

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN090820\
 Data File : BN011732.D
 Acq On : 08 Sep 2020 12:14
 Operator : CG/JU
 Sample : SST0.405
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.405

Quant Time: Sep 08 12:48:24 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN090820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 08 12:47:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1798	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	6489	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	3514	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	7655	0.40	ng/ul	0.00
16) Chrysene-d12	21.32	240	6338	0.40	ng/ul	0.00
20) Perylene-d12	23.57	264	5954	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	4201	0.46	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	8179	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	7766	0.411	ng/ul	100
5) 2-Methylnaphthalene	12.19	142	5210	0.436	ng/ul	100
7) Acenaphthylene	14.09	152	6903	0.390	ng/ul	100
8) Acenaphthene	14.44	153	5604	0.407	ng/ul	100
9) Fluorene	15.43	166	6371	0.393	ng/ul	100
11) Pentachlorophenol	16.77	266	913	0.502	ng/ul	100
12) Phenanthrene	17.16	178	10443	0.403	ng/ul	100
13) Anthracene	17.25	178	9202	0.432	ng/ul	100
15) Fluoranthene	19.18	202	11539	0.387	ng/ul	100
17) Pyrene	19.55	202	11471	0.463	ng/ul	100
18) Benzo(a)anthracene	21.30	228	9923	0.467	ng/ul	100
19) Chrysene	21.35	228	11244	0.392	ng/ul	100
21) Benzo(b)fluoranthene	22.89	252	11157	0.431	ng/ul	100
22) Benzo(k)fluoranthene	22.94	252	11654	0.407	ng/ul	100
23) Benzo(a)pyrene	23.47	252	9870	0.454	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.83	276	13115	0.446	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.84	278	10676	0.445	ng/ul	100
26) Benzo(g,h,i)perylene	26.52	276	11372	0.413	ng/ul	100

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

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