

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN060220\
 Data File : BN010742.D
 Acq On : 02 Jun 2020 12:40
 Operator : CG/JU
 Sample : SST00.206
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.206

Quant Time: Jun 02 14:33:29 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN060220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 02 14:30:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1767	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	5785	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	3015	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	5994	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	5103	0.40	ng/ul	0.00
20) Perylene-d12	23.30	264	4926	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	1907	0.19	ng/ul	0.01
14) Fluoranthene-d10	18.98	212	3515	0.19	ng/ul	0.00

Target Compounds						Qvalue
3) Naphthalene	10.37	128	3699	0.20	ng/ul	96
5) 2-Methylnaphthalene	12.00	142	2326	0.19	ng/ul	99
7) Acenaphthylene	13.91	152	2930	0.18	ng/ul	97
8) Acenaphthene	14.26	153	2417	0.20	ng/ul	97
9) Fluorene	15.25	166	2649	0.18	ng/ul	100
11) Pentachlorophenol	16.61	266	291	0.25	ng/ul	99
12) Phenanthrene	16.98	178	4200	0.19	ng/ul	98
13) Anthracene	17.08	178	3434	0.18	ng/ul	97
15) Fluoranthene	19.01	202	4559	0.17	ng/ul	97
17) Pyrene	19.37	202	4662	0.19	ng/ul#	94
18) Benzo(a)anthracene	21.13	228	3892	0.18	ng/ul	98
19) Chrysene	21.18	228	4993	0.20	ng/ul	98
21) Benzo(b)fluoranthene	22.66	252	4368	0.20	ng/ul	87
22) Benzo(k)fluoranthene	22.70	252	4908	0.21	ng/ul#	87
23) Benzo(a)pyrene	23.20	252	3887	0.20	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.42	276	5357	0.22	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.43	278	4346	0.22	ng/ul	91
26) Benzo(g,h,i)perylene	26.06	276	4761	0.22	ng/ul	97

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

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