

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101420\
 Data File : BN012285.D
 Acq On : 17 Oct 2020 02:07
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 97 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.422

Quant Time: Oct 17 03:22:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 16 23:42:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1836	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	7920	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	4333	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	8469	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	7014	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	7065	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	5019	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	9424	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	9203	0.403	ng/ul	98
5) 2-Methylnaphthalene	12.06	142	6039	0.416	ng/ul	100
7) Acenaphthylene	13.98	152	8675	0.462	ng/ul	98
8) Acenaphthene	14.32	153	6621	0.417	ng/ul	99
9) Fluorene	15.31	166	7223	0.408	ng/ul	98
11) Pentachlorophenol	16.66	266	1142	0.467	ng/ul	100
12) Phenanthrene	17.05	178	11013	0.408	ng/ul	98
13) Anthracene	17.14	178	10637	0.459	ng/ul	98
15) Fluoranthene	19.07	202	13009	0.412	ng/ul	99
17) Pyrene	19.44	202	13365	0.423	ng/ul	98
18) Benzo(a)anthracene	21.19	228	11322	0.436	ng/ul	95
19) Chrysene	21.24	228	11918	0.377	ng/ul	96
21) Benzo(b)fluoranthene	22.74	252	11590	0.406	ng/ul	83
22) Benzo(k)fluoranthene	22.78	252	11914	0.377	ng/ul#	91
23) Benzo(a)pyrene	23.30	252	11237	0.405	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.58	276	14229	0.399	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.59	278	11443	0.401	ng/ul#	93
26) Benzo(g,h,i)perylene	26.25	276	12258	0.375	ng/ul#	89

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

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