SUPPLEMENTARY MATERIALS

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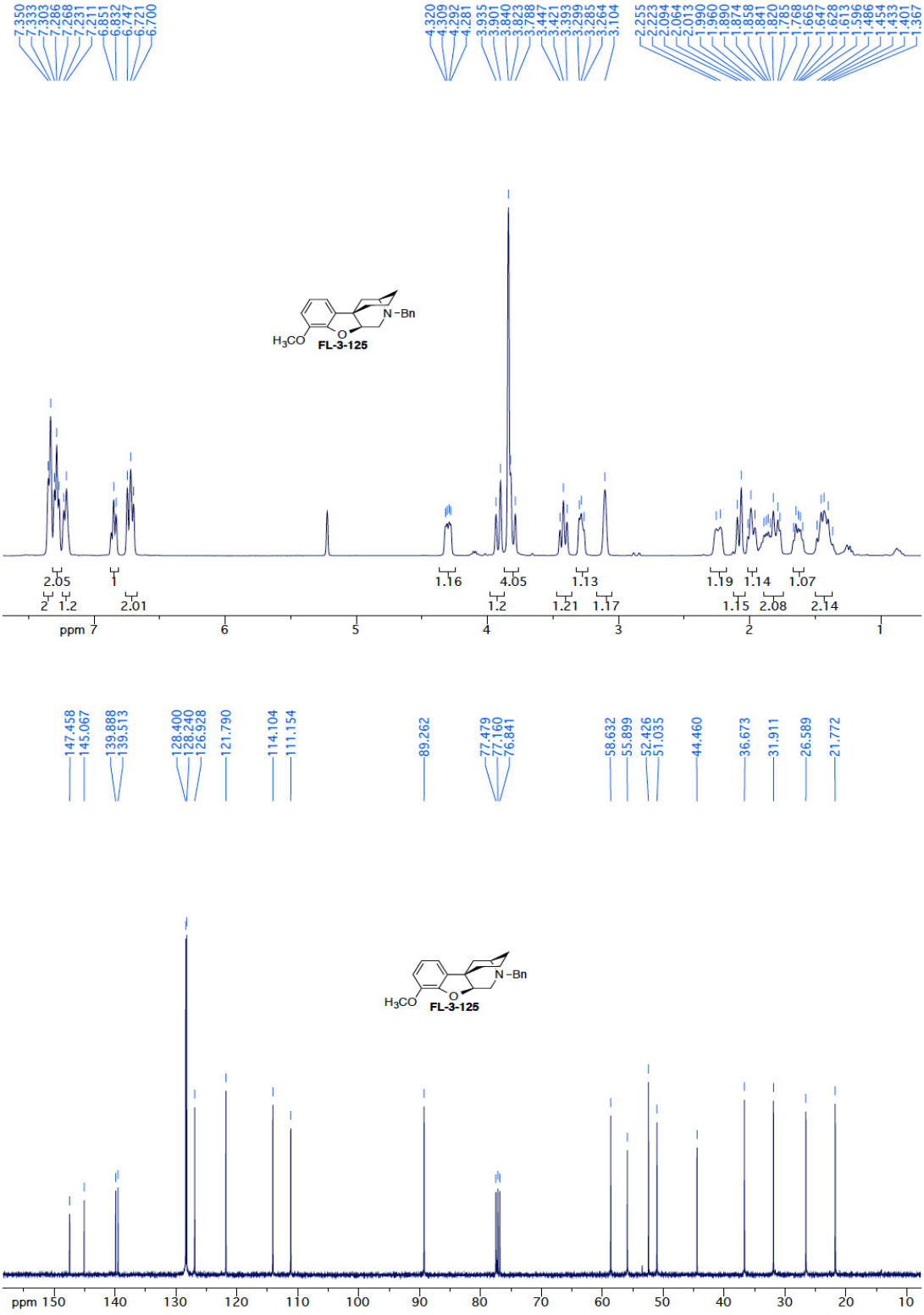
Compound **5**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4.320
4.309
4.292
4.281
3.935
3.901
3.840
3.823
3.788
3.447
3.421
3.393
3.299
3.285
3.264
3.104
2.255
2.223
2.094
2.064
2.013
1.990
1.960
1.890
1.874
1.858
1.841
1.820
1.765
1.668
1.645
1.647
1.628
1.613
1.596
1.486
1.454
1.433
1.401
1.367
N-Bn
O
H3CO
FL-3-125
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1
1.16
4.05
1.13
1.19
1.14
1.07
2
1.2
2.01
1.2
1.21
1.17
1.15
2.08
2.14
ppm 7
6
5
4
3
2
1
147.458
145.067
139.888
139.513
128.400
128.240
126.928
121.790
114.104
111.154
89.262
77.479
77.160
76.841
58.632
55.899
52.426
51.035
44.460
36.673
31.911
26.589
21.772
N-Bn
O
H3CO
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110
100
90
80
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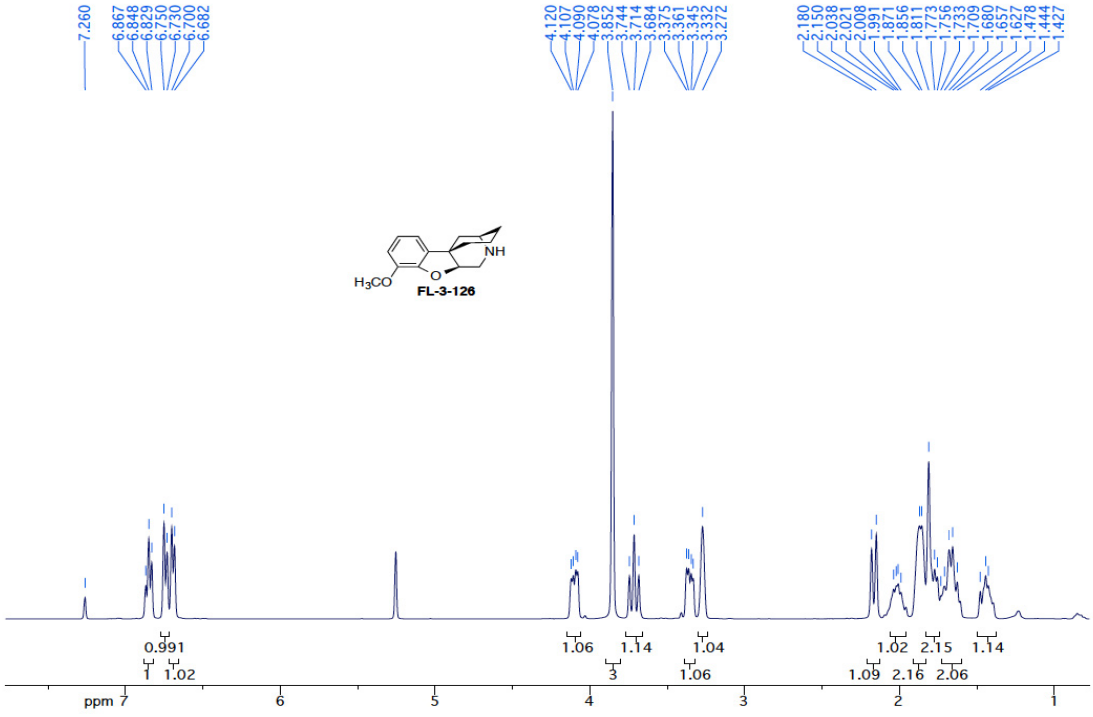
60
50
40
30
20
10

Compound **6**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4.078
3.852
3.744
3.714
3.684
3.375
3.361
3.345
3.332
3.272
2.180
2.150
2.038
2.021
2.008
1.991
1.871
1.856
1.811
1.773
1.756
1.733
1.709
1.680
1.657
1.627
1.478
1.444
1.427
NH
O
H3CO
FL-3-126
0.991
1.02
1.06
1.14
1.04
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1.14
1.09
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4
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2
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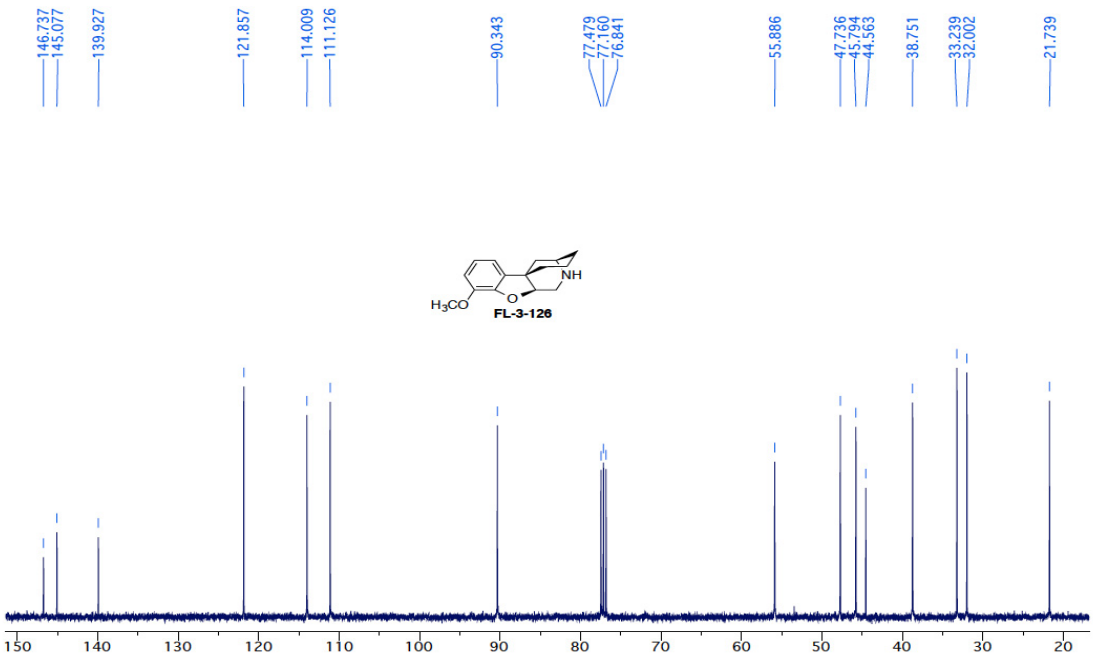
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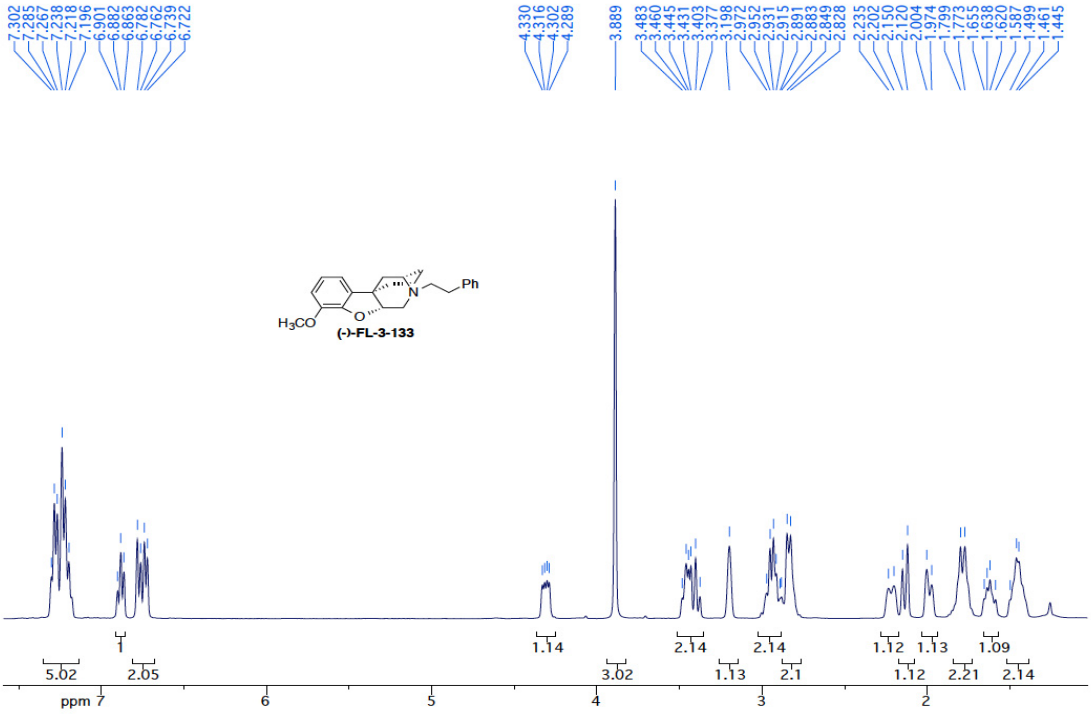
45.794
44.563
38.751
33.239
32.002
21.739
NH
O
H3CO
FL-3-126
150
140
130
120
110
100
90
80
70
60
50
40
30
20

Compound **7**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6.739
6.722
4.330
4.316
4.302
4.289
3.889
3.483
3.460
3.445
3.431
3.403
3.377
3.198
2.972
2.952
2.931
2.915
2.891
2.883
2.849
2.828
2.235
2.202
2.150
2.120
2.004
1.974
1.799
1.773
1.655
1.638
1.620
1.587
1.499
1.461
1.445
Ph
N
O
H3CO
(-)-FL-3-133
1
1.14
2.14
2.14
1.12
1.13
1.09
5.02
2.05
3.02
1.13
2.1
1.12
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2.14
ppm 7
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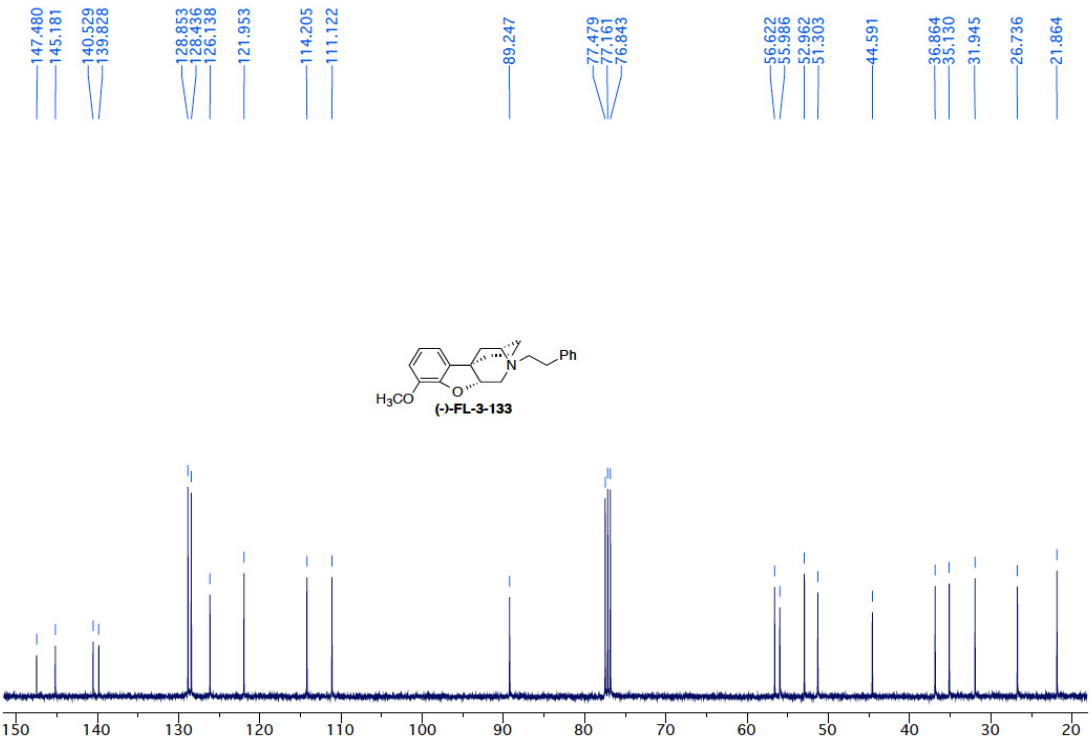
5
4
3
2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128.436
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114.205
111.122
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77.479
77.161
76.843
56.622
55.986
52.962
51.303
44.591
36.864
35.130
31.945
26.736
21.864
Ph
N
O
H3CO
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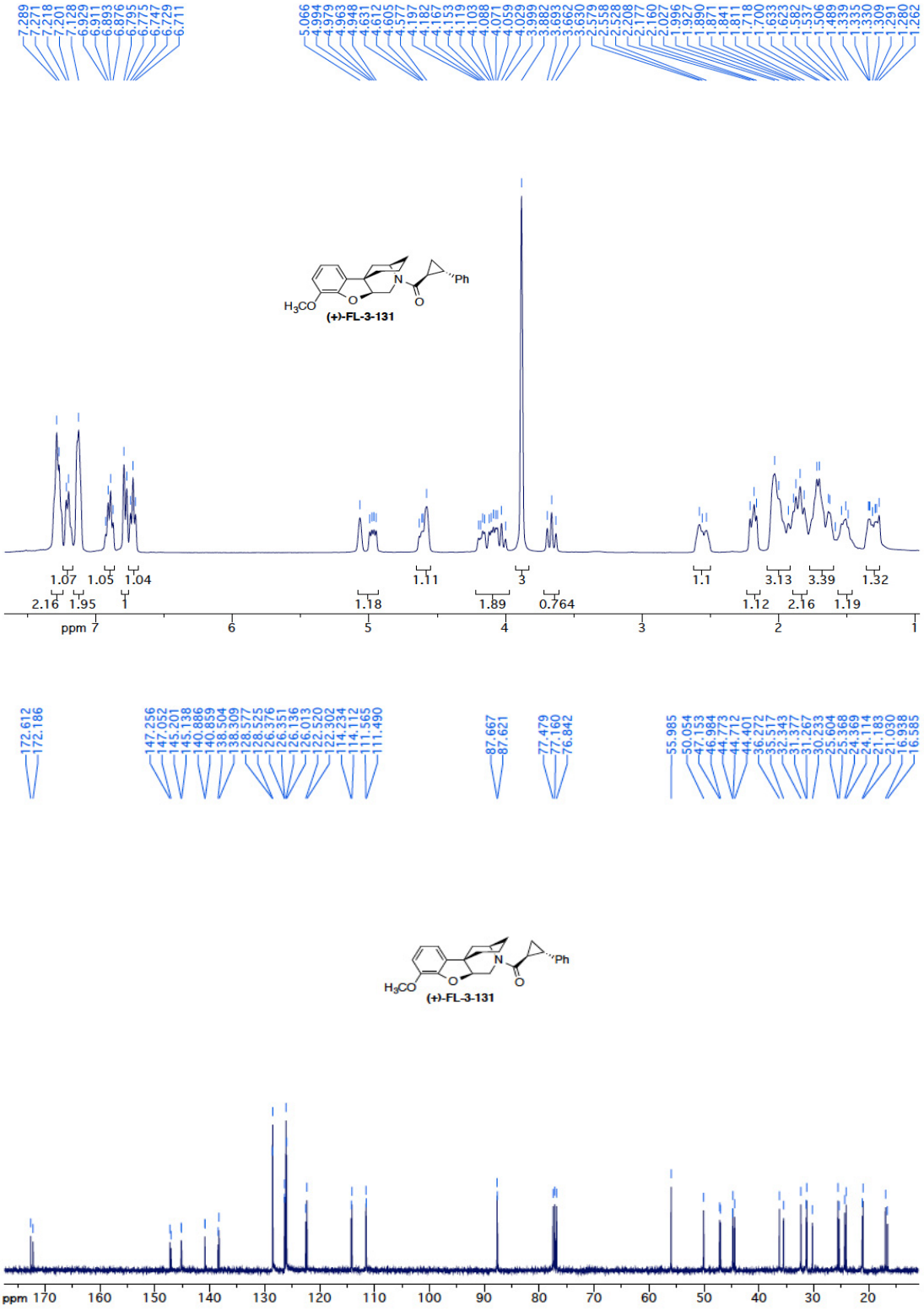
150
140
130
120
110
100
90
80
70
60
50
40
30
20

Compound (+)-8



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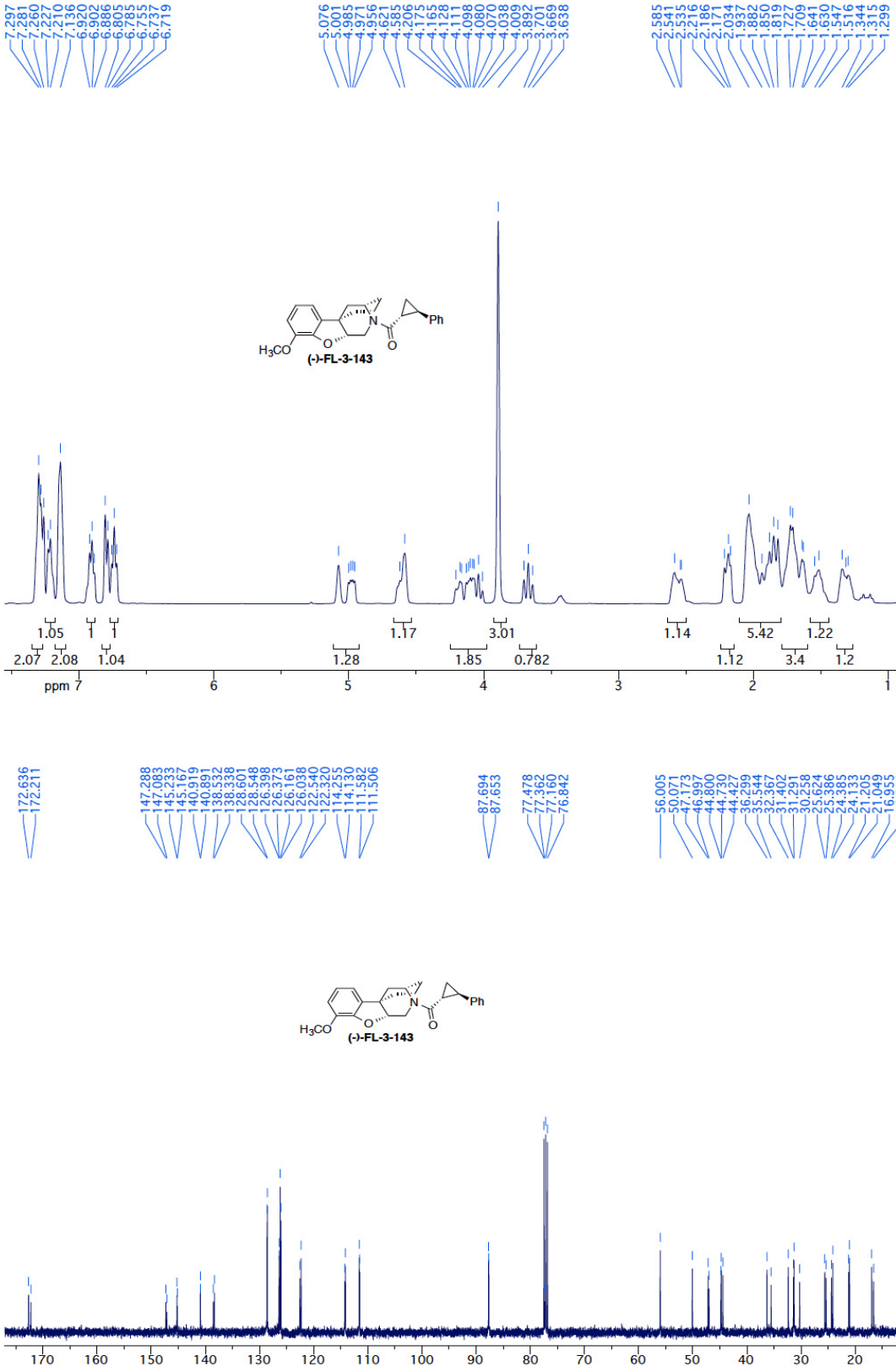
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6.886
6.805
6.785
6.755
6.737
6.719
5.076
5.001
4.995
4.971
4.956
4.621
4.585
4.296
4.175
4.165
4.128
4.111
4.098
4.080
4.070
4.038
4.009
3.892
3.701
3.669
3.638
2.585
2.541
2.535
2.216
2.186
2.171
2.034
1.937
1.882
1.850
1.819
1.727
1.709
1.641
1.630
1.547
1.516
1.344
1.315
1.299
H3CO
O
N
O
Ph
(-)-FL-3-143
1.05
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1
1.17
3.01
1.14
5.42
1.22
2.07
2.08
1.04
1.28
1.85
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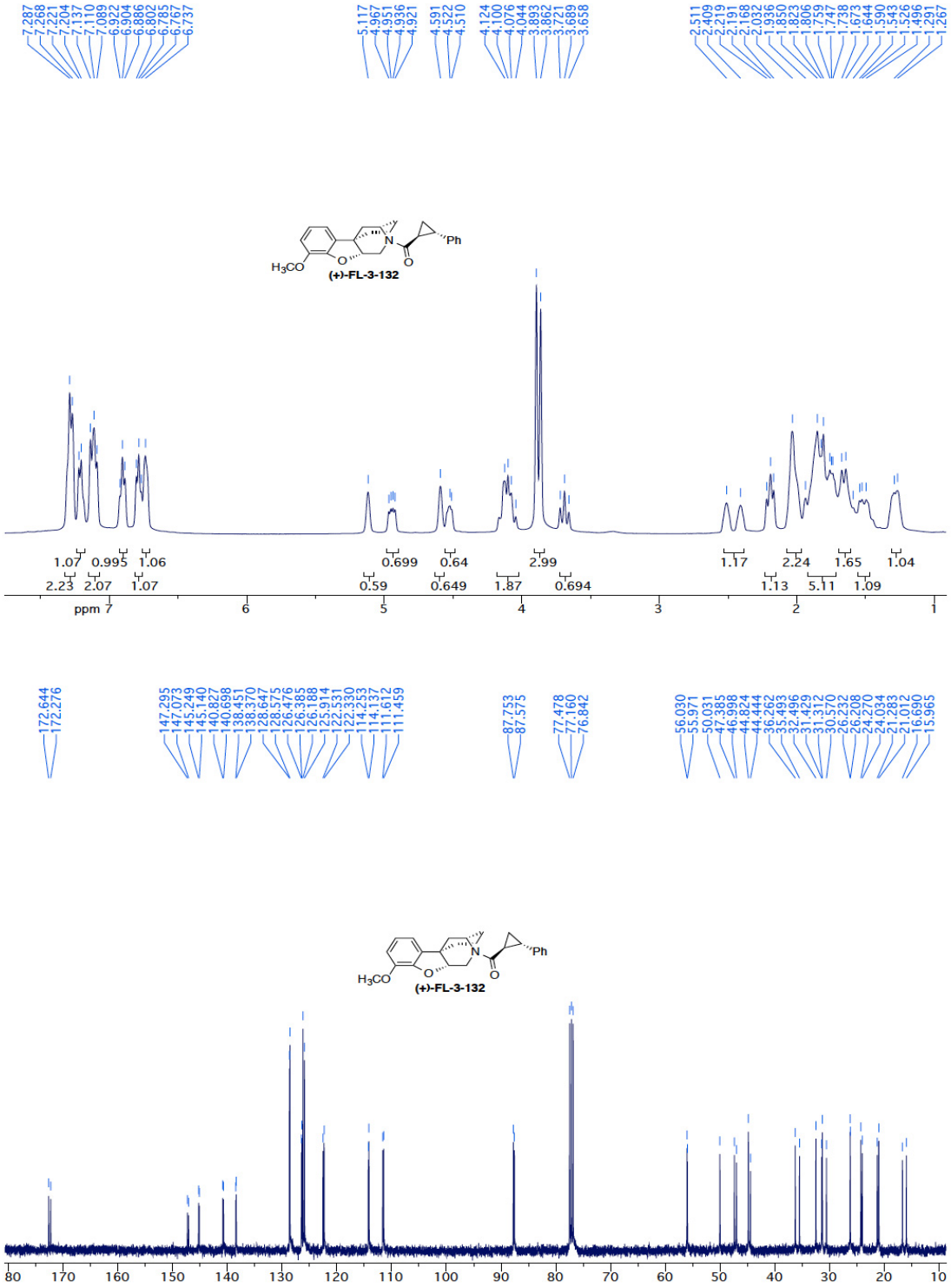
1.12
3.4
1.2
ppm 7
6
5
4
3
2
1
172.636
172.211
147.288
147.083
145.233
145.167
140.919
140.891
138.332
138.338
128.601
126.548
126.398
126.373
126.161
126.038
125.540
125.320
114.255
114.130
111.582
111.506
87.694
87.653
77.478
77.362
77.160
76.842
56.005
50.071
47.173
46.997
44.800
44.730
44.427
36.299
35.544
32.367
31.402
31.291
30.258
25.624
25.386
24.385
24.133
21.205
21.049
16.955
16.610
H3CO
O
N
O
Ph
(-)-FL-3-143
170
160
150
140
130
120
110
100
90
80
70
60
50
40
30
20
10

Compound (+)-**9**



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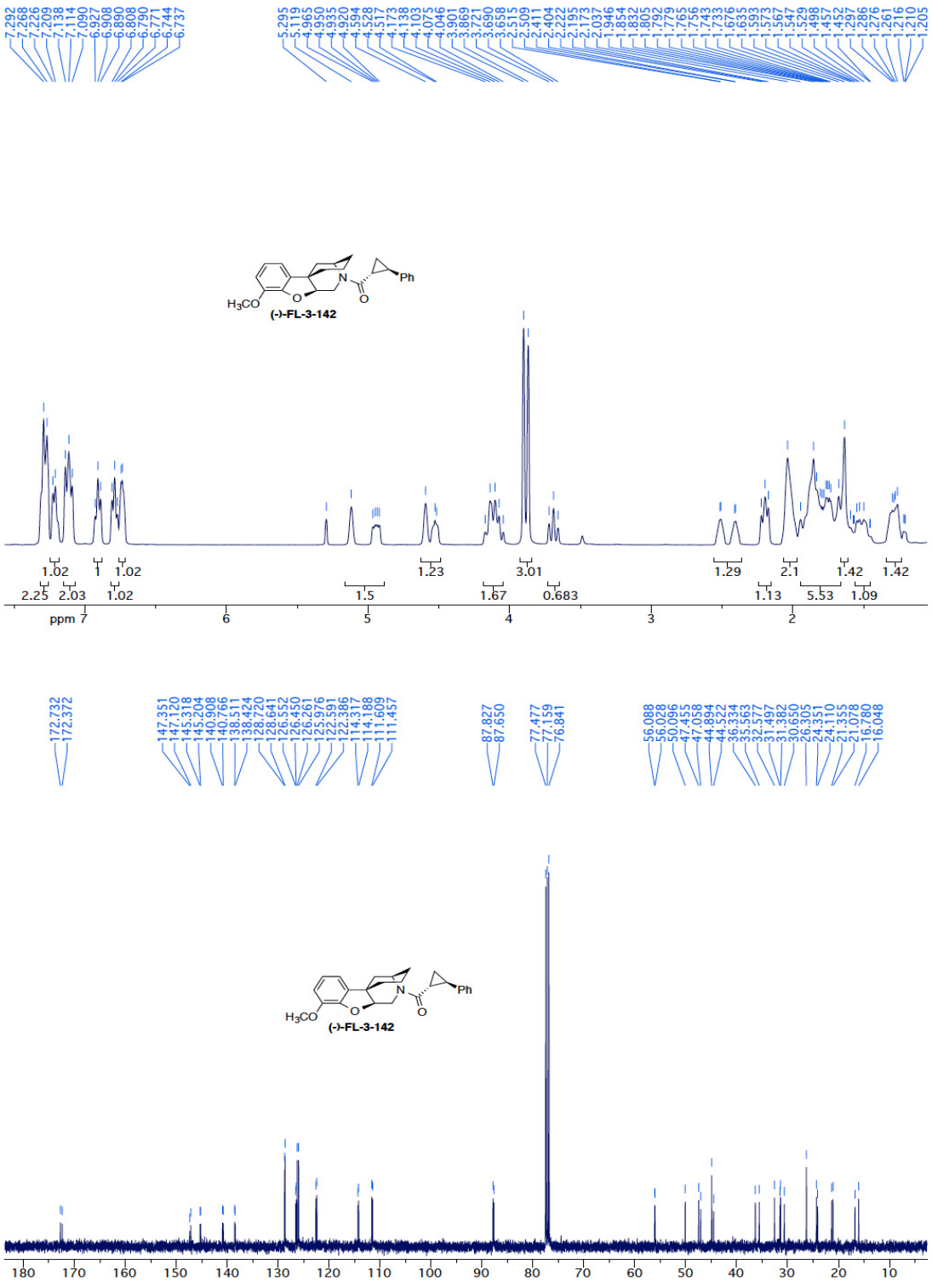
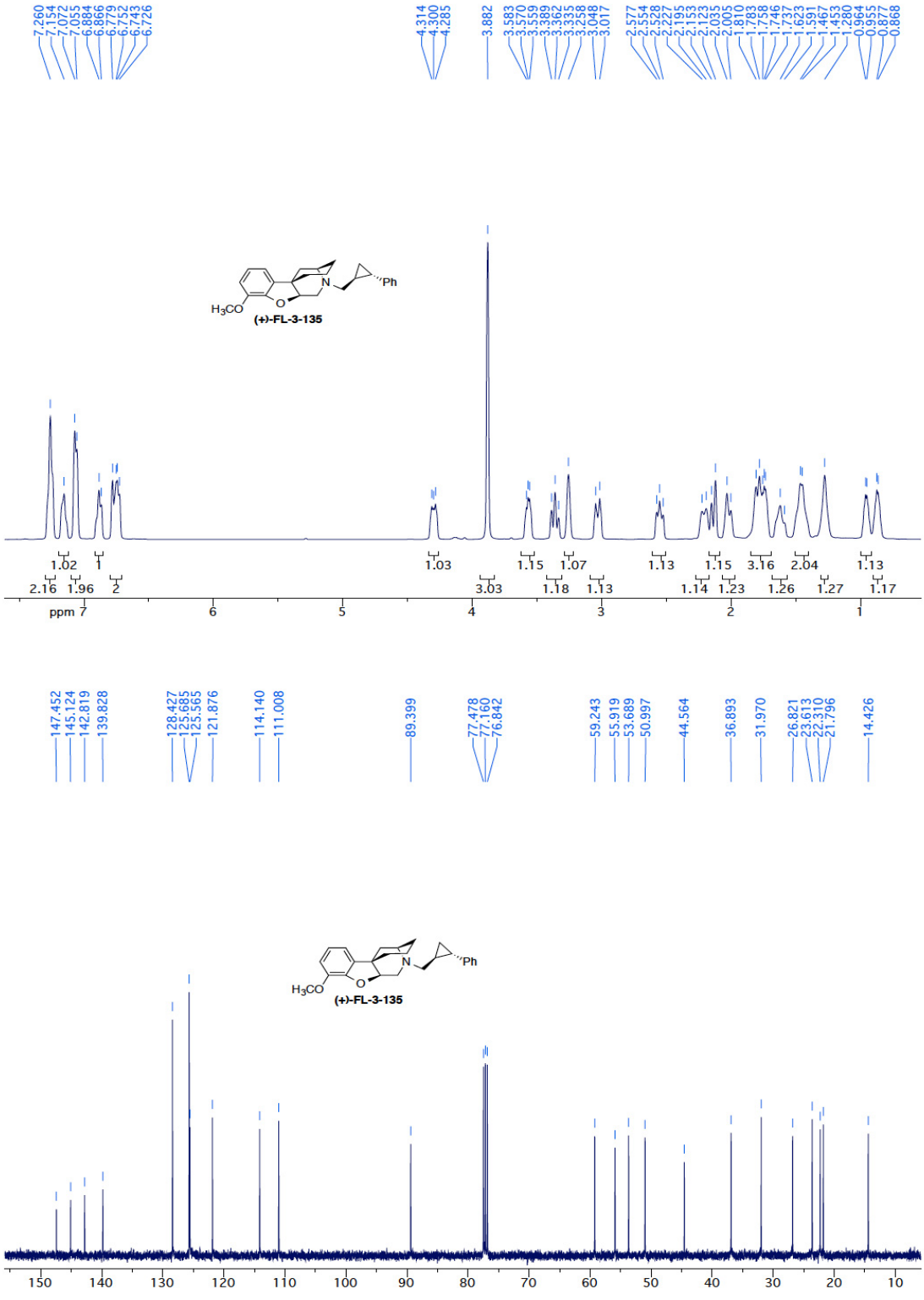


Compound (+)-**12**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7.055
6.884
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6.743
6.726
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4.300
4.285
3.882
3.583
3.570
3.559
3.389
3.362
3.335
3.258
3.048
3.017
2.577
2.554
2.528
2.227
2.195
2.153
2.123
2.035
2.005
1.810
1.783
1.758
1.746
1.737
1.623
1.591
1.467
1.453
1.280
0.964
0.955
0.877
0.868
H3CO
O
N
Ph
(+)-FL-3-135
1.02
1
1.03
1.15
1.07
1.13
1.15
3.16
2.04
1.13
2.16
1.96
2
3.03
1.18
1.13
1.14
1.23
1.26
1.27
1.17
ppm 7
6
5
4
3
2
1
147.452
145.124
142.819
139.828
128.427
125.685
125.565
121.876
114.140
111.008
89.399
77.478
77.160
76.842
59.243
55.919
53.689
50.997
44.564
36.893
31.970
26.821
23.613
22.310
21.796
14.426
H3CO
O
N
Ph
(+)-FL-3-135
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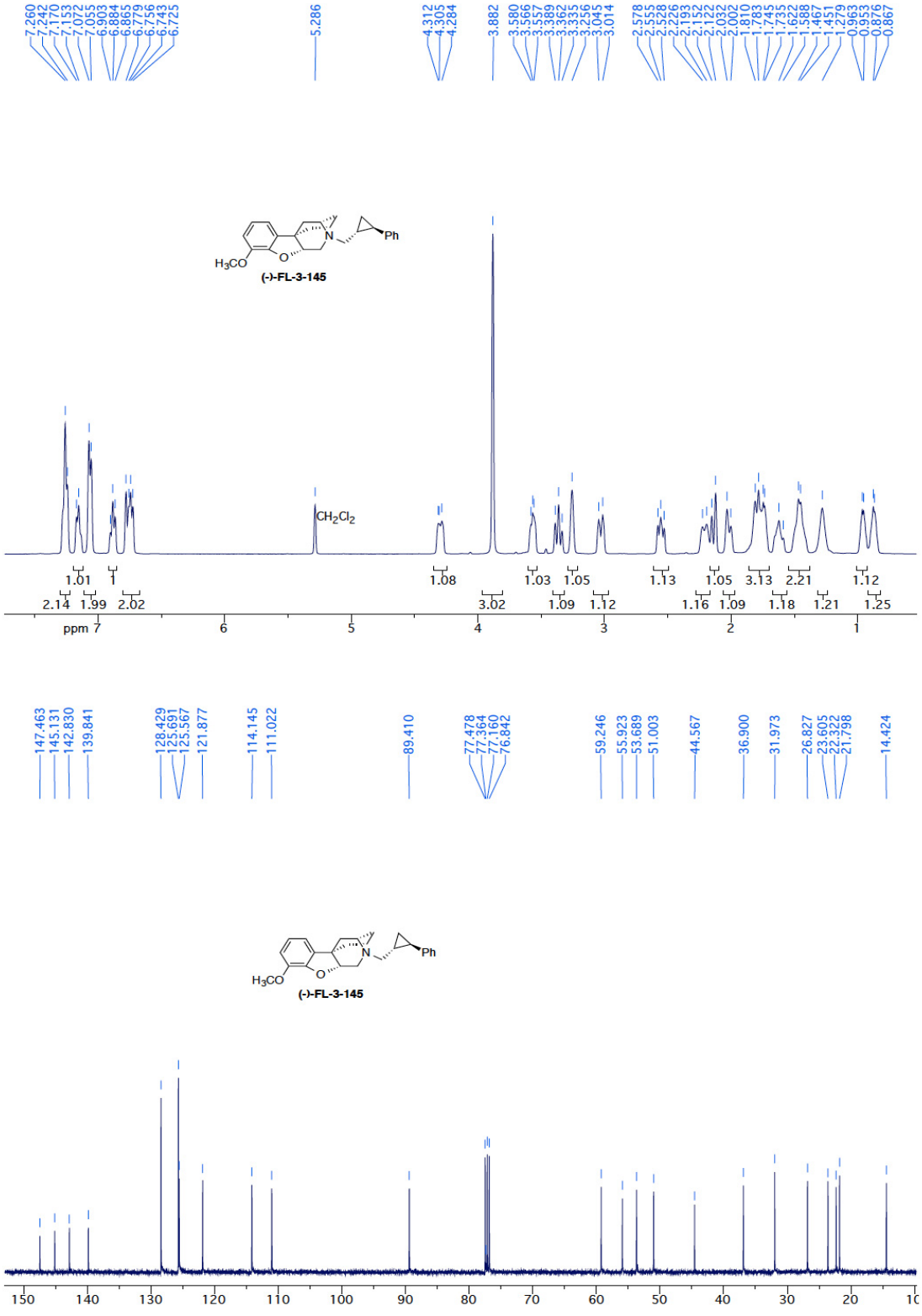
140
130
120
110
100
90
80
70
60
50
40
30
20
10

Compound (-)-**12**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7.153
7.072
7.055
6.903
6.884
6.865
6.779
6.756
6.743
6.725
5.286
4.312
4.305
4.284
3.882
3.580
3.566
3.557
3.389
3.362
3.335
3.256
3.045
3.014
2.578
2.555
2.528
2.226
2.193
2.152
2.122
2.032
2.002
1.810
1.783
1.745
1.735
1.622
1.588
1.467
1.451
1.279
0.963
0.953
0.876
0.867
H3CO
N
O
Ph
(-)-FL-3-145
CH2Cl2
1.01
1
1.08
1.03
1.05
1.13
1.05
3.13
2.21
1.12
2.14
1.99
2.02
3.02
1.09
1.12
1.16
1.09
1.18
1.21
1.25
ppm 7
6
5
4
3
2
1
147.463
145.131
142.830
139.841
128.429
125.691
125.567
121.877
114.145
111.022
89.410
77.478
77.364
77.160
76.842
59.246
55.923
53.689
51.003
44.567
36.900
31.973
26.827
23.605
22.322
21.798
14.424
H3CO
N
O
Ph
(-)-FL-3-145
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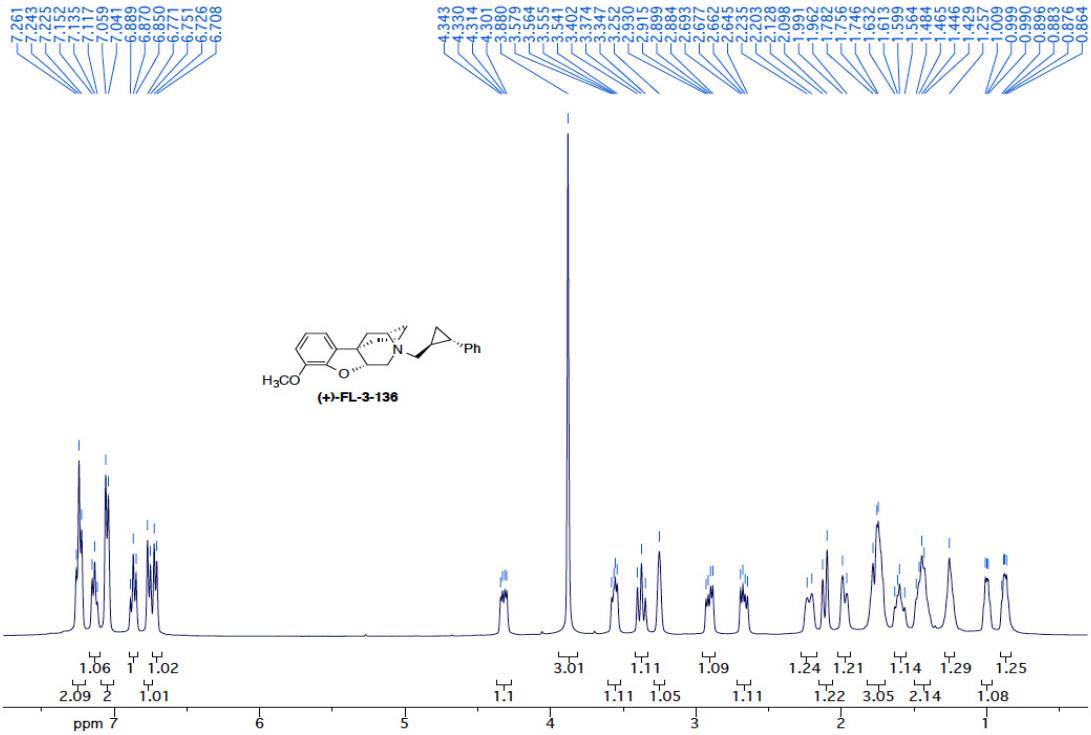
140
130
120
110
100
90
80
70
60
50
40
30
20
10

Compound (+)-**13**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7.135
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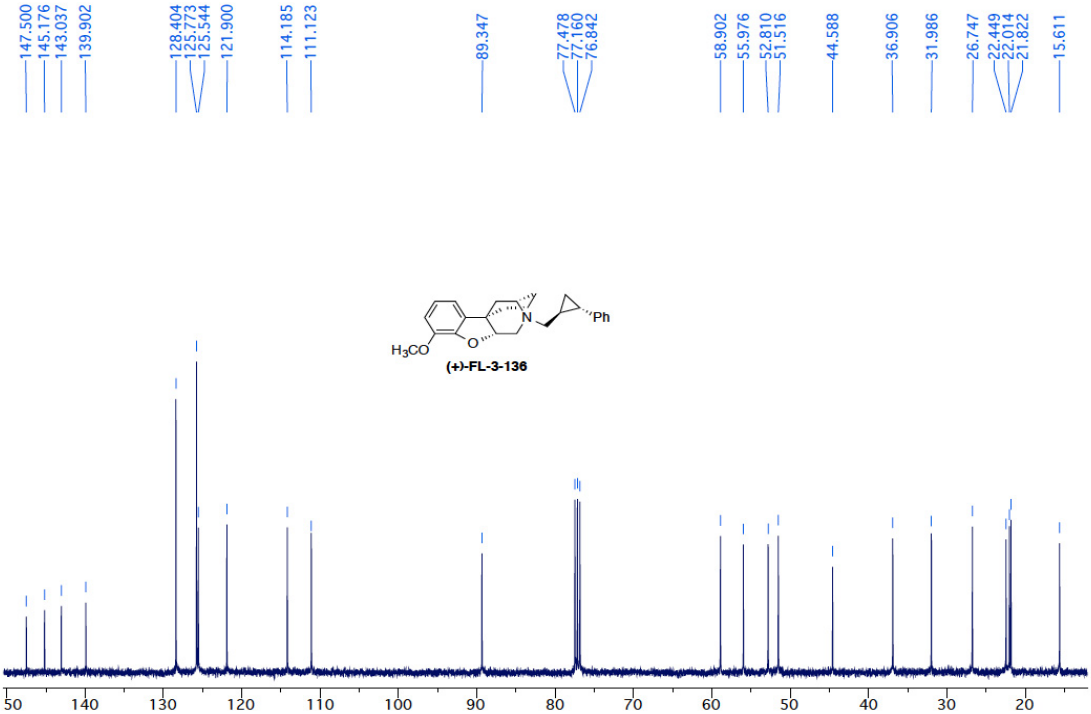
6.870
6.850
6.771
6.751
6.726
6.708
4.343
4.330
4.314
4.301
3.880
3.579
3.564
3.555
3.541
3.402
3.374
3.347
3.252
2.930
2.915
2.899
2.884
2.693
2.677
2.662
2.645
2.235
2.203
2.128
2.098
1.991
1.962
1.782
1.756
1.746
1.632
1.613
1.599
1.564
1.484
1.465
1.446
1.429
1.257
1.009
0.999
0.990
0.896
0.883
0.876
0.864
H3CO
N
O
Ph
(+)-FL-3-136
1.06 1 1.02
2.09 2 1.01
3.01
1.1
1.11 1.11 1.05
1.09 1.11
1.24 1.21 1.14 1.29 1.25
1.22 3.05 2.14 1.08
ppm 7 6 5 4 3 2 1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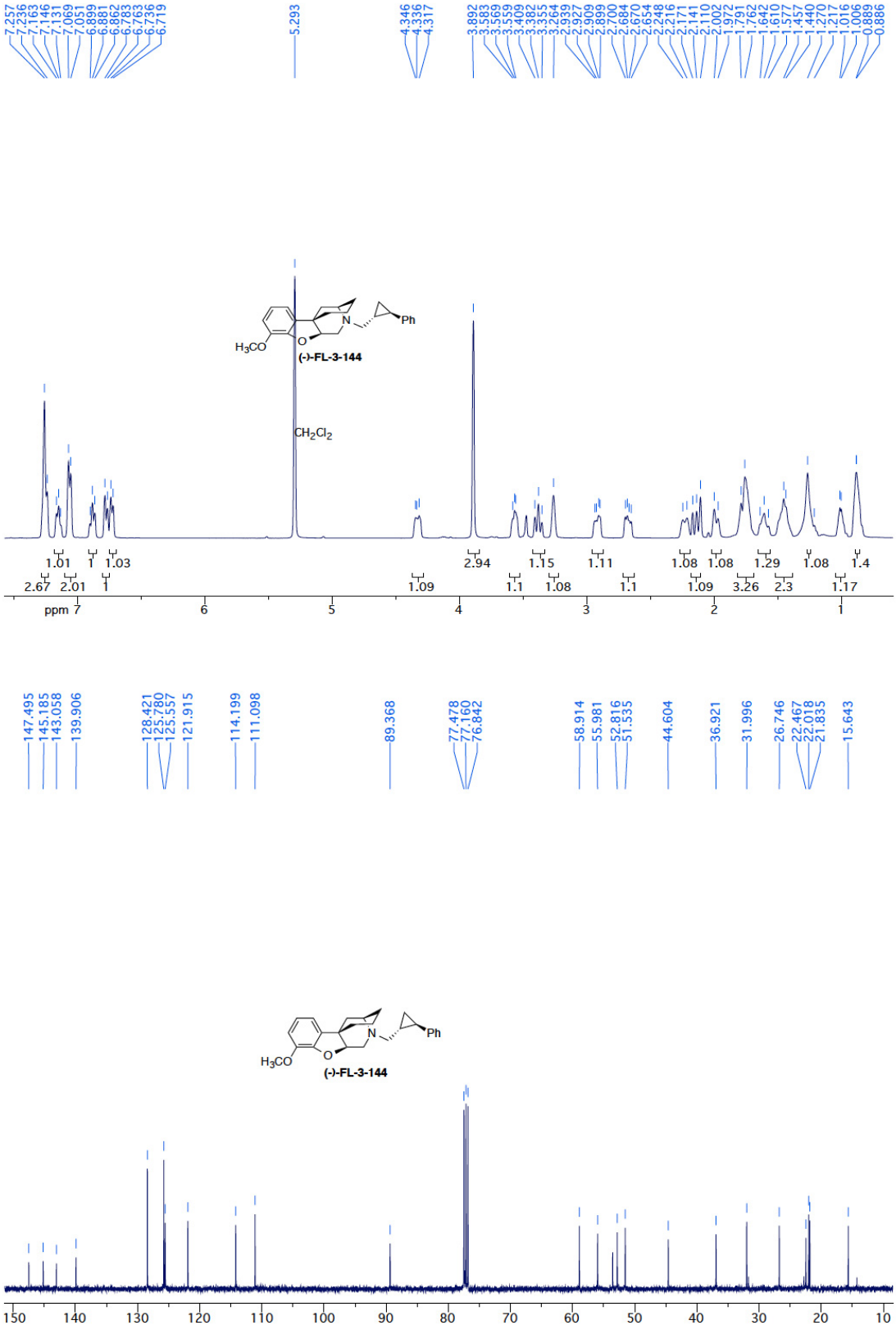
139.902
128.404
125.773
125.544
121.900
114.185
111.123
89.347
77.478
77.160
76.842
58.902
55.976
52.810
51.516
44.588
36.906
31.986
26.747
22.449
22.014
21.822
15.611
H3CO
N
O
Ph
(+)-FL-3-136
50 140 130 120 110 100 90 80 70 60 50 40 30 20

Compound (-)-**13**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7.163
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7.131
7.069
7.051
6.899
6.881
6.862
6.783
6.763
6.736
6.719
5.293
4.346
4.336
4.317
3.892
3.583
3.569
3.559
3.409
3.382
3.355
3.264
3.239
2.927
2.907
2.899
2.700
2.684
2.670
2.654
2.248
2.216
2.171
2.141
2.110
2.002
1.972
1.791
1.762
1.642
1.610
1.577
1.457
1.440
1.270
1.217
1.016
1.006
0.889
0.886
H3CO
N
O
Ph
(-)-FL-3-144
CH2Cl2
1.01
1
1.03
2.94
1.15
1.11
1.08
1.08
1.29
1.08
1.4
2.67
2.01
1
1.09
1.1
1.08
1.1
1.09
3.26
2.3
1.17
ppm 7
6
5
4
3
2
1
147.495
145.185
143.058
139.906
128.421
125.780
125.557
121.915
114.199
111.098
89.368
77.478
77.160
76.842
58.914
55.981
52.816
51.535
44.604
36.921
31.996
26.746
22.467
22.018
21.835
15.643
H3CO
N
O
Ph
(-)-FL-3-144
150
140
130
120
110
100
90
80
70
60
50
40
30
20
10

Compound (+)-**16**


7.260
7.240
7.180
7.164
7.089
7.071
6.748
6.730
6.712
6.694
6.587
6.572
4.022
3.993
3.636
3.623
3.613
3.254
3.132
3.101
3.016
2.990
2.962
2.468
2.443
2.417
2.223
2.191
2.070
2.041
1.894
1.864
1.740
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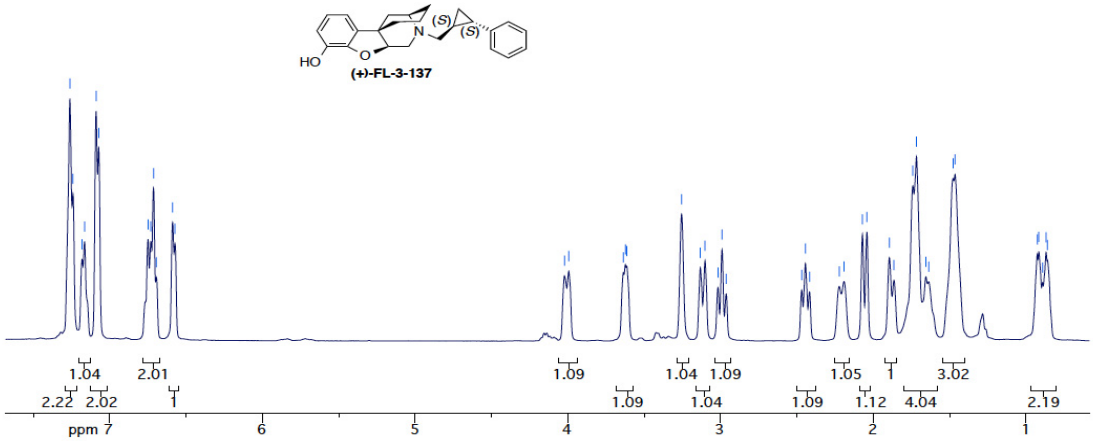
1.637
1.476
1.464
0.925
0.915
0.891
0.868
0.858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O
N
(S)
(S)
(+)-FL-3-137
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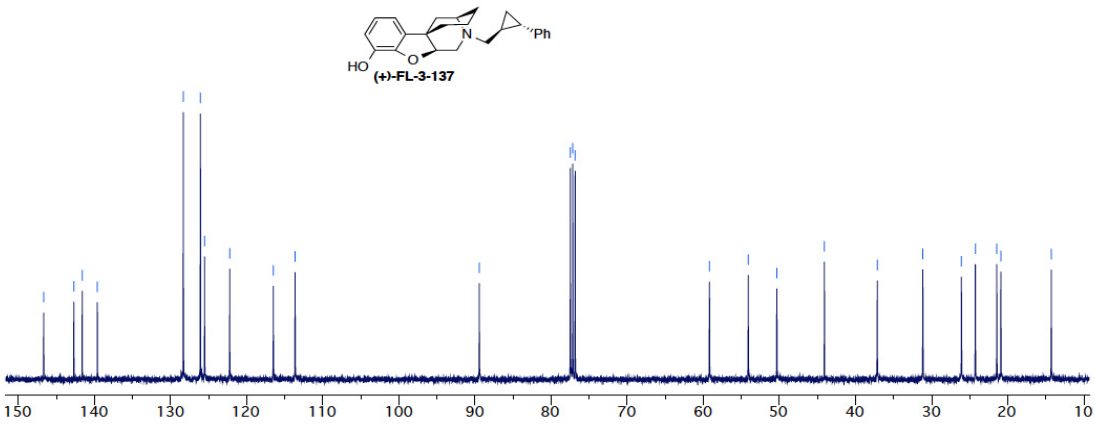
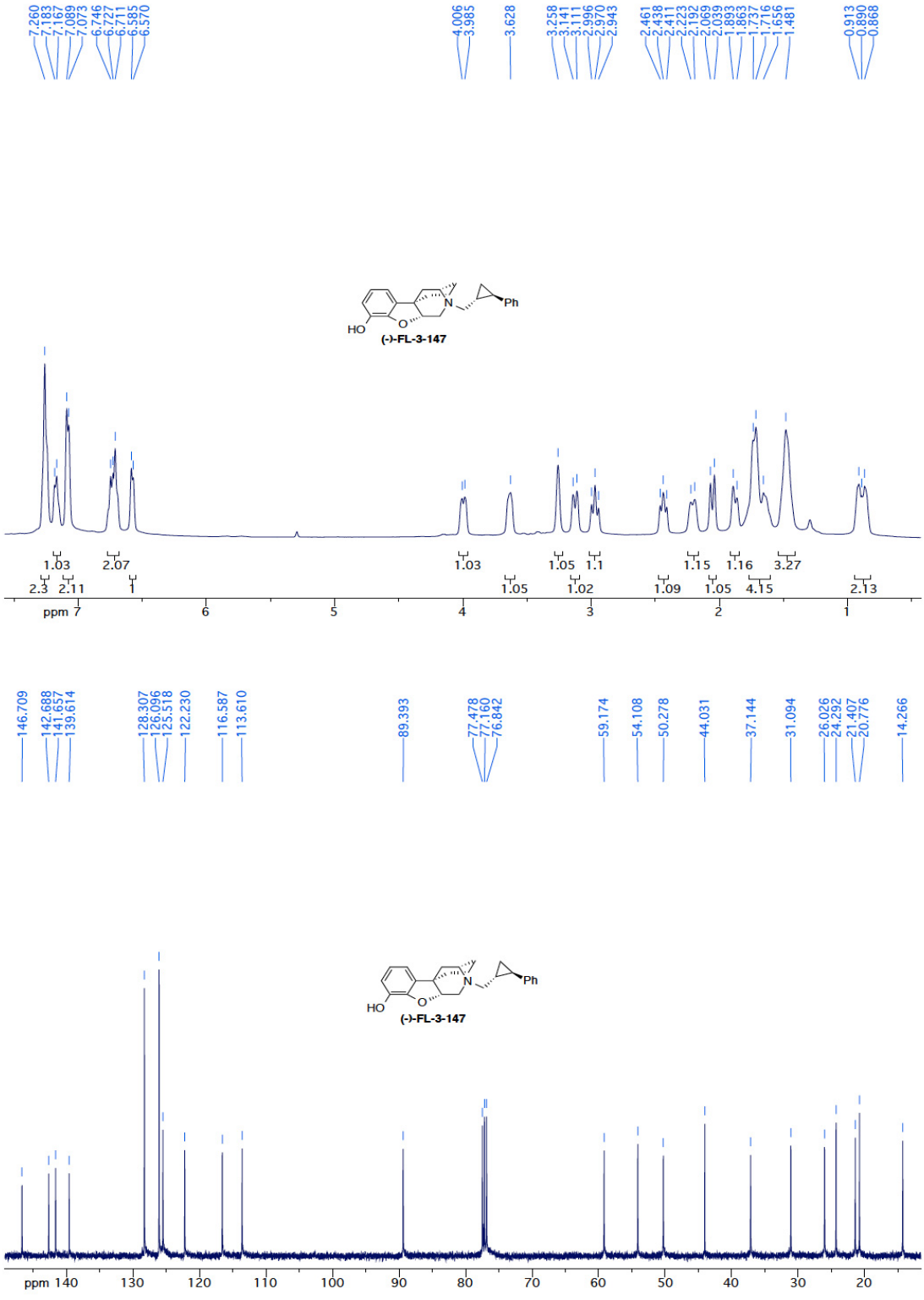
2.01
1.09
1.04
1.09
1.05
1
3.02
2.22
2.02
1
1.09
1.04
1.09
1.12
4.04
2.19
ppm 7
6
5
4
3
2
1


146.683
142.723
141.622
139.629
128.331
126.087
125.538
122.239
116.509
113.647
89.440
77.478
77.160
76.842
59.199
54.103
50.345
44.097
37.150
31.162
26.091
24.260
21.442
20.907
14.281


HO
O
N
Ph
(+)-FL-3-137
150
140
130
120
110
100
90
80
70
60
50
40
30
20
10

Compound (-)-**16**


7.260
7.183
7.167
7.089
7.073
6.746
6.727
6.711
6.585
6.570
4.006
3.985
3.628
3.258
3.141
3.111
2.996
2.970
2.943
2.461
2.438
2.411
2.223
2.192
2.069
2.039
1.893
1.863
1.737
1.716
1.656
1.481
0.913
0.890
0.868
HO
O
N
Ph
(-)-FL-3-147
1.03
2.07
2.3
2.11
1
1.03
1.05
1.1
1.05
1.02
1.15
1.16
3.27
1.09
1.05
4.15
2.13
ppm 7
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4
3
2
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125.518
122.230
116.587
113.610
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26.026
24.292
21.407
20.776
14.266
HO
O
N
Ph
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120
110
100
90
80
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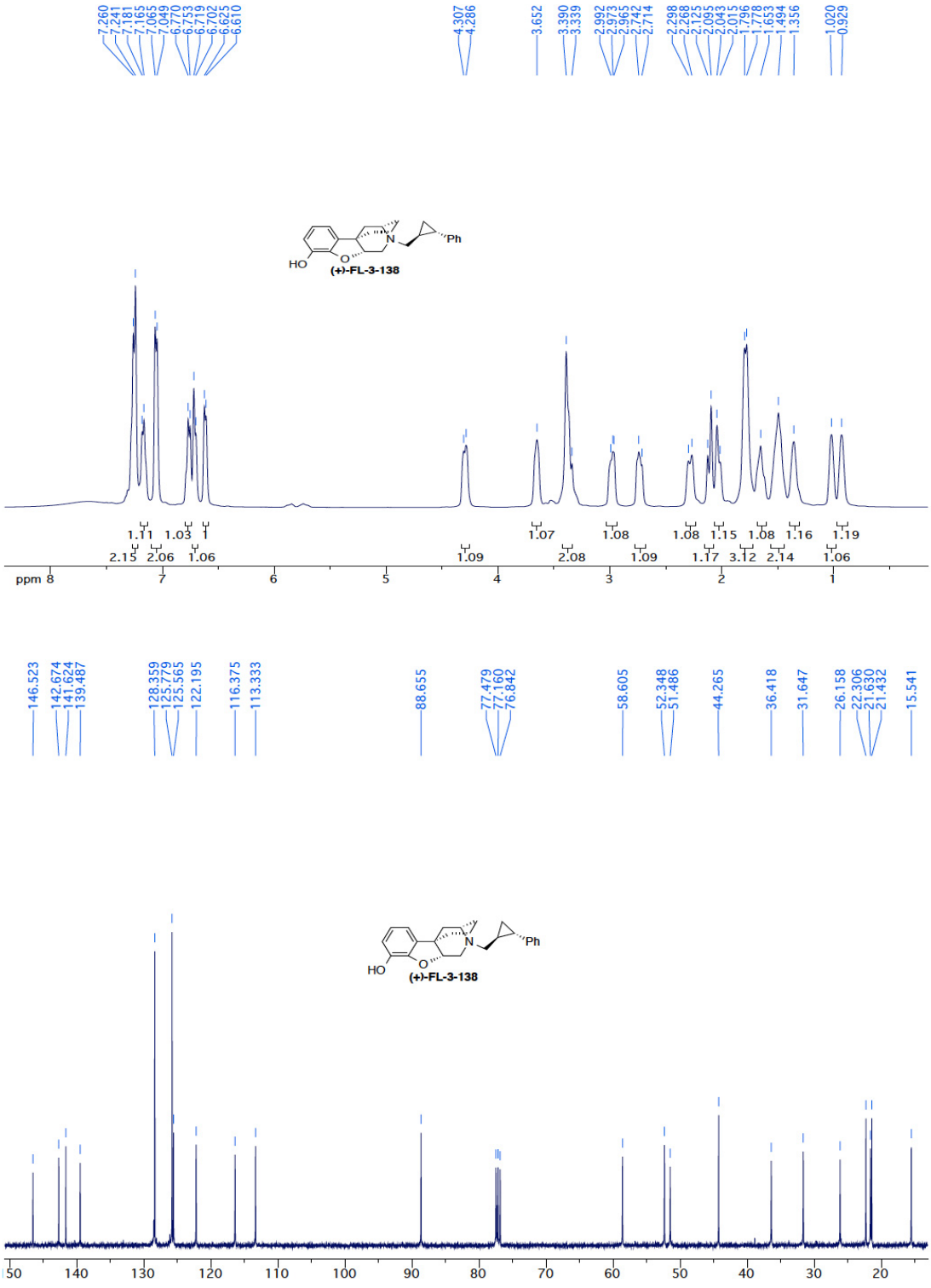
60
50
40
30
20

Compound (+)-**17**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O
N
Ph
(+)-FL-3-138
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1
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1.09
1.07
2.08
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1.09
1.08
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1.08
2.14
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1.19
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3
2
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125.565
122.195
116.375
113.333
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36.418
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26.158
22.306
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21.432
15.541
HO
O
N
Ph
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130
120
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100
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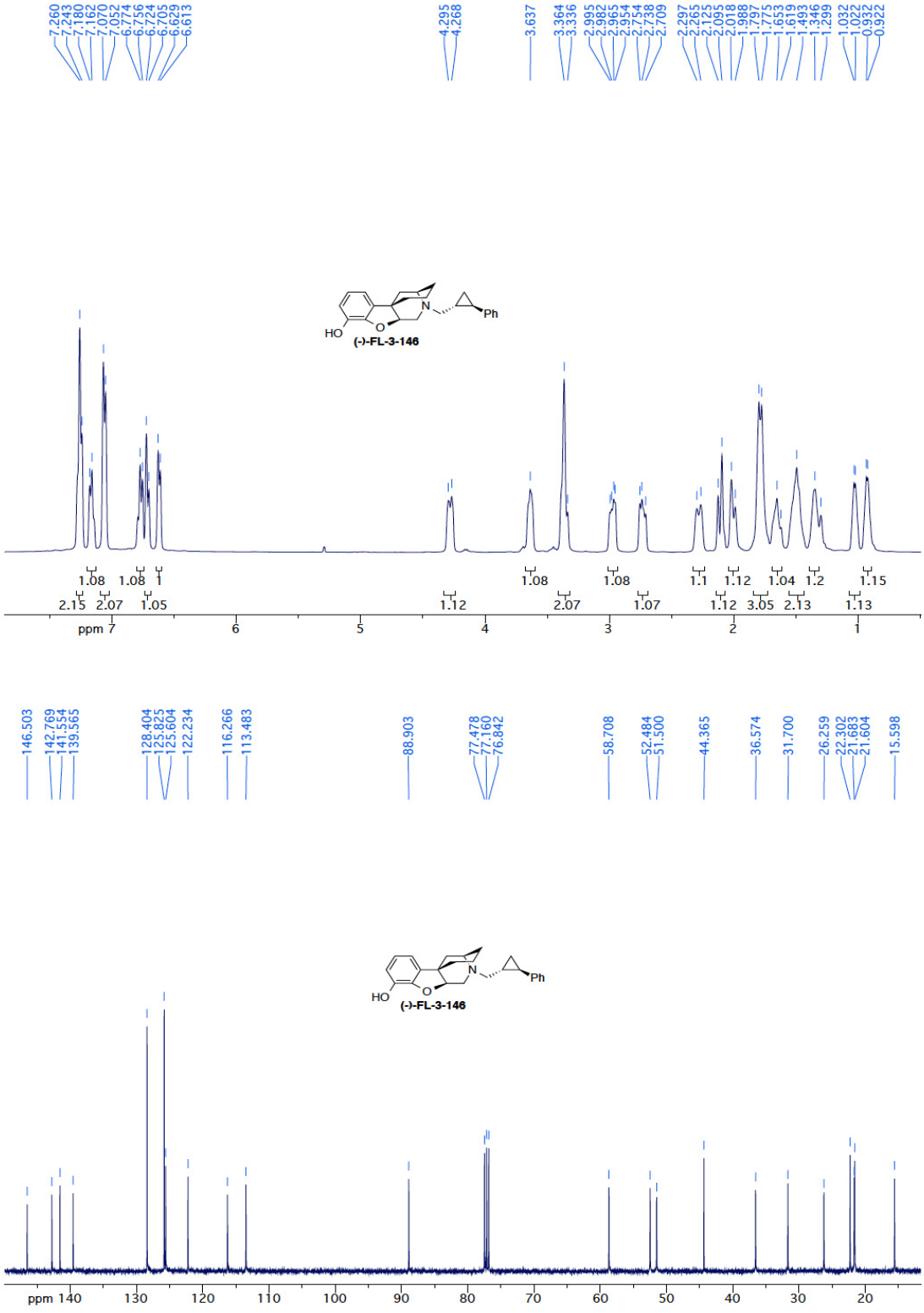
60
50
40
30
20

Compound (-)-**17**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1.299
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1.15
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2.07
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2.13
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88.903
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77.160
76.842
58.708
52.484
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44.365
36.574
31.700
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22.302
21.683
21.604
15.598
HO
O
N
Ph
(-)-FL-3-146
ppm 140
130
120
110
100
90
80
70
60
50
40
30
20

Table S1. Crystal data and structure refinement for **(-)-6**

Identification code	knih103	
Empirical formula	$C_{24}H_{31}NO_6$	
Formula weight	429.50	
Temperature	293(2) K	
Wavelength	1.54178 Å	
Crystal system	Orthorhombic	
Space group	$P2_12_12_1$	
Unit cell dimensions	a = 6.7324(6) Å	⟨= 90°.
	b = 9.4370(6) Å	®= 90°.
	c = 34.612(4) Å	© = 90°.
Volume	2199.0(3) $Å^3$	
Z	4	
Density (20°C)	1.297 Mg/m^3	
Absorption coefficient	0.759 mm^{-1}	
F(000)	920	
Crystal size	0.211 x 0.102 x 0.040 mm^3	
Theta range for data collection	2.553 to 74.508°.	
Index ranges	-6<=h<=8, -11<=k<=10, -40<=l<=42	
Reflections collected	16601	
Independent reflections	4436 [R_{int} = 0.0558]	
Completeness to theta = 25.000°	98.8 %	
Absorption correction	Semi-empirical from equivalents	
Max. and min. transmission	0.7538 and 0.5283	
Refinement method	Full-matrix least-squares on F^2	
Data / restraints / parameters	4436 / 3 / 291	
Goodness-of-fit on F^2	1.091	
Final R indices [I>2sigma(I)]	R_1 = 0.0466, wR_2 = 0.1120	
R indices (all data)	R_1 = 0.0480, wR_2 = 0.1137	
Absolute structure parameter	0.10(8)	
Largest diff. peak and hole	0.298 and -0.388 $e.Å^{-3}$	

Table S2. Atomic coordinates (x 10^4) and equivalent isotropic displacement parameters ($Å^2$x 10^3) for **(-)-6**. U(eq) is defined as one third of the trace of the orthogonalized U^{ij} tensor.

	x	y	z	U(eq)
C(1)	3200(3)	4745(2)	3486(1)	35(1)
N(2)	3254(3)	5165(2)	3067(1)	35(1)
C(3)	3074(3)	6735(2)	2972(1)	38(1)
C(3A)	4569(3)	7609(2)	3206(1)	37(1)
C(4)	966(4)	7266(3)	3039(1)	48(1)
C(5)	362(3)	7318(3)	3464(1)	49(1)
C(6)	2024(3)	7758(2)	3739(1)	38(1)
C(6A)	4167(3)	7304(2)	3631(1)	29(1)
C(6B)	5542(3)	7847(2)	3946(1)	32(1)
C(7)	6423(3)	9149(2)	4017(1)	41(1)
C(8)	7506(4)	9301(3)	4354(1)	48(1)
C(9)	7648(3)	8205(3)	4623(1)	46(1)
C(10)	6713(3)	6909(2)	4559(1)	36(1)
O(10)	6645(3)	5783(2)	4807(1)	46(1)
C(10A)	7629(5)	5932(4)	5167(1)	70(1)
C(11)	5708(3)	6754(2)	4211(1)	31(1)
O(11)	4768(2)	5518(2)	4097(1)	35(1)
C(11A)	4616(3)	5733(2)	3685(1)	29(1)
C(1')	-1562(3)	3347(2)	2951(1)	31(1)
O(1')	200(2)	3430(2)	2839(1)	39(1)
O(1'')	-2911(3)	4195(2)	2871(1)	47(1)
C(2')	-2045(3)	2040(2)	3210(1)	37(1)
O(2')	-4099(3)	1862(2)	3280(1)	49(1)
C(3')	-991(3)	2215(2)	3594(1)	34(1)
C(4')	886(3)	1654(2)	3659(1)	39(1)
C(5')	1885(4)	1908(2)	4002(1)	44(1)
C(6')	1054(4)	2736(3)	4291(1)	45(1)
C(7')	-850(4)	3260(3)	4230(1)	47(1)
C(8')	-1859(4)	3002(2)	3889(1)	42(1)
C(9')	2176(6)	3086(4)	4655(1)	71(1)
O(1S)	3135(3)	1293(3)	2704(1)	62(1)

Table S3. Bond lengths [Å] and angles [°] for **(-)-6**

.

C(1)-C(11A)	1.501(3)	C(1)-N(2)	1.507(3)
C(1)-H(1A)	0.9700	C(1)-H(1B)	0.9700
N(2)-C(3)	1.522(3)	N(2)-H(2A)	0.8900
N(2)-H(2B)	0.8900	C(3)-C(4)	1.523(3)
C(3)-C(3A)	1.533(3)	C(3)-H(3)	0.9800
C(3A)-C(6A)	1.523(3)	C(3A)-H(3A)	0.9700
C(3A)-H(3B)	0.9700	C(4)-C(5)	1.526(4)
C(4)-H(4A)	0.9700	C(4)-H(4B)	0.9700
C(5)-C(6)	1.528(3)	C(5)-H(5A)	0.9700
C(5)-H(5B)	0.9700	C(6)-C(6A)	1.551(3)
C(6)-H(6A)	0.9700	C(6)-H(6B)	0.9700
C(6A)-C(6B)	1.518(3)	C(6A)-C(11A)	1.525(3)
C(6B)-C(7)	1.386(3)	C(6B)-C(11)	1.386(3)
C(7)-C(8)	1.385(4)	C(7)-H(7)	0.9300
C(8)-C(9)	1.395(4)	C(8)-H(8)	0.9300
C(9)-C(10)	1.393(3)	C(9)-H(9)	0.9300
C(10)-O(10)	1.367(3)	C(10)-C(11)	1.390(3)
O(10)-C(10A)	1.419(3)	C(10A)-H(10A)	0.9600
C(10A)-H(10B)	0.9600	C(10A)-H(10C)	0.9600
C(11)-O(11)	1.384(2)	O(11)-C(11A)	1.443(2)
C(11A)-H(11A)	0.9800	C(1')-O(1")	1.242(3)
C(1')-O(1')	1.251(3)	C(1')-C(2')	1.558(3)
C(2')-O(2')	1.414(3)	C(2')-C(3')	1.517(3)
C(2')-H(2')	0.9800	O(2')-H(2C)	0.85(4)
C(3')-C(4')	1.388(3)	C(3')-C(8')	1.392(3)
C(4')-C(5')	1.386(3)	C(4')-H(4')	0.9300
C(5')-C(6')	1.387(4)	C(5')-H(5')	0.9300
C(6')-C(7')	1.390(4)	C(6')-C(9')	1.508(4)
C(7')-C(8')	1.383(4)	C(7')-H(7')	0.9300
C(8')-H(8')	0.9300	C(9')-H(9D)	0.9600
C(9')-H(9E)	0.9600	C(9')-H(9F)	0.9600
O(1S)-H(1S)	0.8199(14)	O(1S)-H(2S)	0.8201(14)

C(11A)-C(1)-N(2)	105.32(17)	C(11A)-C(1)-H(1A)	110.7
N(2)-C(1)-H(1A)	110.7	C(11A)-C(1)-H(1B)	110.7
N(2)-C(1)-H(1B)	110.7	H(1A)-C(1)-H(1B)	108.8
C(1)-N(2)-C(3)	117.49(16)	C(1)-N(2)-H(2A)	107.9
C(3)-N(2)-H(2A)	107.9	C(1)-N(2)-H(2B)	107.9
C(3)-N(2)-H(2B)	107.9	H(2A)-N(2)-H(2B)	107.2
N(2)-C(3)-C(4)	111.2(2)	N(2)-C(3)-C(3A)	110.98(17)
C(4)-C(3)-C(3A)	110.75(19)	N(2)-C(3)-H(3)	107.9
C(4)-C(3)-H(3)	107.9	C(3A)-C(3)-H(3)	107.9
C(6A)-C(3A)-C(3)	106.95(17)	C(6A)-C(3A)-H(3A)	110.3
C(3)-C(3A)-H(3A)	110.3	C(6A)-C(3A)-H(3B)	110.3

Table S3. (continued).

C(3)-C(3A)-H(3B)	110.3	H(3A)-C(3A)-H(3B)	108.6
C(3)-C(4)-C(5)	113.90(19)	C(3)-C(4)-H(4A)	108.8
C(5)-C(4)-H(4A)	108.8	C(3)-C(4)-H(4B)	108.8
C(5)-C(4)-H(4B)	108.8	H(4A)-C(4)-H(4B)	107.7
C(4)-C(5)-C(6)	114.5(2)	C(4)-C(5)-H(5A)	108.6
C(6)-C(5)-H(5A)	108.6	C(4)-C(5)-H(5B)	108.6
C(6)-C(5)-H(5B)	108.6	H(5A)-C(5)-H(5B)	107.6
C(5)-C(6)-C(6A)	117.08(18)	C(5)-C(6)-H(6A)	108.0
C(6A)-C(6)-H(6A)	108.0	C(5)-C(6)-H(6B)	108.0
C(6A)-C(6)-H(6B)	108.0	H(6A)-C(6)-H(6B)	107.3
C(6B)-C(6A)-C(3A)	121.34(17)	C(6B)-C(6A)-C(11A)	96.83(15)
C(3A)-C(6A)-C(11A)	105.57(17)	C(6B)-C(6A)-C(6)	107.49(16)
C(3A)-C(6A)-C(6)	110.25(17)	C(11A)-C(6A)-C(6)	115.05(17)
C(7)-C(6B)-C(11)	120.5(2)	C(7)-C(6B)-C(6A)	133.5(2)
C(11)-C(6B)-C(6A)	105.86(17)	C(8)-C(7)-C(6B)	117.9(2)
C(8)-C(7)-H(7)	121.1	C(6B)-C(7)-H(7)	121.1
C(7)-C(8)-C(9)	121.5(2)	C(7)-C(8)-H(8)	119.3
C(9)-C(8)-H(8)	119.3	C(10)-C(9)-C(8)	120.9(2)
C(10)-C(9)-H(9)	119.5	C(8)-C(9)-H(9)	119.5
O(10)-C(10)-C(11)	116.51(18)	O(10)-C(10)-C(9)	126.7(2)
C(11)-C(10)-C(9)	116.8(2)	C(10)-O(10)-C(10A)	117.3(2)
O(10)-C(10A)-H(10A)	109.5	O(10)-C(10A)-H(10B)	109.5
H(10A)-C(10A)-H(10B)	109.5	O(10)-C(10A)-H(10C)	109.5
H(10A)-C(10A)-H(10C)	109.5	H(10B)-C(10A)-H(10C)	109.5
O(11)-C(11)-C(6B)	113.68(17)	O(11)-C(11)-C(10)	123.91(18)
C(6B)-C(11)-C(10)	122.40(19)	C(11)-O(11)-C(11A)	101.25(15)
O(11)-C(11A)-C(1)	114.29(16)	O(11)-C(11A)-C(6A)	105.82(15)
C(1)-C(11A)-C(6A)	114.90(17)	O(11)-C(11A)-H(11A)	107.1
C(1)-C(11A)-H(11A)	107.1	C(6A)-C(11A)-H(11A)	107.1
O(1")-C(1')-O(1')	125.7(2)	O(1")-C(1')-C(2')	119.09(18)
O(1')-C(1')-C(2')	115.23(19)	O(2')-C(2')-C(3')	108.64(19)
O(2')-C(2')-C(1')	113.41(18)	C(3')-C(2')-C(1')	108.72(16)
O(2')-C(2')-H(2')	108.7	C(3')-C(2')-H(2')	108.7
C(1')-C(2')-H(2')	108.7	C(2')-O(2')-H(2C)	107(3)
C(4')-C(3')-C(8')	117.9(2)	C(4')-C(3')-C(2')	121.7(2)
C(8')-C(3')-C(2')	120.4(2)	C(5')-C(4')-C(3')	120.9(2)
C(5')-C(4')-H(4')	119.5	C(3')-C(4')-H(4')	119.5
C(4')-C(5')-C(6')	121.3(2)	C(4')-C(5')-H(5')	119.4
C(6')-C(5')-H(5')	119.4	C(5')-C(6')-C(7')	117.7(2)
C(5')-C(6')-C(9')	121.7(3)	C(7')-C(6')-C(9')	120.6(3)
C(8')-C(7')-C(6')	121.2(2)	C(8')-C(7')-H(7')	119.4
C(6')-C(7')-H(7')	119.4	C(7')-C(8')-C(3')	120.9(2)
C(7')-C(8')-H(8')	119.5	C(3')-C(8')-H(8')	119.5

Table S3. (continued).

C(6')-C(9')-H(9D)	109.5	C(6')-C(9')-H(9E)	109.5
H(9D)-C(9')-H(9E)	109.5	C(6')-C(9')-H(9F)	109.5
H(9D)-C(9')-H(9F)	109.5	H(9E)-C(9')-H(9F)	109.5
H(1S)-O(1S)-H(2S)	104.5(2)		

Table S4. Anisotropic displacement parameters ($Å^2 x\ 10^3$) for **(-)-6**. The anisotropic displacement factor exponent takes the form: $-2\pi^2[h^2 a^{*2} U^{11} + ... + 2\ h\ k\ a^*\ b^*\ U^{12}]$

	U^{11}	U^{22}	U^{33}	U^{23}	U^{13}	U^{12}
C(1)	38(1)	32(1)	37(1)	-1(1)	-6(1)	-4(1)
N(2)	29(1)	43(1)	34(1)	-7(1)	-1(1)	-1(1)
C(3)	39(1)	45(1)	31(1)	5(1)	0(1)	-1(1)
C(3A)	39(1)	37(1)	36(1)	7(1)	4(1)	-4(1)
C(4)	43(1)	56(1)	45(1)	4(1)	-12(1)	11(1)
C(5)	31(1)	64(2)	52(1)	-3(1)	-2(1)	13(1)
C(6)	36(1)	38(1)	42(1)	-3(1)	2(1)	7(1)
C(6A)	29(1)	26(1)	33(1)	2(1)	1(1)	-1(1)
C(6B)	30(1)	29(1)	37(1)	-3(1)	2(1)	-1(1)
C(7)	39(1)	31(1)	52(1)	-3(1)	6(1)	-7(1)
C(8)	41(1)	41(1)	63(2)	-13(1)	1(1)	-14(1)
C(9)	38(1)	51(1)	49(1)	-15(1)	-6(1)	-6(1)
C(10)	32(1)	41(1)	35(1)	-5(1)	-2(1)	1(1)
O(10)	53(1)	50(1)	36(1)	1(1)	-10(1)	0(1)
C(10A)	74(2)	82(2)	53(2)	10(2)	-28(2)	-6(2)
C(11)	28(1)	31(1)	35(1)	-4(1)	-1(1)	-2(1)
O(11)	43(1)	28(1)	32(1)	2(1)	-7(1)	-4(1)
C(11A)	31(1)	26(1)	32(1)	0(1)	-1(1)	1(1)
C(1')	31(1)	34(1)	29(1)	-4(1)	-3(1)	-2(1)
O(1')	34(1)	46(1)	38(1)	-4(1)	0(1)	-3(1)
O(1'')	41(1)	49(1)	51(1)	15(1)	7(1)	10(1)
C(2')	40(1)	32(1)	39(1)	-1(1)	-4(1)	-3(1)
O(2')	42(1)	60(1)	45(1)	5(1)	-7(1)	-18(1)
C(3')	39(1)	27(1)	37(1)	4(1)	-3(1)	-1(1)
C(4')	40(1)	35(1)	41(1)	5(1)	0(1)	4(1)
C(5')	39(1)	44(1)	48(1)	13(1)	-6(1)	3(1)
C(6')	55(1)	40(1)	39(1)	9(1)	-10(1)	-7(1)
C(7')	61(2)	42(1)	39(1)	-3(1)	-4(1)	6(1)
C(8')	44(1)	39(1)	43(1)	0(1)	-4(1)	11(1)
C(9')	84(2)	77(2)	50(2)	8(1)	-25(2)	-15(2)
O(1S)	43(1)	83(1)	62(1)	-31(1)	-15(1)	14(1)

Table S5. Hydrogen coordinates (x 10^4) and isotropic displacement parameters (Å^{2}x 10^3) for **(-)-6**.

	x	y	z	U(eq)
H(1A)	1870	4849	3590	42
H(1B)	3619	3769	3518	42
H(2A)	2274	4712	2946	42
H(2B)	4390	4852	2967	42
H(3)	3385	6858	2697	46
H(3A)	5917	7341	3139	45
H(3B)	4401	8611	3153	45
H(4A)	50	6653	2901	57
H(4B)	845	8209	2930	57
H(5A)	-115	6388	3539	58
H(5B)	-733	7978	3492	58
H(6A)	1722	7374	3992	46
H(6B)	2001	8782	3762	46
H(7)	6291	9896	3843	49
H(8)	8153	10153	4403	58
H(9)	8376	8342	4849	55
H(10A)	9034	6015	5125	104
H(10B)	7367	5116	5325	104
H(10C)	7154	6767	5296	104
H(11A)	5936	5559	3576	35
H(2')	-1538	1186	3083	44
H(2C)	-4640(60)	1690(40)	3062(12)	74
H(4')	1480	1100	3469	46
H(5')	3138	1516	4039	52
H(7')	-1456	3795	4423	57
H(8')	-3136	3359	3857	50
H(9D)	2006	4072	4715	106
H(9E)	3562	2888	4619	106
H(9F)	1675	2522	4864	106
H(1S)	2710(50)	650(20)	2569(9)	94
H(2S)	2220(30)	1860(20)	2717(12)	94

Table S6. Torsion angles [°] for **(-)-6**.

C(11A)-C(1)-N(2)-C(3)	-47.6(2)	C(1)-N(2)-C(3)-C(4)	-72.5(2)
C(1)-N(2)-C(3)-C(3A)	51.3(3)	N(2)-C(3)-C(3A)-C(6A)	-57.3(2)
C(4)-C(3)-C(3A)-C(6A)	66.7(2)	N(2)-C(3)-C(4)-C(5)	69.3(3)
C(3A)-C(3)-C(4)-C(5)	-54.6(3)	C(3)-C(4)-C(5)-C(6)	36.8(3)
C(4)-C(5)-C(6)-C(6A)	-33.1(3)	C(3)-C(3A)-C(6A)-C(6B)	172.16(18)
C(3)-C(3A)-C(6A)-C(11A)	63.9(2)	C(3)-C(3A)-C(6A)-C(6)	-61.0(2)
C(5)-C(6)-C(6A)-C(6B)	-179.6(2)	C(5)-C(6)-C(6A)-C(3A)	46.1(3)
C(5)-C(6)-C(6A)-C(11A)	-73.1(3)	C(3A)-C(6A)-C(6B)-C(7)	46.1(3)
C(11A)-C(6A)-C(6B)-C(7)	159.0(2)	C(6)-C(6A)-C(6B)-C(7)	-82.0(3)
C(3A)-C(6A)-C(6B)-C(11)	-139.50(19)	C(11A)-C(6A)-C(6B)-C(11)	-26.59(19)
C(6)-C(6A)-C(6B)-C(11)	92.42(19)	C(11)-C(6B)-C(7)-C(8)	1.7(3)
C(6A)-C(6B)-C(7)-C(8)	175.4(2)	C(6B)-C(7)-C(8)-C(9)	-2.6(4)
C(7)-C(8)-C(9)-C(10)	0.5(4)	C(8)-C(9)-C(10)-O(10)	-176.2(2)
C(8)-C(9)-C(10)-C(11)	2.3(3)	C(11)-C(10)-O(10)-C(10A)	-178.8(2)
C(9)-C(10)-O(10)-C(10A)	-0.3(4)	C(7)-C(6B)-C(11)-O(11)	-179.36(19)
C(6A)-C(6B)-C(11)-O(11)	5.4(2)	C(7)-C(6B)-C(11)-C(10)	1.3(3)
C(6A)-C(6B)-C(11)-C(10)	-174.00(19)	O(10)-C(10)-C(11)-O(11)	-3.9(3)
C(9)-C(10)-C(11)-O(11)	177.47(19)	O(10)-C(10)-C(11)-C(6B)	175.42(19)
C(9)-C(10)-C(11)-C(6B)	-3.2(3)	C(6B)-C(11)-O(11)-C(11A)	20.1(2)
C(10)-C(11)-O(11)-C(11A)	-160.52(19)	C(11)-O(11)-C(11A)-C(1)	-165.13(17)
C(11)-O(11)-C(11A)-C(6A)	-37.68(19)	N(2)-C(1)-C(11A)-O(11)	179.24(16)
N(2)-C(1)-C(11A)-C(6A)	56.6(2)	C(6B)-C(6A)-C(11A)-O(11)	39.56(18)
C(3A)-C(6A)-C(11A)-O(11)	164.81(16)	C(6)-C(6A)-C(11A)-O(11)	-73.4(2)
C(6B)-C(6A)-C(11A)-C(1)	166.64(17)	C(3A)-C(6A)-C(11A)-C(1)	-68.1(2)
C(6)-C(6A)-C(11A)-C(1)	53.7(2)	O(1")-C(1')-C(2')-O(2')	8.8(3)
O(1')-C(1')-C(2')-O(2')	-170.61(19)	O(1")-C(1')-C(2')-C(3')	-112.2(2)
O(1')-C(1')-C(2')-C(3')	68.4(2)	O(2')-C(2')-C(3')-C(4')	143.6(2)
C(1')-C(2')-C(3')-C(4')	-92.5(2)	O(2')-C(2')-C(3')-C(8')	-39.1(3)
C(1')-C(2')-C(3')-C(8')	84.8(2)	C(8')-C(3')-C(4')-C(5')	-2.0(3)
C(2')-C(3')-C(4')-C(5')	175.4(2)	C(3')-C(4')-C(5')-C(6')	-0.4(3)
C(4')-C(5')-C(6')-C(7')	2.3(3)	C(4')-C(5')-C(6')-C(9')	-176.5(2)
C(5')-C(6')-C(7')-C(8')	-1.9(4)	C(9')-C(6')-C(7')-C(8')	176.9(3)
C(6')-C(7')-C(8')-C(3')	-0.5(4)	C(4')-C(3')-C(8')-C(7')	2.4(3)
C(2')-C(3')-C(8')-C(7')	-175.0(2)		

Table S7. Hydrogen bonds for **(-)-6** [Å and °].

D-H...A	d(D-H)	d(H...A)	d(D...A)	<(DHA)
C(1)-H(1B)...O(2')#1	0.97	2.51	3.349(3)	145.4
N(2)-H(2A)...O(1')	0.89	1.88	2.744(2)	161.8
N(2)-H(2B)...O(1")#1	0.89	1.95	2.822(2)	166.8
C(9)-H(9)...O(10)#2	0.93	2.63	3.469(3)	149.6
O(2')-H(2C)...O(1S)#3	0.85(4)	1.98(4)	2.781(3)	156(4)
O(1S)-H(1S)...O(1")#4	0.8199(14)	2.052(17)	2.810(3)	154(4)
O(1S)-H(2S)...O(1')	0.8201(14)	2.053(7)	2.862(3)	169(3)

Symmetry transformations used to generate equivalent atoms:
#1 x+1,y,z #2 x+1/2,-y+3/2,-z+1 #3 x-1,y,z
#4 -x,y-1/2,-z+1/2