

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN090820\
 Data File : BN011733.D
 Acq On : 08 Sep 2020 12:51
 Operator : CG/JU
 Sample : SST0.806
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.806

Quant Time: Sep 08 13:22:17 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	1488	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	5349	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	2893	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	6062	0.40	ng/ul	0.00
16) Chrysene-d12	21.32	240	5336	0.40	ng/ul	0.00
20) Perylene-d12	23.57	264	4825	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	7047	0.93	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	13520	0.79	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	12901	0.827	ng/ul	99
5) 2-Methylnaphthalene	12.19	142	8683	0.881	ng/ul	100
7) Acenaphthylene	14.09	152	11525	0.791	ng/ul	98
8) Acenaphthene	14.44	153	9156	0.808	ng/ul	98
9) Fluorene	15.43	166	10641	0.796	ng/ul	99
11) Pentachlorophenol	16.77	266	1522	1.058	ng/ul	98
12) Phenanthrene	17.16	178	16953	0.827	ng/ul	99
13) Anthracene	17.25	178	15314	0.908	ng/ul	100
15) Fluoranthene	19.18	202	19197	0.814	ng/ul	99
17) Pyrene	19.55	202	18833	0.903	ng/ul	100
18) Benzo(a)anthracene	21.30	228	17292	0.966	ng/ul	99
19) Chrysene	21.35	228	19222	0.796	ng/ul	98
21) Benzo(b)fluoranthene	22.89	252	18945	0.903	ng/ul	97
22) Benzo(k)fluoranthene	22.94	252	19691	0.849	ng/ul	95
23) Benzo(a)pyrene	23.47	252	16249	0.922	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	25.83	276	21705	0.911	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.85	278	17863	0.920	ng/ul	96
26) Benzo(g,h,i)perylene	26.52	276	18903	0.848	ng/ul	99

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

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