

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012721\
 Data File : BN013490.D
 Acq On : 26 Jan 2021 15:51
 Operator : CG/JU
 Sample : SST00.202
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.202

Quant Time: Jan 26 17:07:39 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN012721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 26 17:05:29 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	4163	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	14785	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	7566	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	15734	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	12901	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	11807	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	4222	0.19	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	7730	0.18	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	8326	0.195	ng/ul	98
5) 2-Methylnaphthalene	12.00	142	5226	0.186	ng/ul	99
7) Acenaphthylene	13.92	152	5946	0.166	ng/ul	99
8) Acenaphthene	14.26	153	5322	0.181	ng/ul	100
9) Fluorene	15.25	166	5881	0.174	ng/ul	100
11) Pentachlorophenol	16.61	266	369	0.081	ng/ul	98
12) Phenanthrene	16.99	178	9898	0.188	ng/ul	99
13) Anthracene	17.08	178	7905	0.176	ng/ul	98
15) Fluoranthene	19.02	202	10426	0.166	ng/ul	100
17) Pyrene	19.38	202	10608	0.174	ng/ul	100
18) Benzo(a)anthracene	21.14	228	8084	0.151	ng/ul	99
19) Chrysene	21.19	228	10527	0.179	ng/ul	100
21) Benzo(b)fluoranthene	22.68	252	9689	0.177	ng/ul	94
22) Benzo(k)fluoranthene	22.72	252	10350	0.171	ng/ul	94
23) Benzo(a)pyrene	23.23	252	8192	0.162	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.48	276	10515	0.158	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.50	278	8440	0.160	ng/ul	95
26) Benzo(g,h,i)perylene	26.13	276	9687	0.161	ng/ul	98

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

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