

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080219\
 Data File : BM021765.D
 Acq On : 02 Aug 2019 19:46
 Operator : HP/JU
 Sample : SSTD1.617
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.617

Quant Time: Aug 02 23:41:31 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM080219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 02 23:36:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	1055	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	4427	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.30	164	1839	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	3505	0.40	ng/ul	-0.02
16) Chrysene-d12	21.25	240	2643	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	2302	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	10262	1.94	ng/ul	-0.03
14) Fluoranthene-d10	19.09	212	12691	1.26	ng/ul	-0.01

Target Compounds				Ovalue		
3) Naphthalene	10.48	128	19975	1.787	ng/ul#	93
5) 2-Methylnaphthalene	12.11	142	12682	2.149	ng/ul	97
7) Acenaphthylene	14.03	152	14811	1.474	ng/ul	96
8) Acenaphthene	14.37	153	12782	1.726	ng/ul	95
9) Fluorene	15.36	166	13800	1.660	ng/ul	97
11) Pentachlorophenol	16.73	266	807	0.674	ng/ul#	89
12) Phenanthrene	17.10	178	19669	2.029	ng/ul	96
13) Anthracene	17.20	178	16355	1.960	ng/ul	94
15) Fluoranthene	19.13	202	18200	1.502	ng/ul	96
17) Pyrene	19.49	202	18012	2.406	ng/ul	95
18) Benzo(a)anthracene	21.24	228	12787	1.553	ng/ul	96
19) Chrysene	21.29	228	16647	1.604	ng/ul	98
21) Benzo(b)fluoranthene	22.81	252	14397	1.694	ng/ul	84
22) Benzo(k)fluoranthene	22.86	252	15905	1.847	ng/ul#	81
23) Benzo(a)pyrene	23.38	252	12723	1.666	ng/ul#	71
24) Indeno(1,2,3-cd)pyrene	25.71	276	15785	1.544	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.72	278	12702	1.586	ng/ul#	80
26) Benzo(g,h,i)perylene	26.39	276	14445	1.475	ng/ul	92

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

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