

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082318\
 Data File : BN002460.D
 Acq On : 23 Aug 2018 11:51
 Operator : JU/SJ
 Sample : SST0.410
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST0.410

Quant Time: Aug 23 13:12:21 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 23 13:07:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	1552	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	6305	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	3636	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	9194	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	8870	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	7528	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	4004	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	10672	0.38	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.50	128	6478	0.399	ng/ul#	97
5) 2-Methylnaphthalene	12.13	142	4519	0.389	ng/ul	100
7) Acenaphthylene	14.03	152	6883	0.387	ng/ul	97
8) Acenaphthene	14.38	153	5283	0.404	ng/ul	95
9) Fluorene	15.37	166	6176	0.405	ng/ul#	93
11) Pentachlorophenol	16.74	266	389	0.223	ng/ul	97
12) Phenanthrene	17.10	178	10895	0.397	ng/ul#	89
13) Anthracene	17.20	178	9349	0.360	ng/ul#	92
15) Fluoranthene	19.12	202	12607	0.387	ng/ul	97
17) Pyrene	19.49	202	12165	0.377	ng/ul	96
18) Benzo(a)anthracene	21.24	228	10666	0.348	ng/ul#	85
19) Chrysene	21.30	228	12247	0.401	ng/ul	97
21) Benzo(b)fluoranthene	22.82	252	11217	0.411	ng/ul	85
22) Benzo(k)fluoranthene	22.86	252	11436	0.420	ng/ul#	88
23) Benzo(a)pyrene	23.39	252	10001	0.394	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.71	276	10194	0.358	ng/ul#	84
25) Dibenzo(a,h)anthracene	25.72	278	8013	0.353	ng/ul	91
26) Benzo(g,h,i)perylene	26.39	276	8808	0.364	ng/ul	96

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

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