

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022121\
 Data File : BN013697.D
 Acq On : 20 Feb 2021 20:27
 Operator : CG/JU
 Sample : SST00.806
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.806

Quant Time: Feb 20 23:11:10 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN022121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 20 23:07:21 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	3908	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	14761	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	8168	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	17657	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	16550	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	15354	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	16707	0.79	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	36567	0.84	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.35	128	29712	0.773	ng/ul	99
5) 2-Methylnaphthalene	11.97	142	20072	0.779	ng/ul	100
7) Acenaphthylene	13.90	152	23949	0.787	ng/ul	99
8) Acenaphthene	14.25	153	20971	0.774	ng/ul	99
9) Fluorene	15.24	166	24200	0.790	ng/ul	99
11) Pentachlorophenol	16.59	266	3048	1.418	ng/ul	98
12) Phenanthrene	16.97	178	39768	0.757	ng/ul	100
13) Anthracene	17.06	178	35169	0.807	ng/ul	99
15) Fluoranthene	19.00	202	45139	0.796	ng/ul	100
17) Pyrene	19.37	202	46363	0.779	ng/ul	96
18) Benzo(a)anthracene	21.12	228	43072	0.802	ng/ul	99
19) Chrysene	21.18	228	44955	0.769	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	44873	0.715	ng/ul	95
22) Benzo(k)fluoranthene	22.70	252	46070	0.708	ng/ul	94
23) Benzo(a)pyrene	23.21	252	39371	0.780	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	25.45	276	51157	0.762	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.47	278	41656	0.756	ng/ul	97
26) Benzo(g,h,i)perylene	26.10	276	43844	0.743	ng/ul	99

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

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