

Data Path : Z:\HPCHEM1\BNA M\DATA\BM030317\
 Data File : BM009067.D
 Acq On : 03 Mar 2017 14:15
 Operator : SJ/MA
 Sample : SSTD0.235
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.235

Quant Time: Mar 03 15:30:32 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	125	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	351	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	205	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.28	188	452	0.40	ng/ul	0.00
16) Chrysene-d12	21.45	240	523	0.40	ng/ul	0.00
20) Perylene-d12	23.78	264	490	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.28	152	125	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.30	212	287	0.22	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.74	128	204	0.21	ng/ul#	48
5) 2-Methylnaphthalene	12.35	142	150	0.21	ng/ul	92
7) Acenaphthylene	14.25	152	205	0.23	ng/ul#	59
8) Acenaphthene	14.59	153	152	0.24	ng/ul#	86
9) Fluorene	15.58	166	188	0.23	ng/ul#	89
11) Pentachlorophenol	16.92	266	42	0.20	ng/ul#	83
12) Phenanthrene	17.31	178	276	0.20	ng/ul#	68
13) Anthracene	17.40	178	275	0.20	ng/ul#	67
15) Fluoranthene	19.33	202	367	0.21	ng/ul	73
17) Pyrene	19.69	202	354	0.17	ng/ul#	74
18) Benzo(a)anthracene	21.43	228	345	0.18	ng/ul#	79
19) Chrysene	21.48	228	320	0.17	ng/ul#	79
21) Benzo(b)fluoranthene	23.08	252	378	0.18	ng/ul	66
22) Benzo(k)fluoranthene	23.12	252	350	0.19	ng/ul#	62
23) Benzo(a)pyrene	23.68	252	350	0.18	ng/ul#	68
24) Indeno(1,2,3-cd)pyrene	26.17	276	441	0.20	ng/ul#	86
25) Dibenzo(a,h)anthracene	26.18	278	369	0.21	ng/ul#	63
26) Benzo(g,h,i)perylene	26.90	276	383	0.20	ng/ul#	68

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

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