18.89	202	8582	0.386	ng/ul	97
17) Pyrene	19.25	202	8777	0.409	ng/ul	99
18) Benzo(a)anthracene	21.02	228	7641	0.406	ng/ul	99
19) Chrysene	21.07	228	8342	0.402	ng/ul	99
21) Benzo(b)fluoranthene	22.52	252	8454	0.399	ng/ul	95
22) Benzo(k)fluoranthene	22.56	252	9260	0.395	ng/ul	95
23) Benzo(a)pyrene	23.05	252	7854	0.399	ng/ul	91
24) Indeno(1,2,3-cd)pyrene	25.19	276	9117	0.353	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.21	278	7322	0.358	ng/ul	94
26) Benzo(g,h,i)perylene	25.82	276	7741	0.331	ng/ul	98

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

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