

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012721\
 Data File : BN013489.D
 Acq On : 26 Jan 2021 15:16
 Operator : CG/JU
 Sample : SST00.101
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.101

Quant Time: Jan 26 17:08:45 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	4055	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	14463	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	7393	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	15199	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	12651	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	12355	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	2068	0.10	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	3726	0.09	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.38	128	4052	0.097	ng/ul	96
5) 2-Methylnaphthalene	12.00	142	2526	0.092	ng/ul	100
7) Acenaphthylene	13.92	152	2880	0.082	ng/ul	96
8) Acenaphthene	14.27	153	2605	0.091	ng/ul	97
9) Fluorene	15.26	166	2816	0.085	ng/ul	99
12) Phenanthrene	16.99	178	4731	0.093	ng/ul	98
13) Anthracene	17.08	178	3753	0.087	ng/ul	94
15) Fluoranthene	19.02	202	4921	0.081	ng/ul	97
17) Pyrene	19.38	202	5124	0.086	ng/ul	97
18) Benzo(a)anthracene	21.14	228	3971	0.076	ng/ul	99
19) Chrysene	21.19	228	5143	0.089	ng/ul	99
21) Benzo(b)fluoranthene	22.68	252	4757	0.083	ng/ul	83
22) Benzo(k)fluoranthene	22.72	252	5178	0.082	ng/ul#	83
23) Benzo(a)pyrene	23.23	252	4101	0.077	ng/ul#	72
24) Indeno(1,2,3-cd)pyrene	25.48	276	5273	0.076	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.50	278	4202	0.076	ng/ul#	90
26) Benzo(a,h,i)perylene	26.14	276	4853	0.077	ng/ul	94

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

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