

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022121\
 Data File : BN013696.D
 Acq On : 20 Feb 2021 19:51
 Operator : CG/JU
 Sample : SST0.405
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.405

Quant Time: Feb 20 23:08:22 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	3607	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	13387	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	7532	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	16422	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	15465	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	14745	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	8528	0.45	ng/ul	0.00
14) Fluoranthene-d10	18.97	212	18906	0.47	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	15692	0.450	ng/ul	100
5) 2-Methylnaphthalene	11.97	142	10318	0.442	ng/ul	100
7) Acenaphthylene	13.90	152	12742	0.454	ng/ul	100
8) Acenaphthene	14.25	153	11057	0.443	ng/ul	100
9) Fluorene	15.24	166	12697	0.449	ng/ul	100
11) Pentachlorophenol	16.59	266	1588	0.794	ng/ul	100
12) Phenanthrene	16.98	178	21496	0.440	ng/ul	100
13) Anthracene	17.07	178	18697	0.461	ng/ul	100
15) Fluoranthene	19.00	202	24690	0.468	ng/ul	100
17) Pyrene	19.36	202	25700	0.462	ng/ul	100
18) Benzo(a)anthracene	21.12	228	23118	0.461	ng/ul	100
19) Chrysene	21.18	228	24376	0.446	ng/ul	100
21) Benzo(b)fluoranthene	22.66	252	24461	0.406	ng/ul	100
22) Benzo(k)fluoranthene	22.70	252	25195	0.403	ng/ul	100
23) Benzo(a)pyrene	23.21	252	22748	0.470	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.45	276	28272	0.439	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.47	278	23105	0.437	ng/ul	100
26) Benzo(g,h,i)perylene	26.11	276	24568	0.434	ng/ul	100

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

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