

Data Path : Z:\HPCHEM1\BNA M\DATA\BM030317\
 Data File : BM009066.D
 Acq On : 03 Mar 2017 13:39
 Operator : SJ/MA
 Sample : SSTD0.134
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.134

Quant Time: Mar 03 15:30:00 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 03 15:23:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	161	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	476	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	263	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.28	188	606	0.40	ng/ul	0.00
16) Chrysene-d12	21.45	240	696	0.40	ng/ul	0.00
20) Perylene-d12	23.78	264	643	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	76	0.10	ng/ul	0.00
14) Fluoranthene-d10	19.30	212	185	0.10	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.74	128	139	0.11	ng/ul#	32
5) 2-Methylnaphthalene	12.35	142	101	0.11	ng/ul	94
7) Acenaphthylene	14.25	152	138	0.12	ng/ul#	42
8) Acenaphthene	14.59	153	120	0.14	ng/ul#	89
9) Fluorene	15.58	166	119	0.11	ng/ul#	85
12) Phenanthrene	17.31	178	193	0.10	ng/ul#	57
13) Anthracene	17.40	178	186	0.10	ng/ul#	51
15) Fluoranthene	19.33	202	255	0.11	ng/ul	65
17) Pyrene	19.69	202	248	0.09	ng/ul#	62
18) Benzo(a)anthracene	21.43	228	239	0.09	ng/ul#	69
19) Chrysene	21.48	228	233	0.10	ng/ul#	71
21) Benzo(b)fluoranthene	23.08	252	280	0.10	ng/ul#	49
22) Benzo(k)fluoranthene	23.12	252	247	0.10	ng/ul#	45
23) Benzo(a)pyrene	23.68	252	242	0.10	ng/ul#	52
24) Indeno(1,2,3-cd)pyrene	26.17	276	301	0.10	ng/ul#	86
25) Dibenzo(a,h)anthracene	26.19	278	253	0.11	ng/ul#	48
26) Benzo(a,h,i)perylene	26.91	276	252	0.10	ng/ul#	54

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

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