

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN020421\
 Data File : BN013567.D
 Acq On : 04 Feb 2021 12:32
 Operator : CG/JU
 Sample : SST0.404
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.404

Quant Time: Feb 04 13:02:20 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 : Thu Feb 04 13:01:49 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	4520	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	16356	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	8758	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	18531	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	17001	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	13289	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	9785	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	18971	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	17982	0.380	ng/ul	100
5) 2-Methylnaphthalene	11.99	142	11926	0.397	ng/ul	100
7) Acenaphthylene	13.91	152	13432	0.372	ng/ul	100
8) Acenaphthene	14.25	153	12018	0.374	ng/ul	100
9) Fluorene	15.24	166	13549	0.382	ng/ul	100
11) Pentachlorophenol	16.60	266	835	0.306	ng/ul	100
12) Phenanthrene	16.98	178	22624	0.373	ng/ul	100
13) Anthracene	17.07	178	18800	0.378	ng/ul	100
15) Fluoranthene	19.01	202	24159	0.364	ng/ul	100
17) Pyrene	19.37	202	24670	0.355	ng/ul	100
18) Benzo(a)anthracene	21.13	228	22683	0.404	ng/ul	100
19) Chrysene	21.18	228	24982	0.357	ng/ul	100
21) Benzo(b)fluoranthene	22.66	252	22315	0.400	ng/ul	100
22) Benzo(k)fluoranthene	22.70	252	23130	0.389	ng/ul	100
23) Benzo(a)pyrene	23.21	252	17938	0.379	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.46	276	23808	0.379	ng/ul	100
25) Dibenzo(a,h)anthracene	25.47	278	19508	0.384	ng/ul	100
26) Benzo(g,h,i)perylene	26.12	276	20982	0.377	ng/ul	100

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

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