

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022021\
 Data File : BN013685.D
 Acq On : 20 Feb 2021 10:34
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.415

Quant Time: Feb 21 06:56:51 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN020421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 16 16:31:21 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	4961	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	18940	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	9666	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	19564	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	15866	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	13822	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	11541	0.43	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	20182	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.35	128	22002	0.446	ng/ul	100
5) 2-Methylnaphthalene	11.98	142	14122	0.427	ng/ul	100
7) Acenaphthylene	13.89	152	16027	0.445	ng/ul	99
8) Acenaphthene	14.24	153	14310	0.446	ng/ul	99
9) Fluorene	15.24	166	15642	0.431	ng/ul	99
11) Pentachlorophenol	16.59	266	1099	0.461	ng/ul	100
12) Phenanthrene	16.97	178	25462	0.437	ng/ul	100
13) Anthracene	17.06	178	20944	0.434	ng/ul	100
15) Fluoranthene	19.00	202	26668	0.425	ng/ul	99
17) Pyrene	19.36	202	26721	0.468	ng/ul	98
18) Benzo(a)anthracene	21.12	228	21304	0.414	ng/ul	99
19) Chrysene	21.17	228	25909	0.462	ng/ul	100
21) Benzo(b)fluoranthene	22.65	252	24284	0.430	ng/ul	98
22) Benzo(k)fluoranthene	22.70	252	25906	0.443	ng/ul	99
23) Benzo(a)pyrene	23.20	252	20953	0.461	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.44	276	26378	0.437	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.46	278	20847	0.420	ng/ul	99
26) Benzo(g,h,i)perylene	26.10	276	23870	0.450	ng/ul	100

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

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