Supporting Information

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Table S1 ^{1}H NMR Spectroscopic Data (500 MHz, $CDCl_3$) for Compounds **9-14** (δ_H in ppm, J in Hz)

no.	**9**	**10**	**11**	**12**	**13**	**14**
2	5.76, s	5.78, s	5.80, s	5.80, s	5.80, s	5.79, s
4	4.85, t (5.0)	4.87, t (5.0)	4.88, t (5.0)	4.88, t (5.0)	4.87, t (4.5)	4.87, t (5.0)
5	2.62, m[a]	2.65, m	2.68, m	2.68, m	2.68, m	2.67, m
6	5.02, t (7.0)	5.00, t (7.5)	5.06, t (7.5)	5.05, t (7.0)	5.03, t (3.0)	5.05, t (7.0)
8	1.97, m[a]	1.97, m[a]	2.02, m	2.02, m	2.05, m[a]	2.05, m[a]
9	1.54, m[a]	1.54, m[a]	1.54, m	1.54, m	1.68, m[a]	1.68, m[a]
10a	1.83, m[a]	1.83, m[a]	1.83, m	1.81, m	1.98, m[a]	1.98, m[a]
10b	2.01, m[a]	2.01, m[a]	2.01, m	2.05, m	2.32, m[a]	2.32, m[a]
11	2.58, m[a]	2.58, m	2.48, m	2.48, m	2.56, m[a]	2.56, m[a]
16	2.03, s	2.03, s	2.05, s	2.05, s	2.05, s	2.05, s
17	1.58, s	1.58, s	1.64, s	1.64, s	1.65, s	1.65, s
18	1.45, m[a]	1.85, m[a]	1.82, m	1.70, m	1.41, m[a]	1.41, m[a]
18	1.85, m[a]	1.45, m[a]	1.70, m	1.82, m	1.71, m[a]	1.71, m[a]
19	1.37, s	1.37, s	1.37, s	1.37, s	1.37, s	1.37, s
20	1.37, s	1.37, s	1.37, s	1.37, s	1.36, s	1.36, s
2′	5.82, s	5.82, s	5.80, s	5.80, s	5.80, s	5.80, s
4′a	2.38, dd	2.38, dd	2.37, dd	2.37, dd	2.37, dd	2.37, dd
	(11.5, 17.5)	(11.0, 17.5)	(11.5, 17.5)	(11.5, 18.0)	(11.0, 17.5)	(11.5, 18.5)
4′b	2.21, dd	2.22, dd	2.21, dd	2.24, dd	2.23, dd	2.21, dd
	(4.0, 17.5)	(4.0, 18.0)	(4.0, 18.0)	(4.0, 18.0)	(4.0, 18.0)	(4.0, 18.0)
5′	5.11, m	5.12, m	5.09, m	5.10, m	5.09, m	5.11, m
6′	5.32, d (7.5)	5.33, d (8.5)	5.27, d (8.5)	5.31, d (8.5)	5.30, d (8.5)	5.28, d (8.5)
8′	2.04, m[a]	2.04, m[a]	2.08, dd	2.08, dd	2.05, m[a]	2.08, dd
			(6.0, 14.0)	(6.0, 14.0)		(6.0, 14.0)
9′	1.95, m[a]	1.95, m[a]	1.74, 1.81, m	1.74,1.81, m	1.96, m[a]	1.74, 1.81, m[a]
10′a	1.61, m[a]	1.61, m[a]	1.89, m[a]	1.89, m[a]	1.80, m[a]	1.89, m[a]
10′b	2.01, m[a]	2.01, m[a]	1.77, m[a]	1.77, m[a]	2.15, m[a]	1.77, m[a]
13′	5.22, s	5.21, s	5.26, s	5.24, s	5.20, s	5.22, s
16′	1.99, s	1.99, s	2.05, s	2.05, s	1.98, s	2.04, s
17′	1.70, s	1.58, s	1.64, s	1.64, s	1.61, s	1.61, s
18′α	2.12, m[a]	1.82, m[a]	1.72, m[a]	1.82, m[a]	2.18, m[a]	1.72, m[a]
18′β	1.82, m[a]	2.12, m[a]	1.82, m[a]	1.72, m[a]	1.67, m[a]	1.82, m[a]
19′	1.36, s	1.36, s	1.35, s	1.36, s	1.36, s	1.36, s
20′	1.34, s	1.34, s	1.32, s	1.32, s	1.34, s	1.35, s

[a]Overlapping signals

Table S2 ^{13}C NMR Spectroscopic Data (125 MHz, $CDCl_3$) for Compounds **9**, **11** and **13**

no.	**9**	**11**	**13**	no.	**9**	**11**	**13**
1	172.9	172.9	172.9	1′	165.0	165.1	165.1
2	117.2	116.6	117.3	2′	116.3	116.5	116.4
3	168.0	165.0	167.9	3′	156.7	156.7	156.8
4	84.2	87.5	84.1	4'	35.0	35.0	35.0
5	30.2	30.3	30.2	5′	74.0	74.1	74.1
6	116.6	116.6	116.6	6′	122.3	122.4	122.3
7	139.4	141.7	139.2	7′	141.9	141.9	142.1
8	39.1	39.3	39.1	8′	39.6	39.3	39.4
9	25.0	25.0	25.0	9′	29.6	30.3	29.6
10	31.2	31.3	31.2	10′	29.0	28.5	34.0
11	35.4	35.0	35.4	11′	42.0	41.7	42.0
12	189.6	189.7	189.9	12′	194.9	194.3	194.8
13	111.3	110.9	111.1	13′	101.8	101.5	101.8
14	204.1	203.9	204.1	14′	207.1	207	207.1
15	87.7	87.5	87.7	15′	88.9	88.7	88.9
16	23.1	23.1	23.1	16′	13.8	16.6	13.8
17	16.6	16.7	16.4	17′	16.1	16.7	16.1
18	23.8	22.3	22.3	18′	23.8	23.7	23.8
19	22.8	22.8	22.8	19′	22.7	22.8	22.7
20	22.7	22.7	22.7	20′	22.7	22.8	22.7

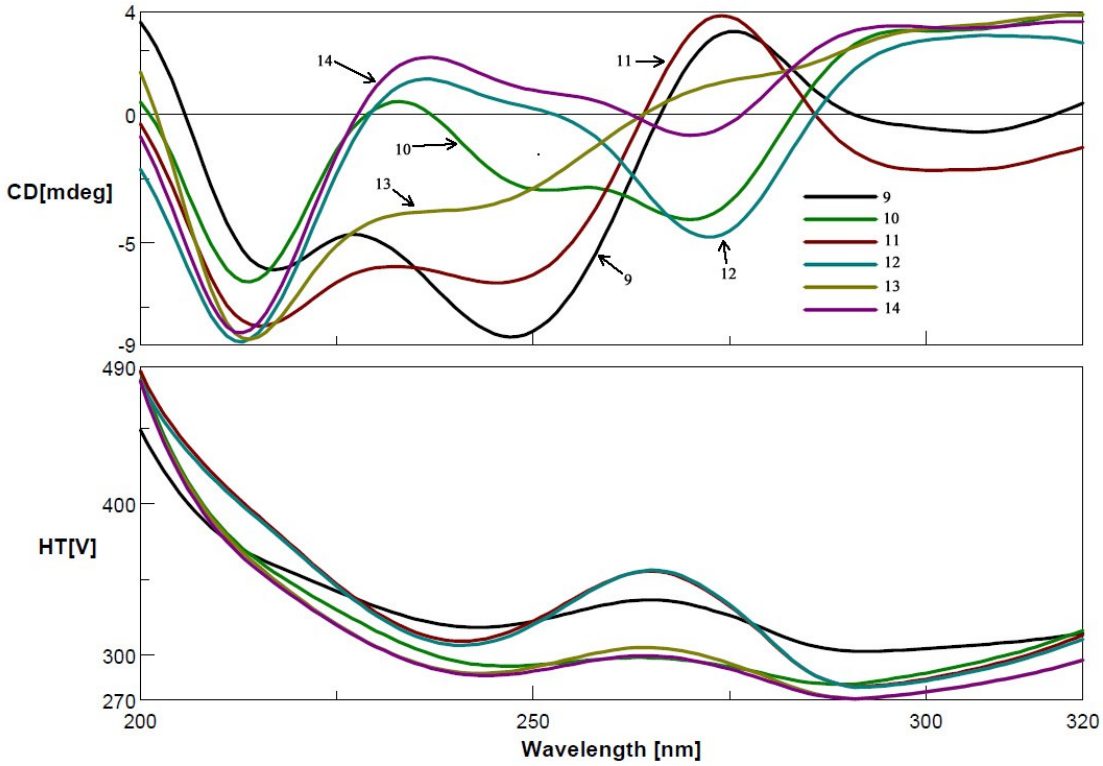
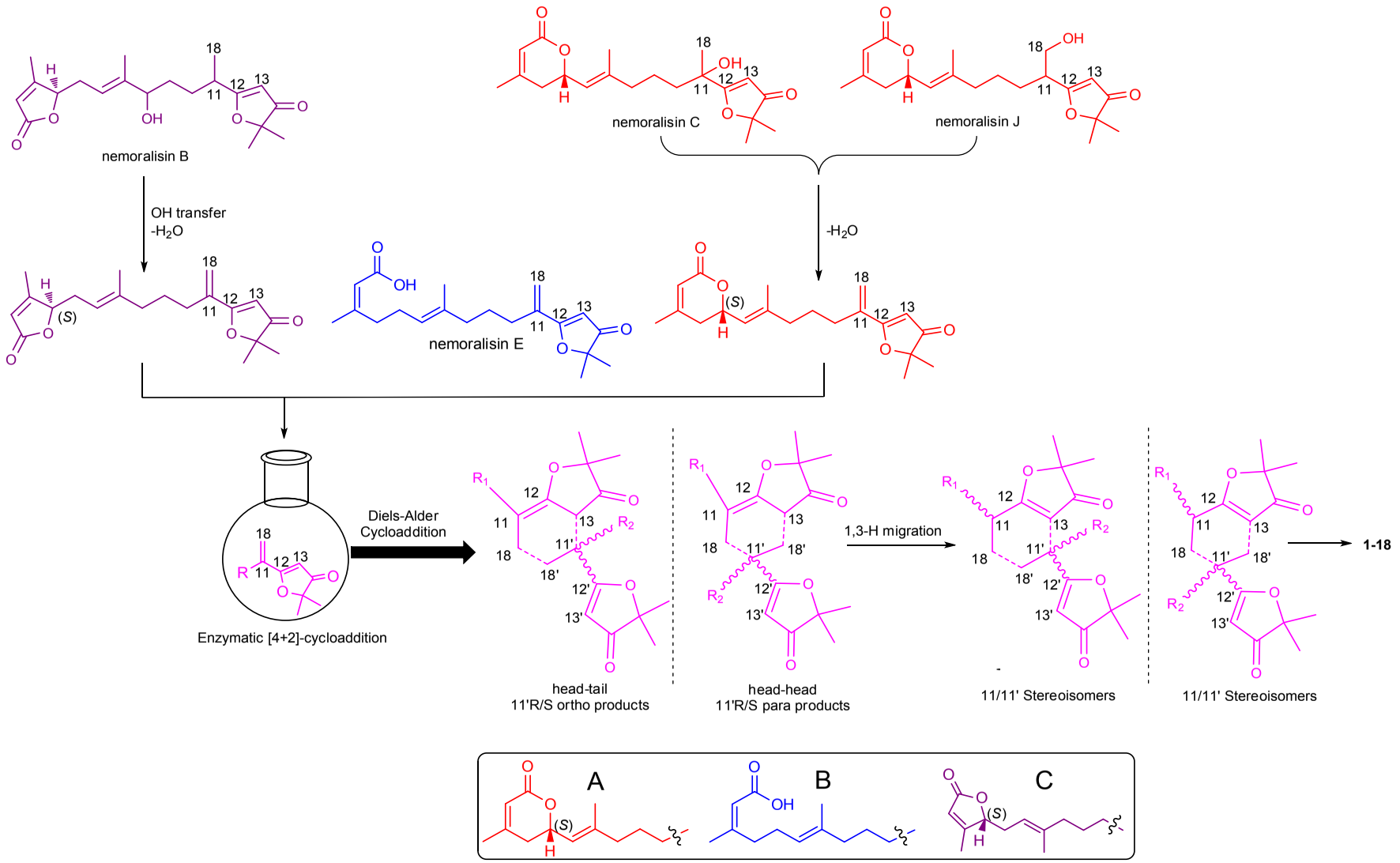
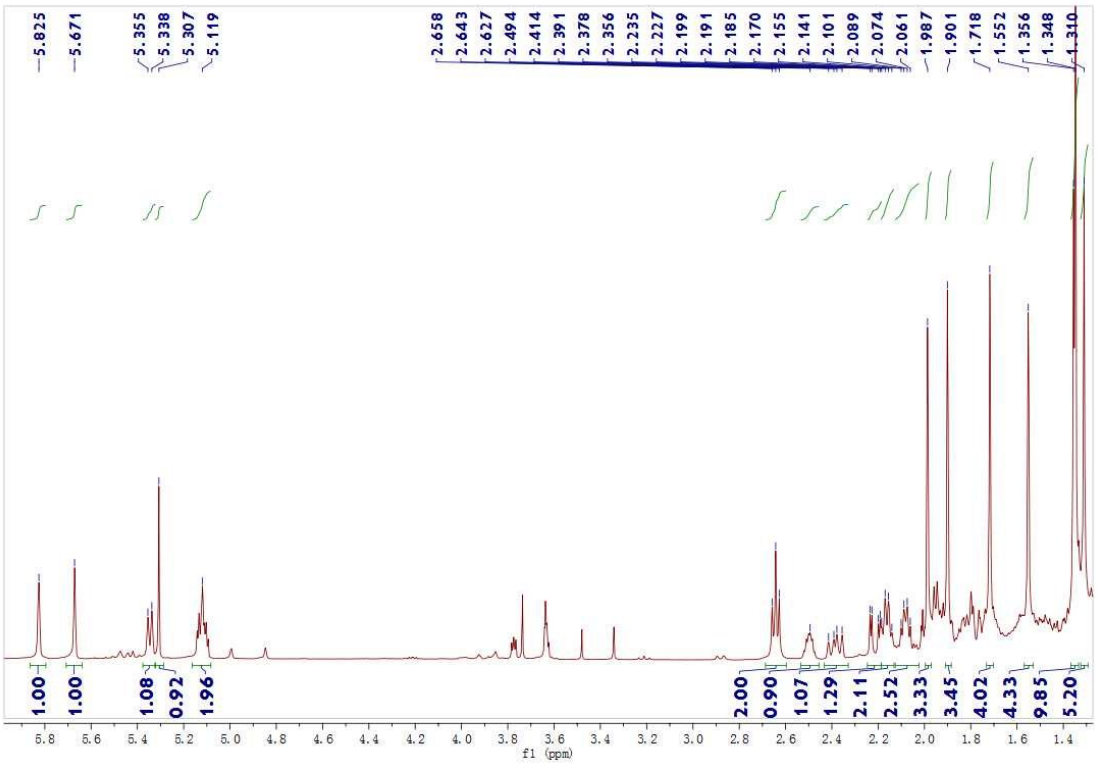


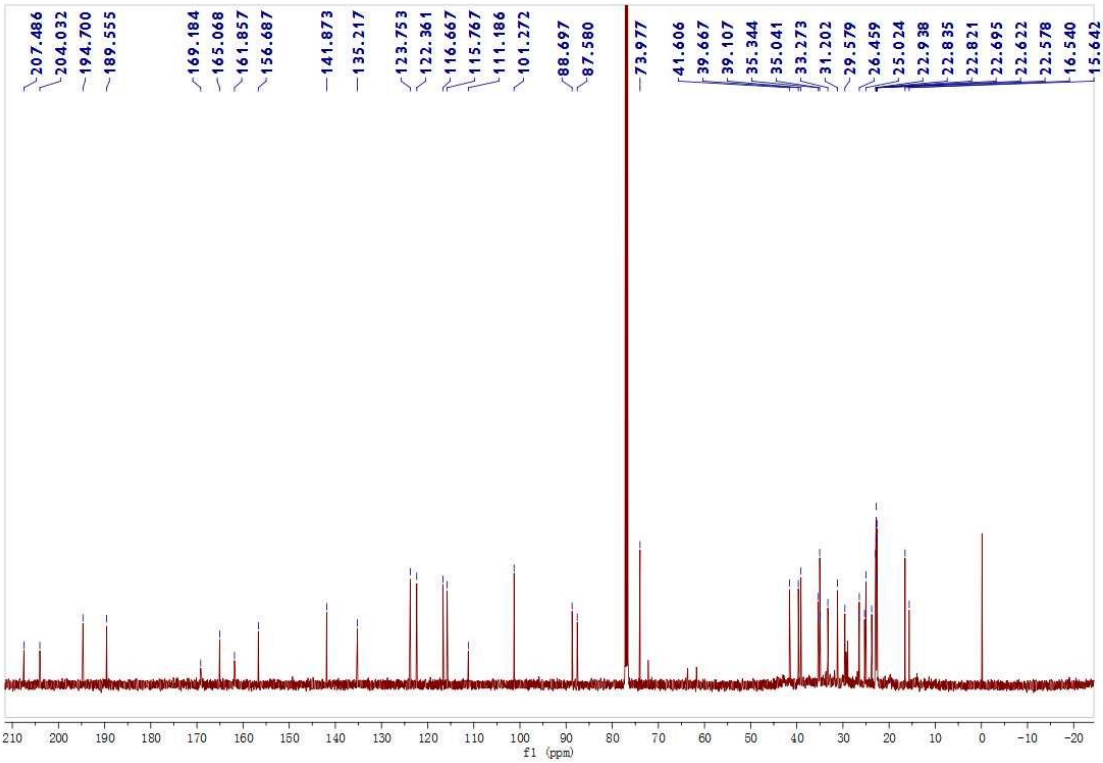
Fig. S1 CD spectra of **9-14**.



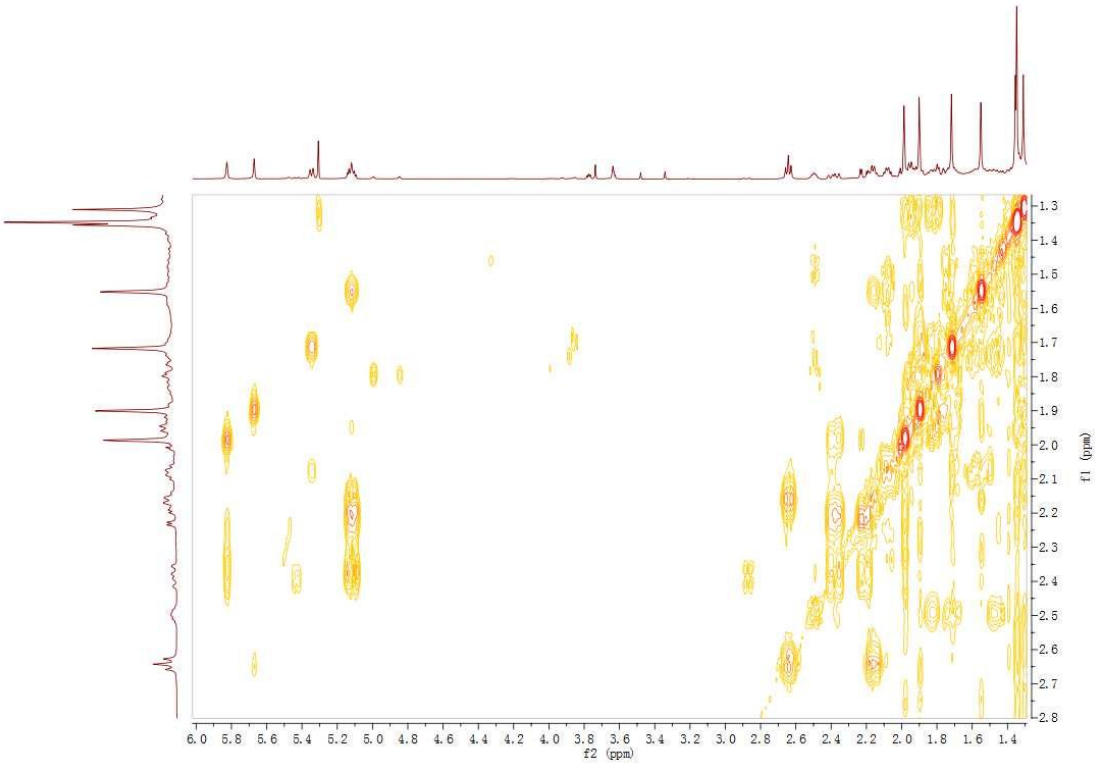
Scheme S1 Plausible biosynthetic pathway for **1-18.**



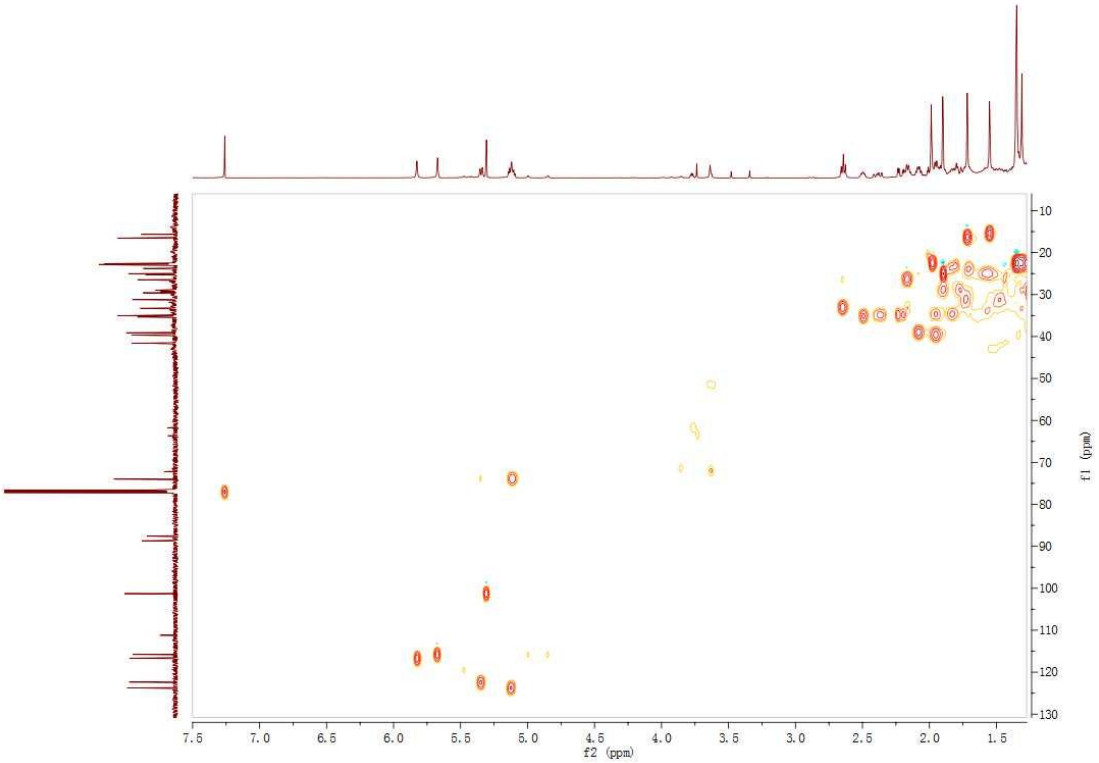
S1. ^{1}H NMR spectrum of compound aphanamene C (**1**) in $CDCl_3$



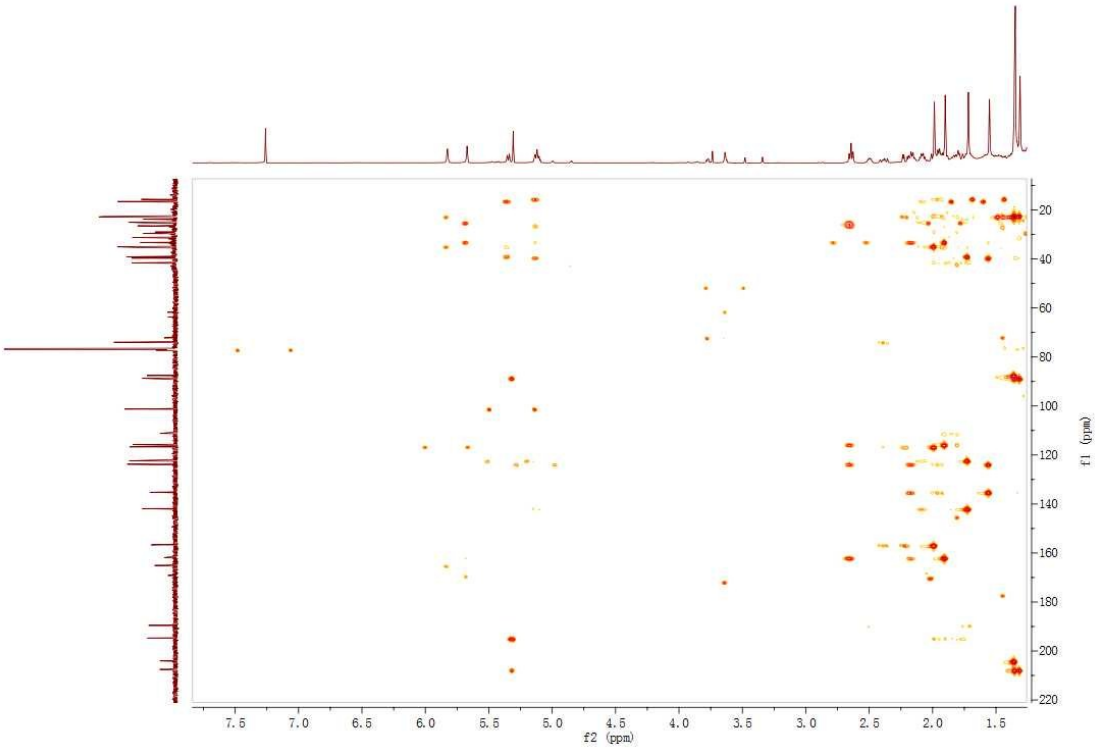
S2. ^{13}C NMR spectrum of compound aphanamene C (**1**) in $CDCl_3$



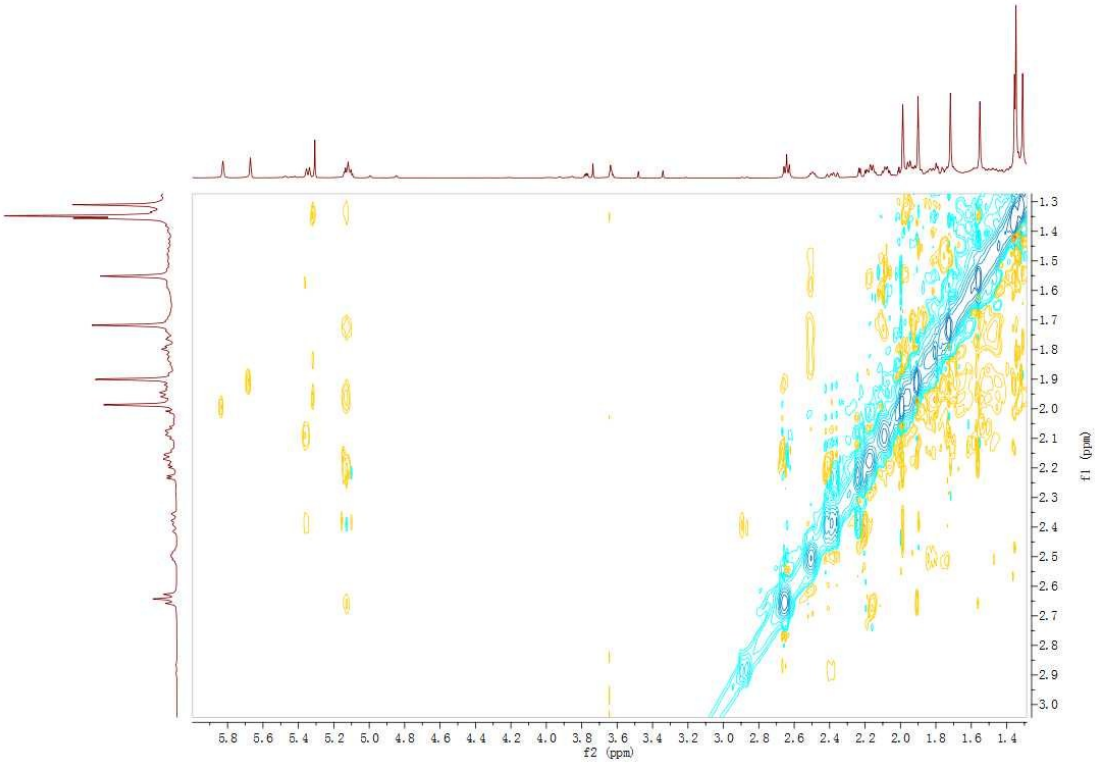
S3. ^{1}H-^{1}HCOSY spectrum of compound aphanamene C (**1**) in $CDCl_3$



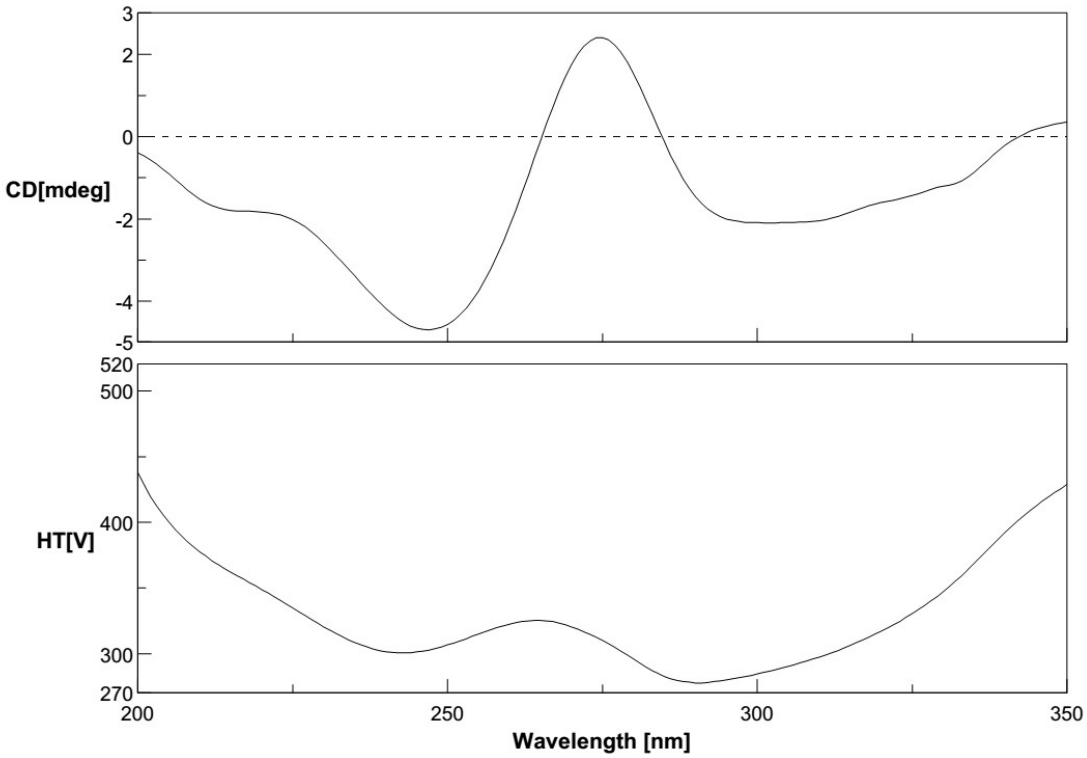
S4. HSQC spectrum of compound aphanamene C (**1**) in $CDCl_3$



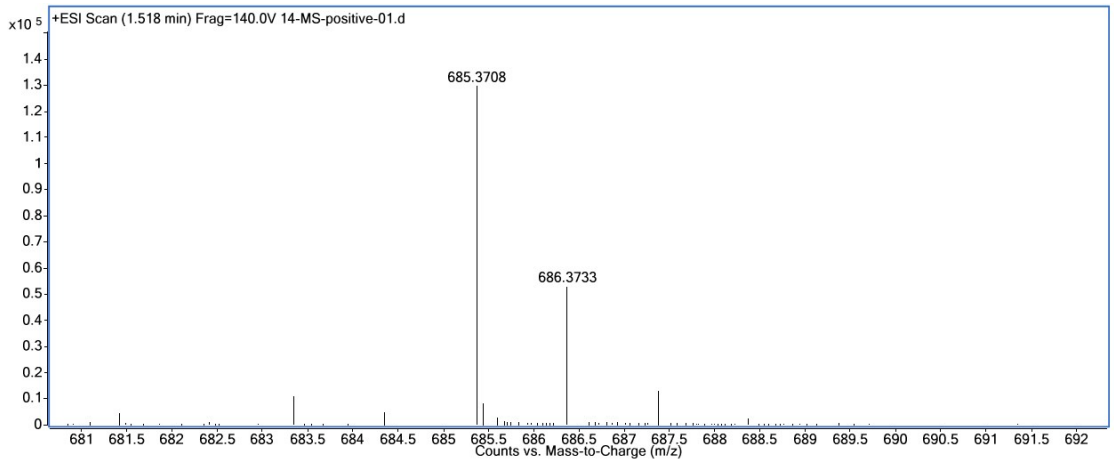
S5. HMBC spectrum of compound aphanamene C (**1**) in $CDCl_3$



S6. ROESY spectrum of compound aphanamene C (**1**) in $CDCl_3$



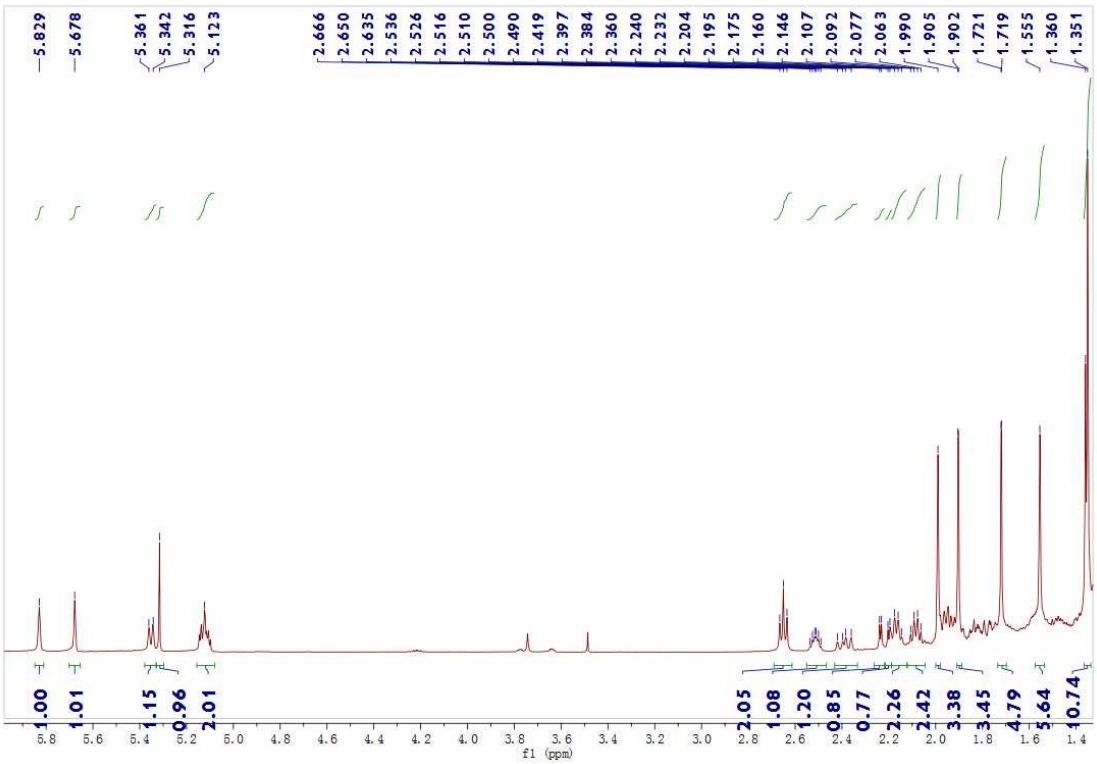
S7. CD spectrum of compound aphanamene C (**1**) in CH_3OH



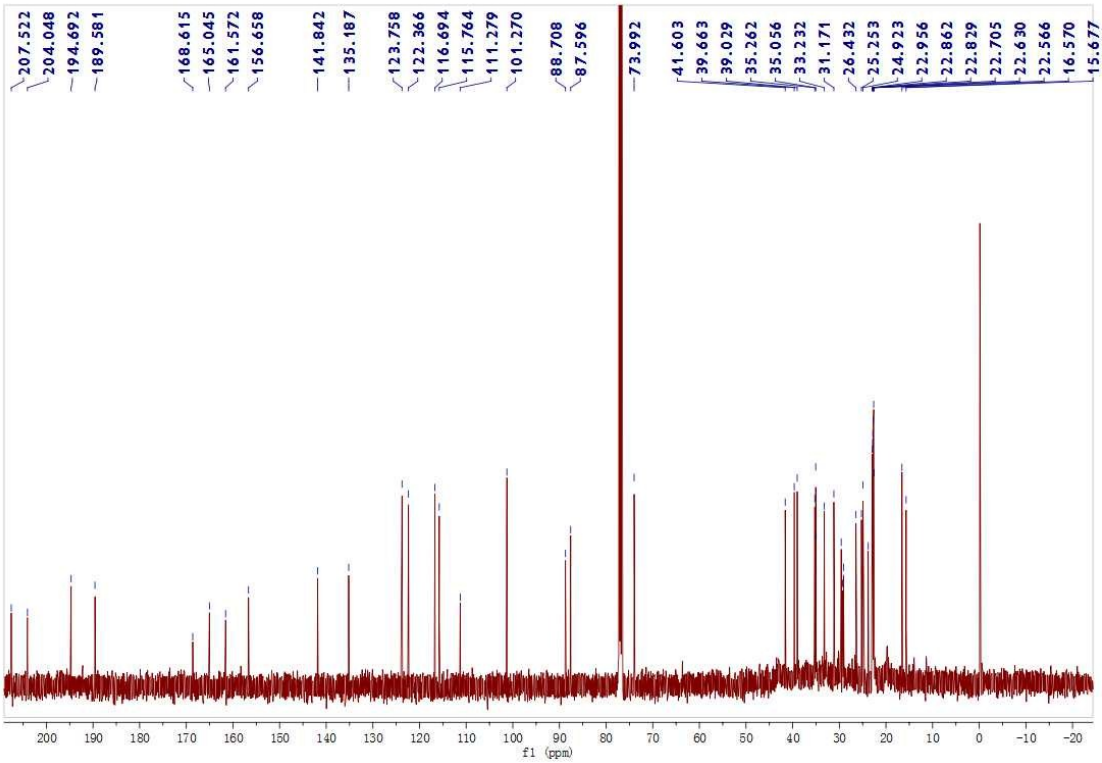
Elemental Composition Calculator

Target m/z:	685.3708	**Result type:**	Positive ions	**Species:**	$[M+Na]^+$
Elements:		C (0-80); H (0-120); O (0-30); N (0-10); Na (0-5); Cl (0-5)			
Ion Formula		**Calculated m/z**		**PPM Error**	
C40H52NaO8		685.3711		0.47	

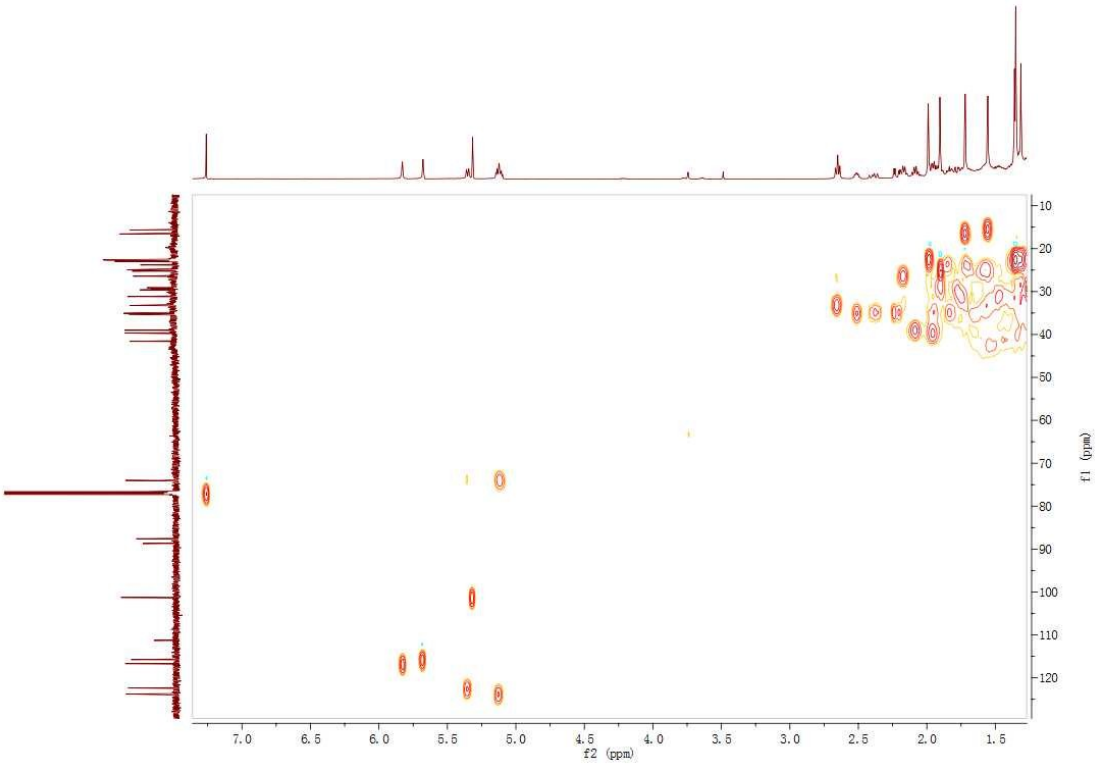
S8. HRESIMS spectrum of compound aphanamene C (**1**) in CH_3OH



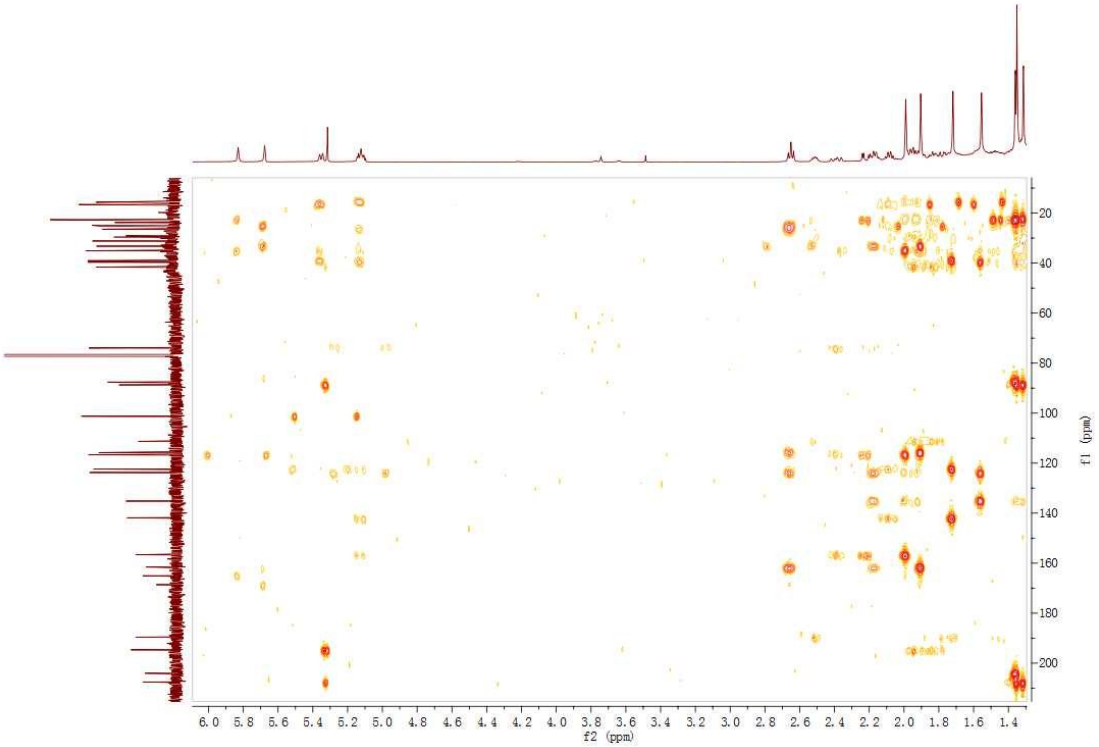
S9. ^{1}H NMR spectrum of compound aphanamene D (**2**) in $CDCl_3$



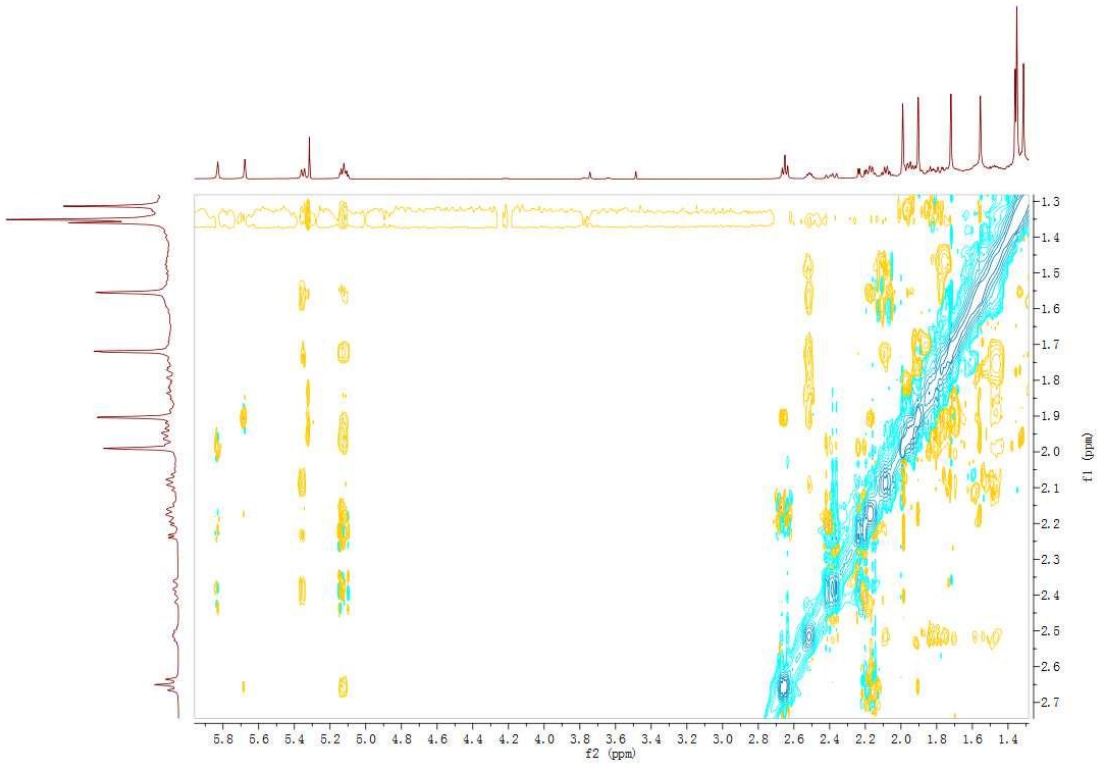
S10. ^{13}C NMR spectrum of compound aphanamene D (**2**) in $CDCl_3$



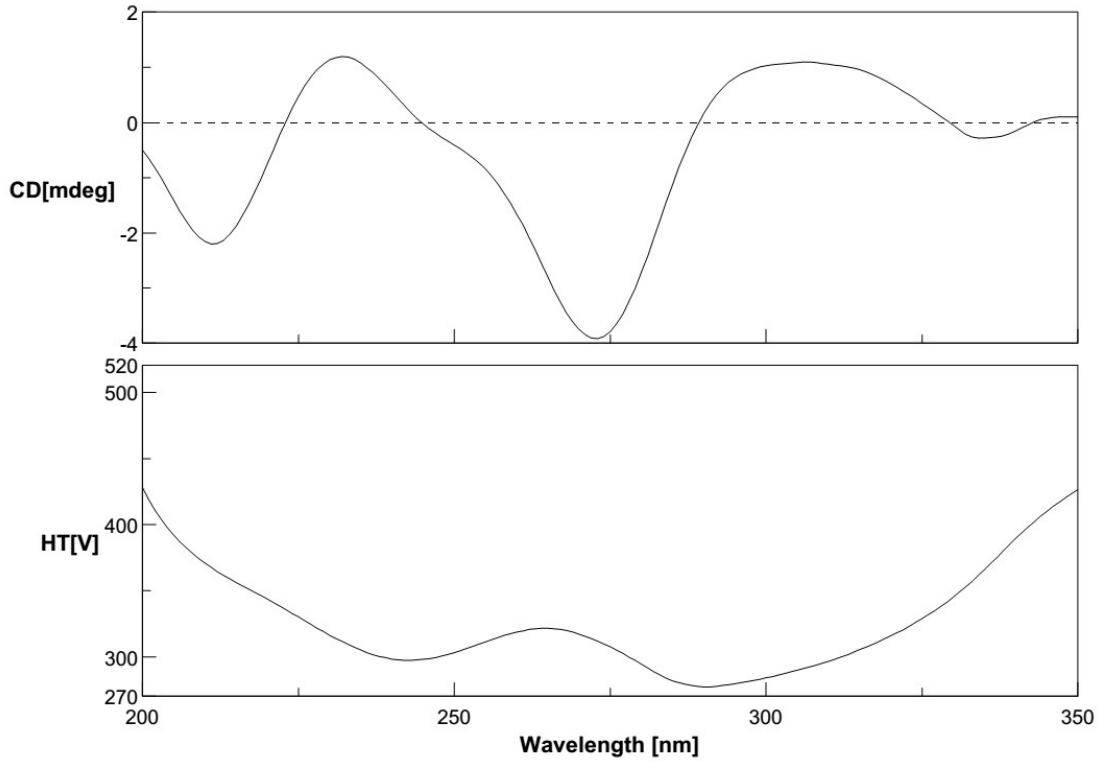
S11. HSQC spectrum of compound aphanamene D (**2**) in $CDCl_3$



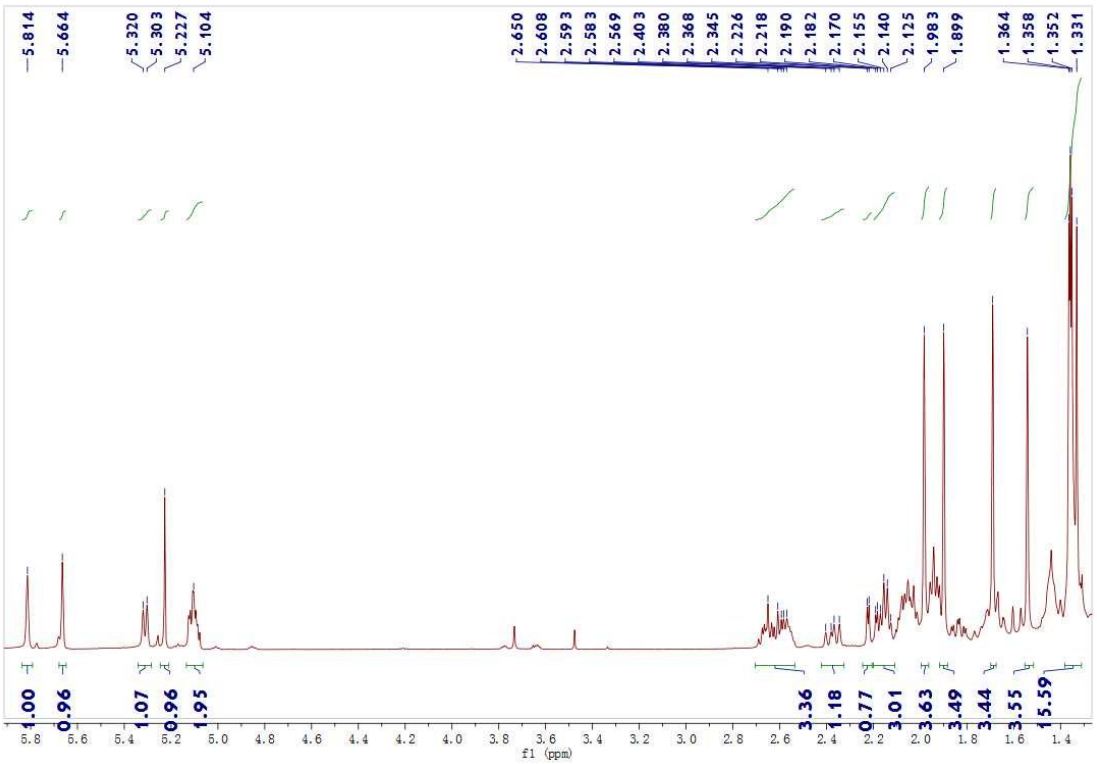
S12. HMBC spectrum of compound aphanamene D (**2**) in $CDCl_3$



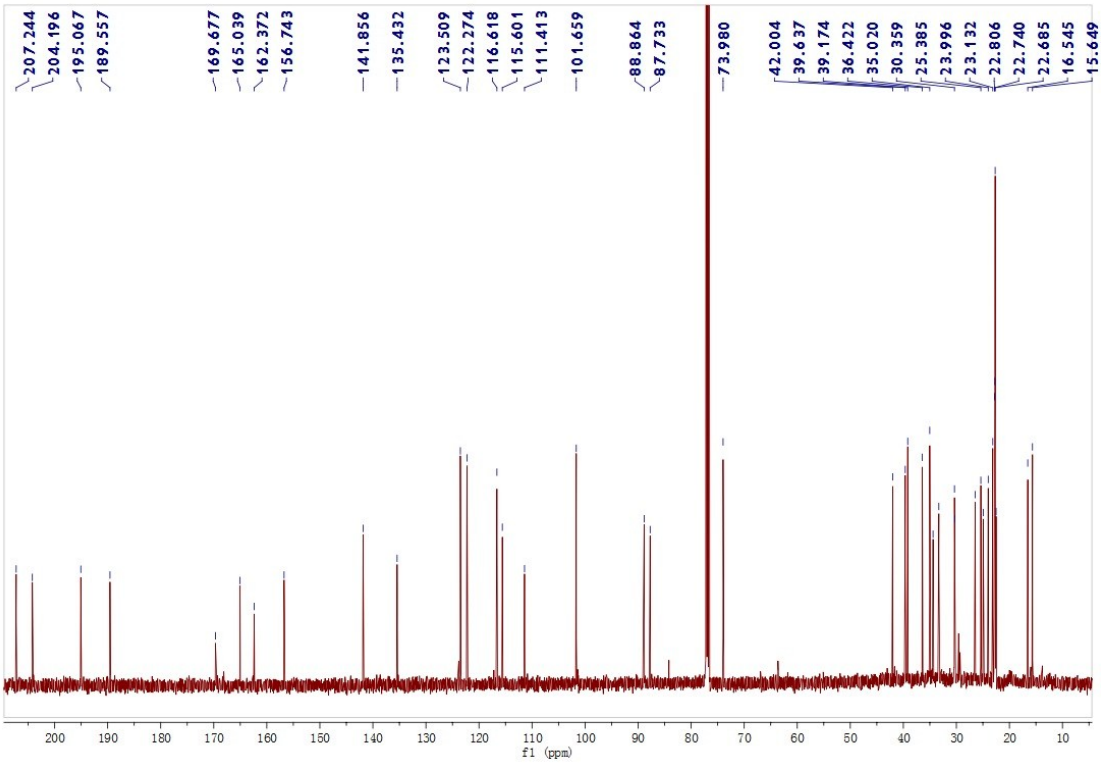
S13. ROESY spectrum of compound aphanamene D (**2**) in $CDCl_3$



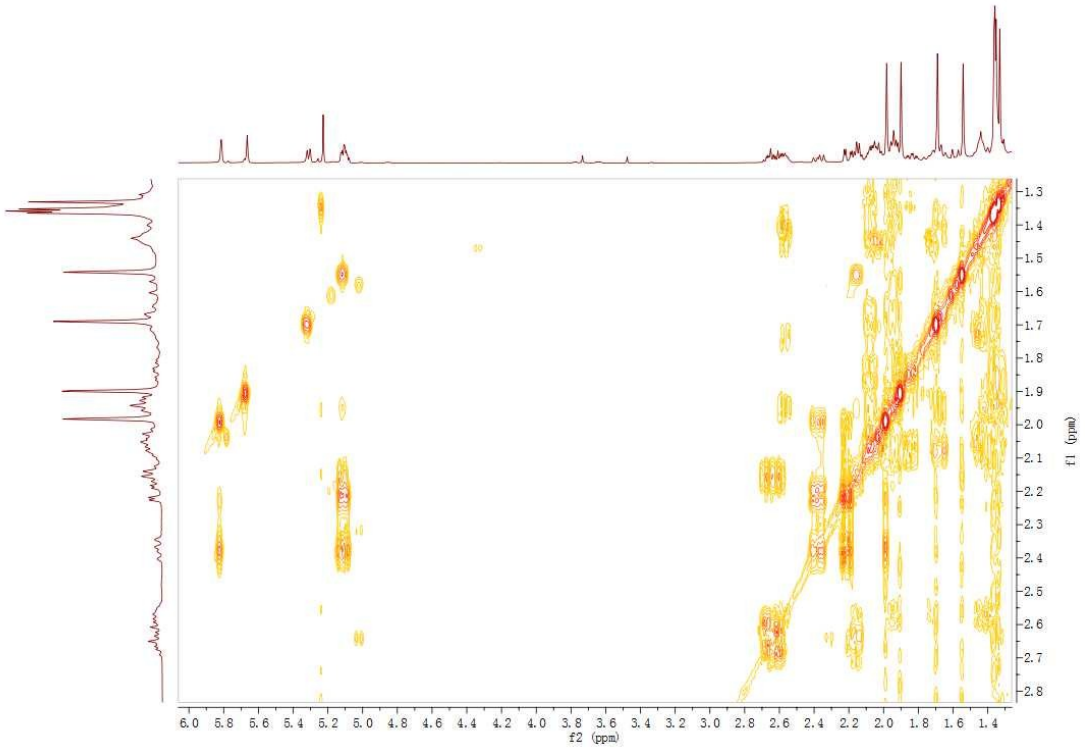
S14. CD spectrum of compound aphanamene D (**2**) in CH_3OH



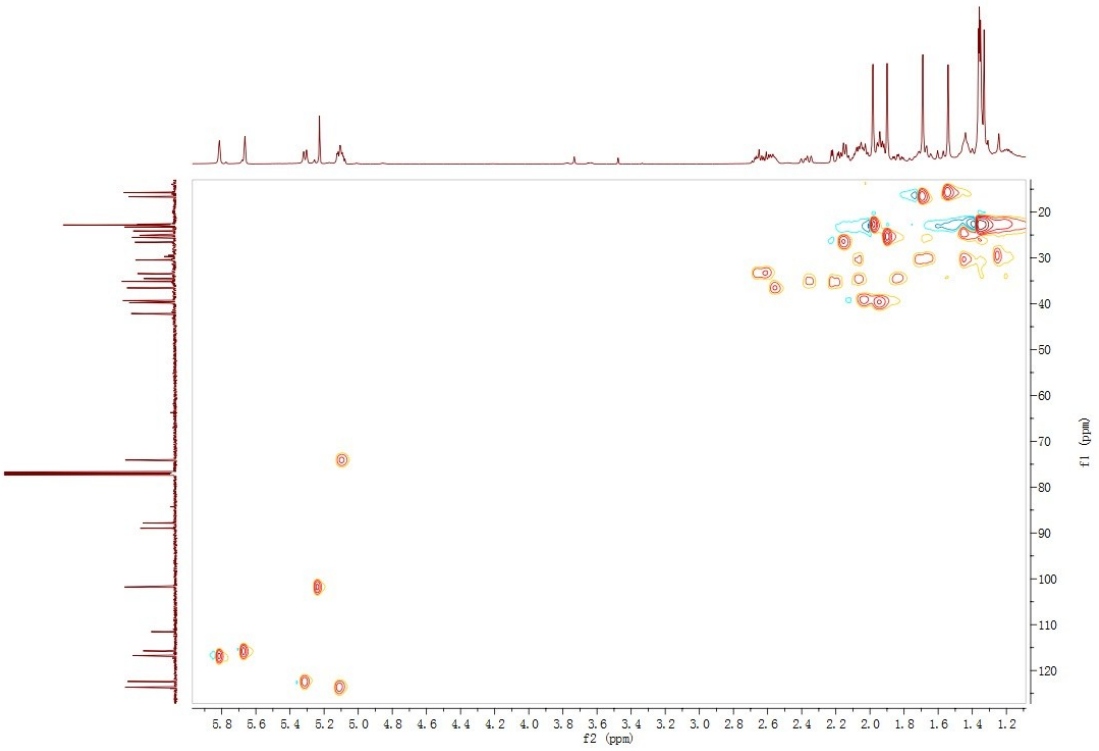
S15. ^{1}H NMR spectrum of compound aphanamene E (**3**) in $CDCl_3$



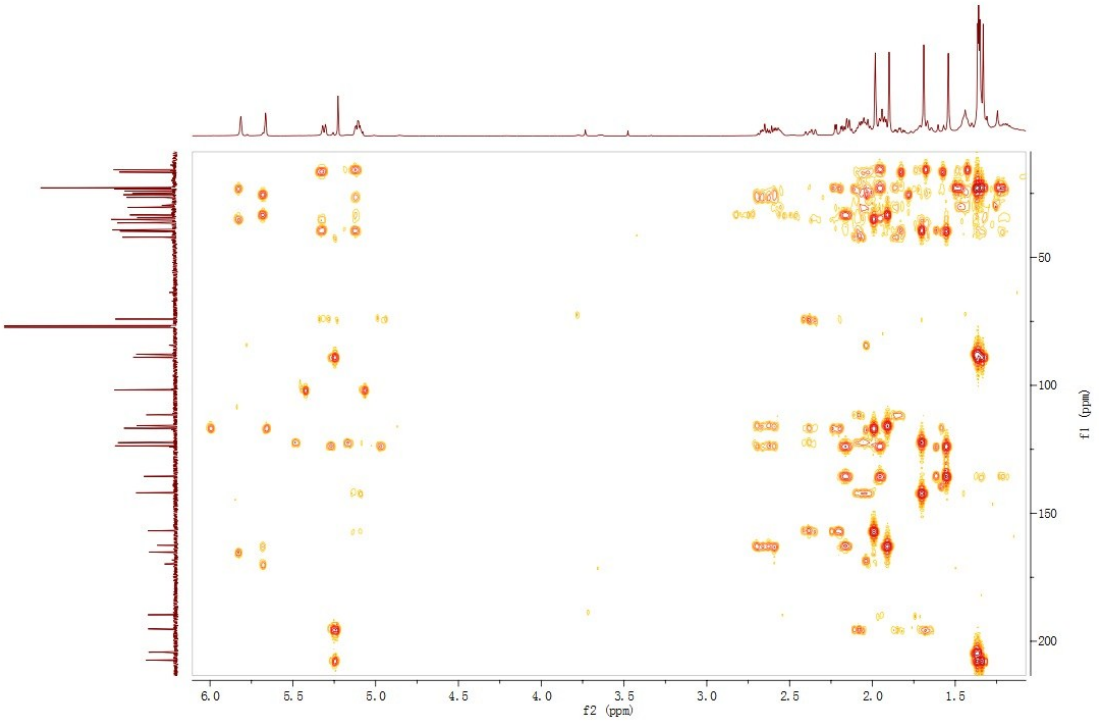
S16. ^{13}C NMR spectrum of compound aphanamene E (**3**) in $CDCl_3$



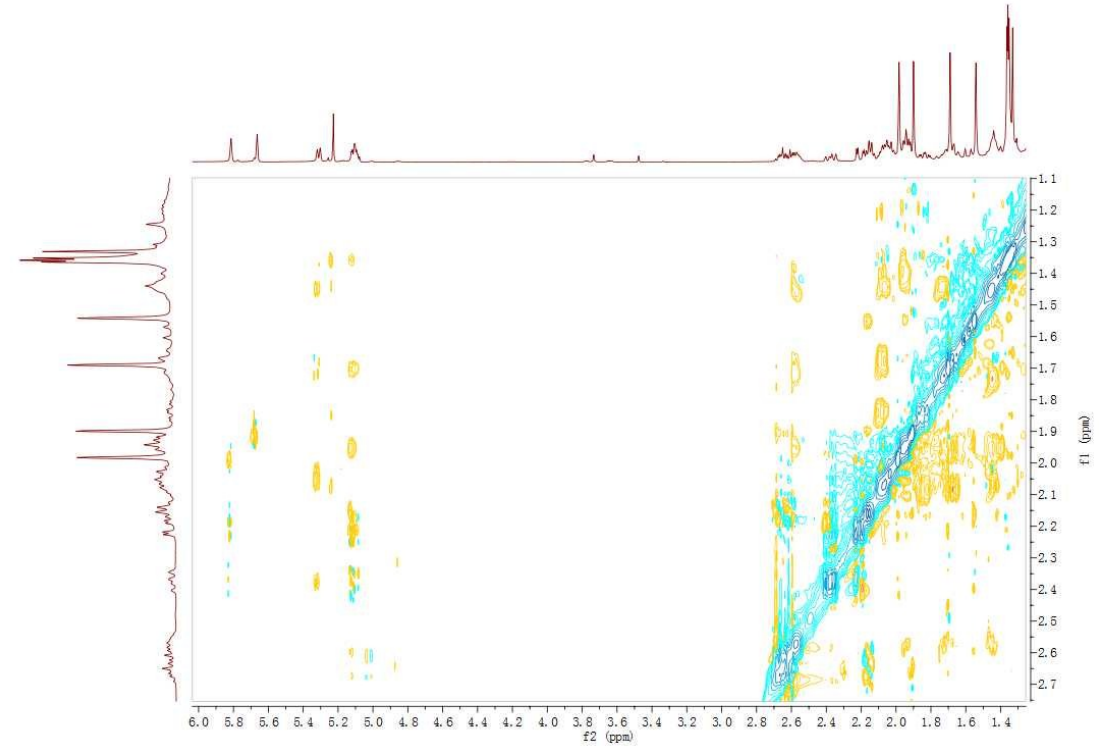
S17. ^{1}H-^{1}H NMR spectrum of compound aphanamene E (**3**) in $CDCl_3$



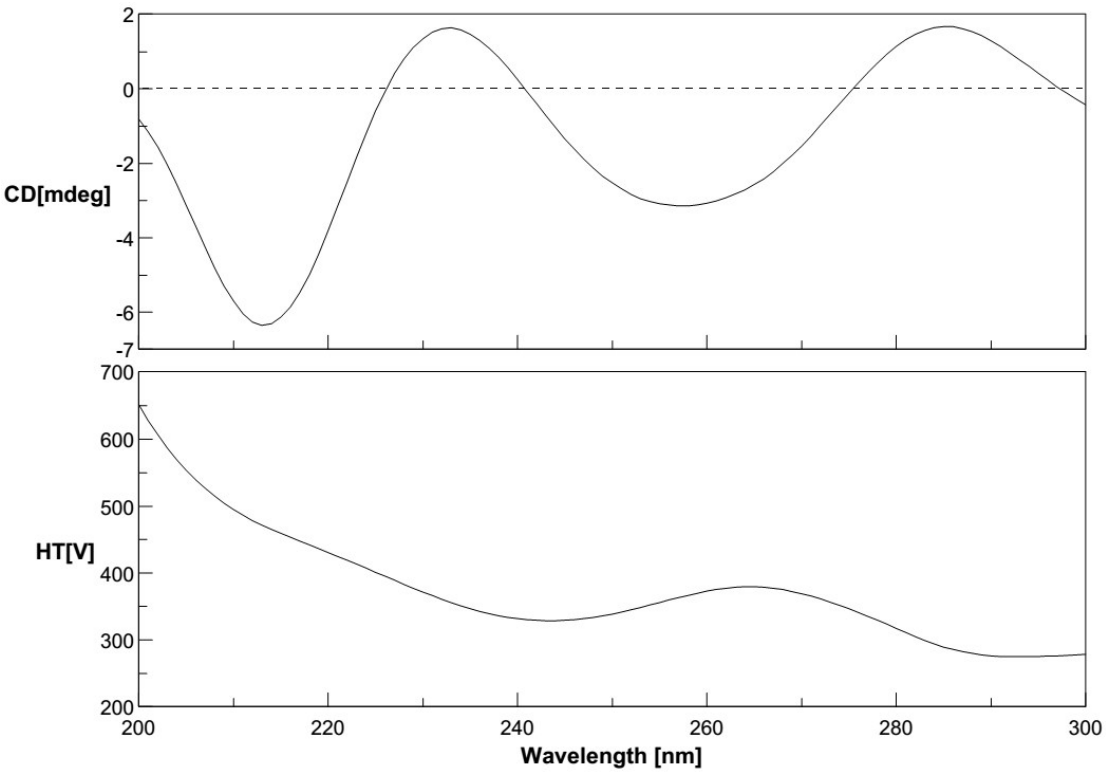
S18. HSQC spectrum of compound aphanamene E (**3**) in $CDCl_3$



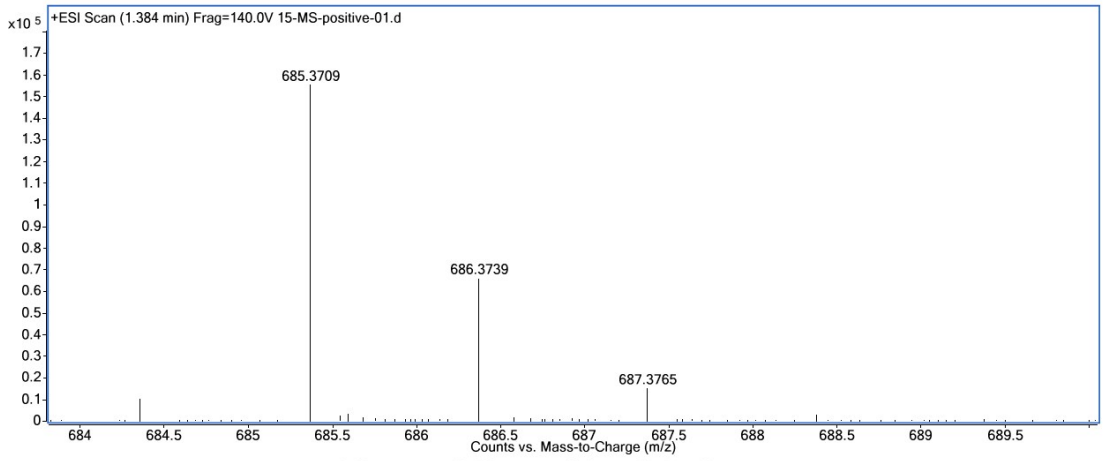
S19. HMBC spectrum of compound aphanamene E (**3**) in $CDCl_3$



S20. ROESY spectrum of compound aphanamene E (**3**) in $CDCl_3$



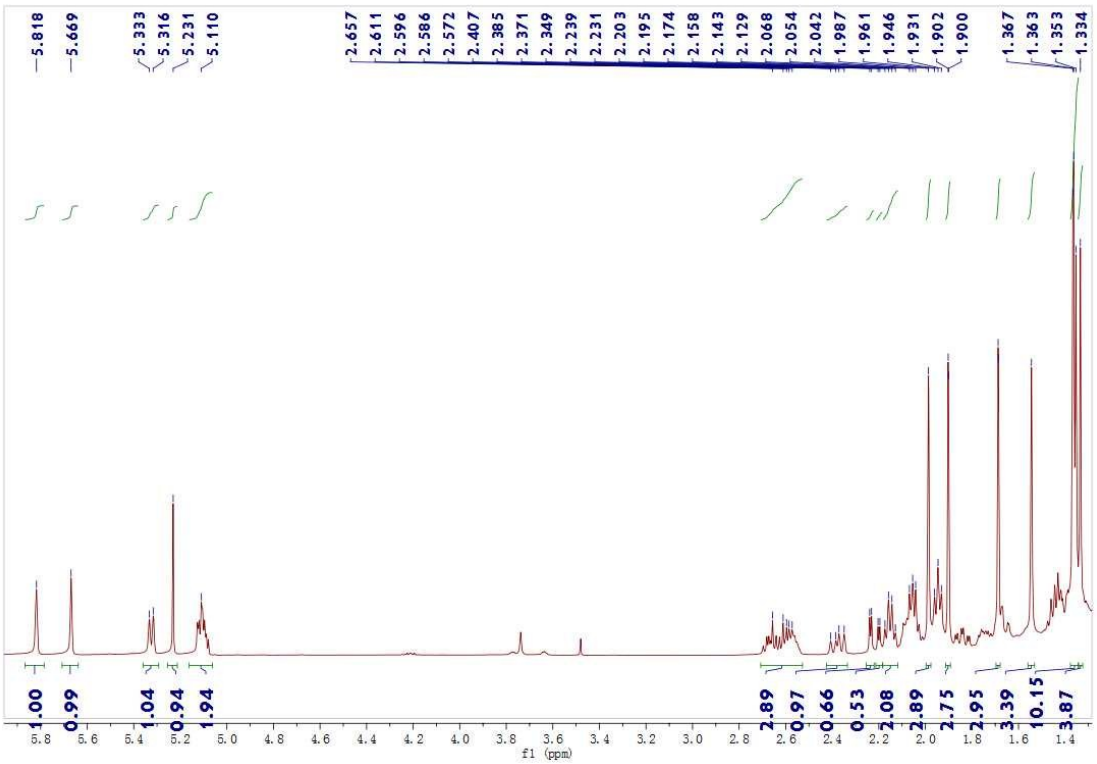
S21. CD spectrum of compound aphanamene E (**3**) in CH_3OH



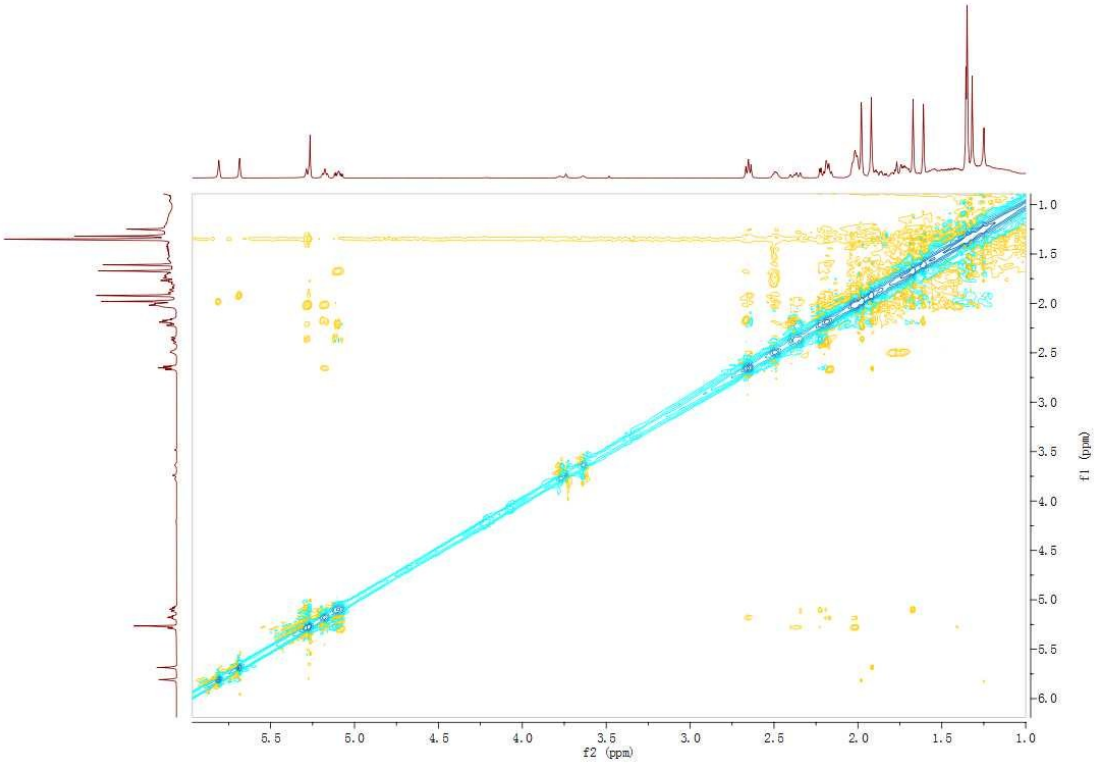
Elemental Composition Calculator

Target m/z:	685.3709	**Result type:**	Positive ions	**Species:**	$[M+Na]^+$
Elements:		C (0-80); H (0-120); O (0-30); N (0-10); Na (0-5); Cl (0-5)			
Ion Formula		**Calculated m/z**		**PPM Error**	
C40H52NaO8		685.3711		0.21	

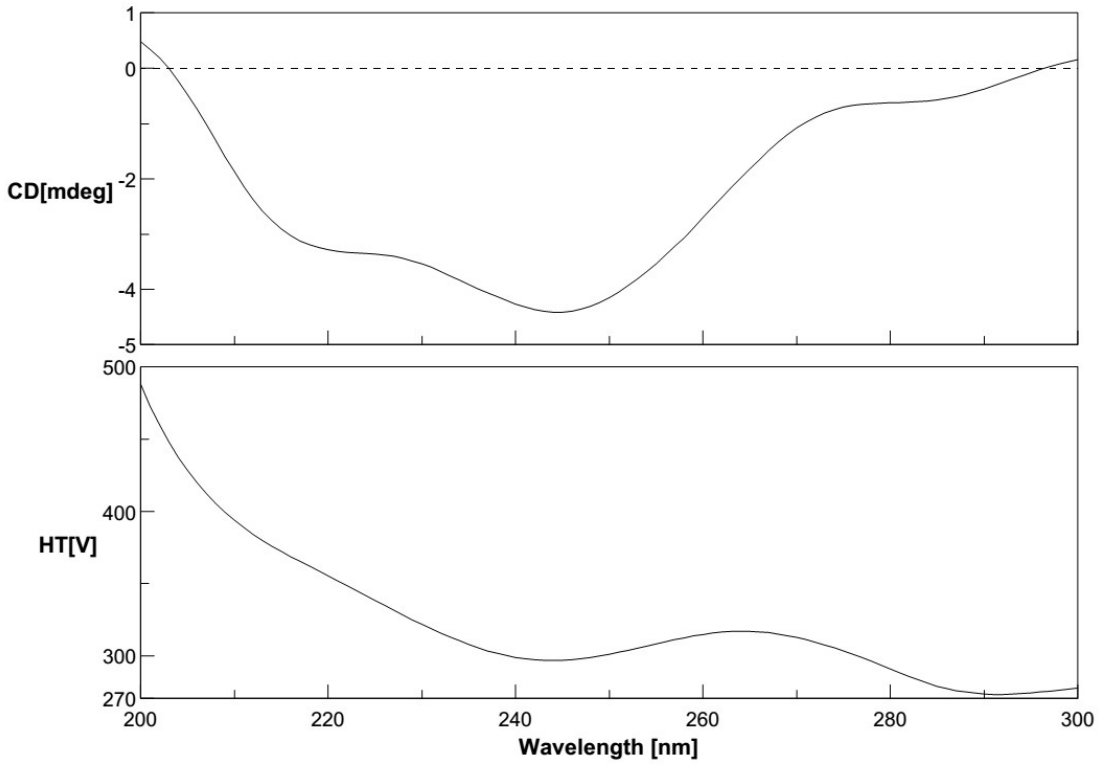
S22. HRESIMS spectrum of compound aphanamene E (**3**) in CH_3OH



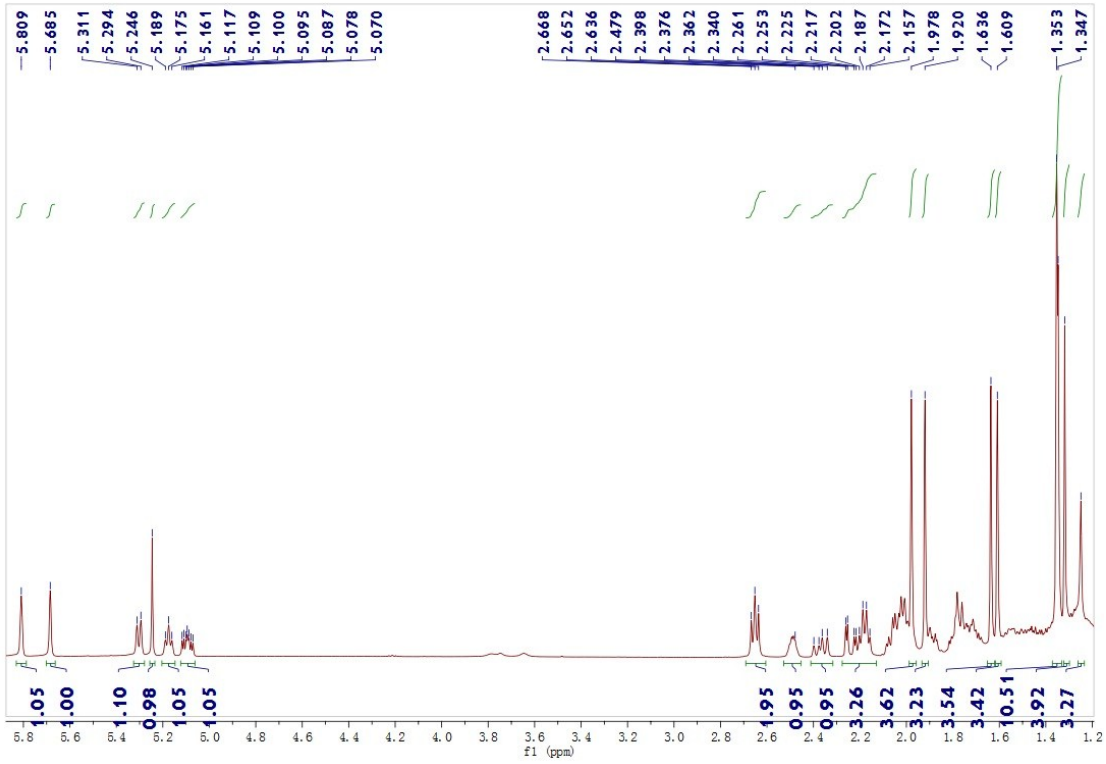
S23. ^{1}H NMR spectrum of compound aphanamene F (**4**) in $CDCl_3$



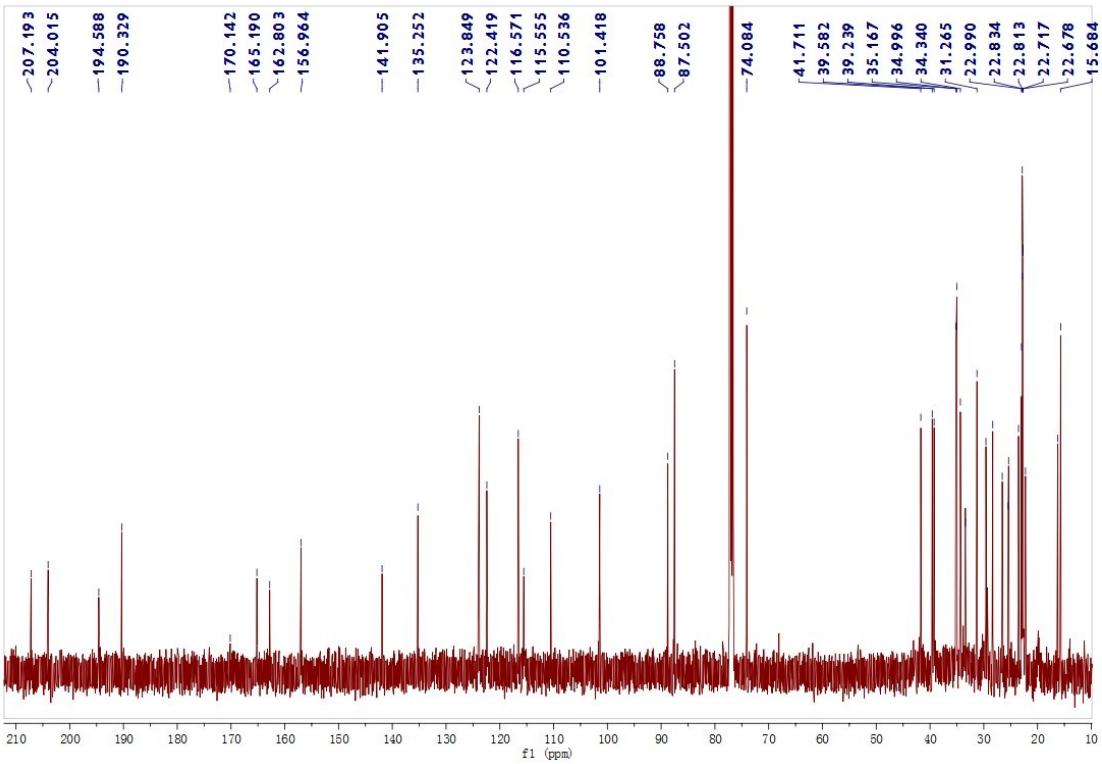
S24. ROESY spectrum of compound aphanamene F (**4**) in $CDCl_3$



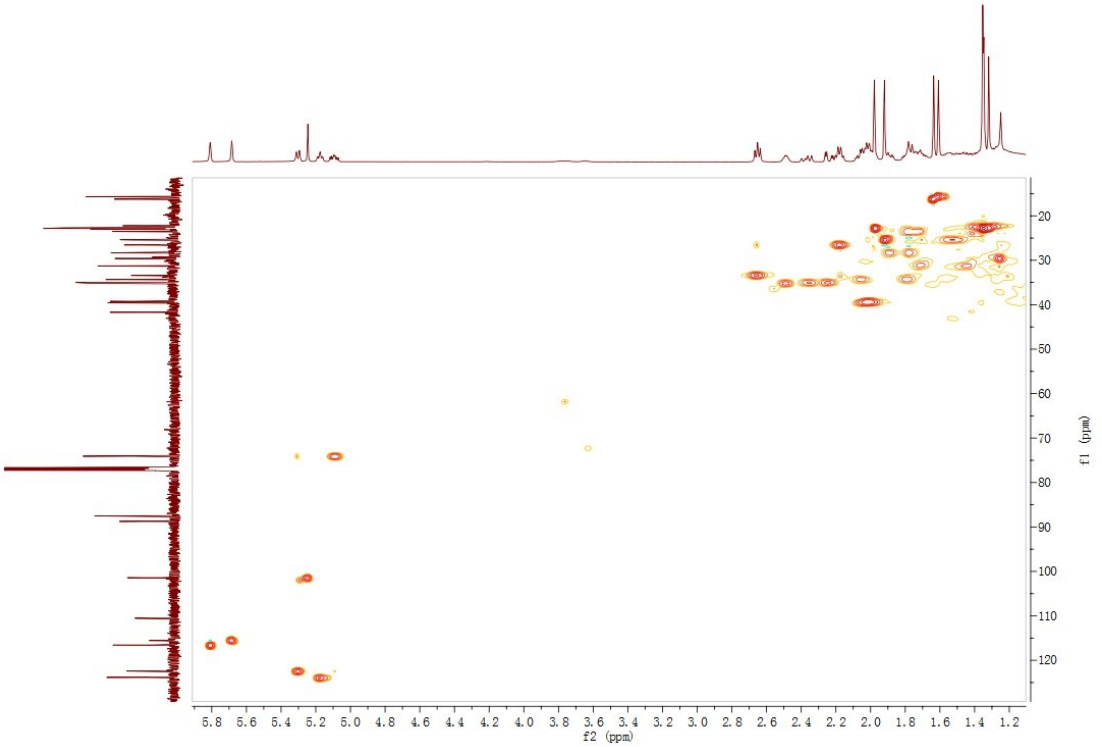
S25. CD spectrum of compound aphanamene F (**4**) in CH_3OH



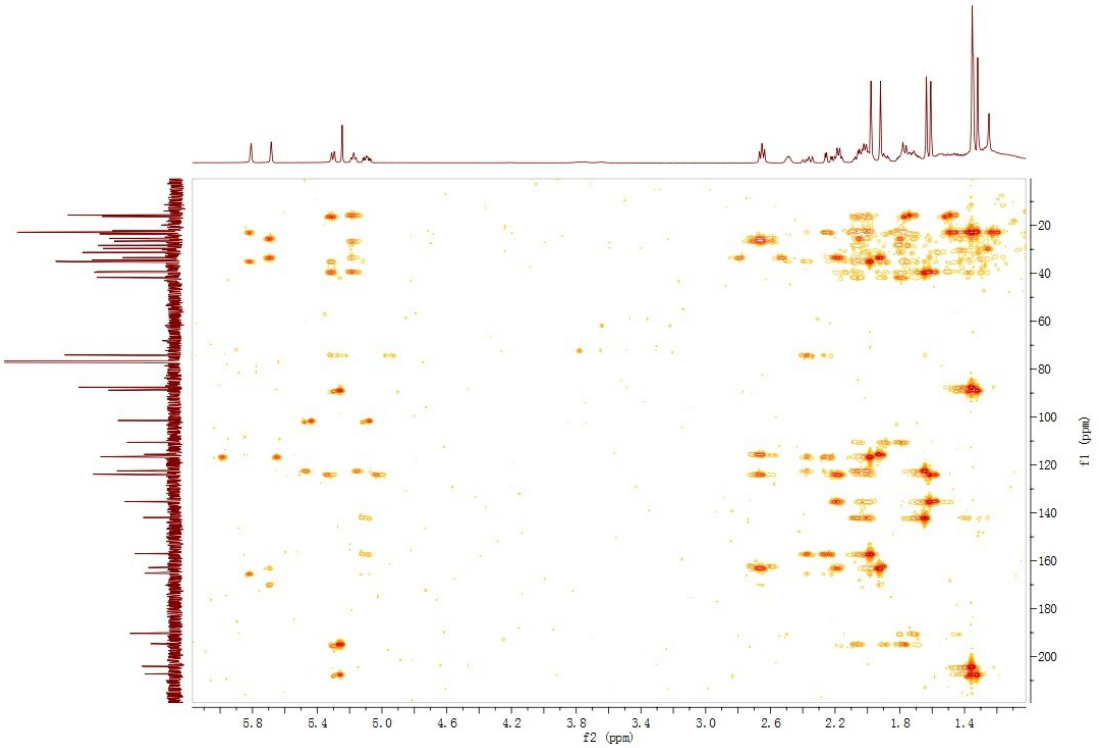
S26. ^1H NMR spectrum of compound aphanamene G (**5**) in $CDCl_3$



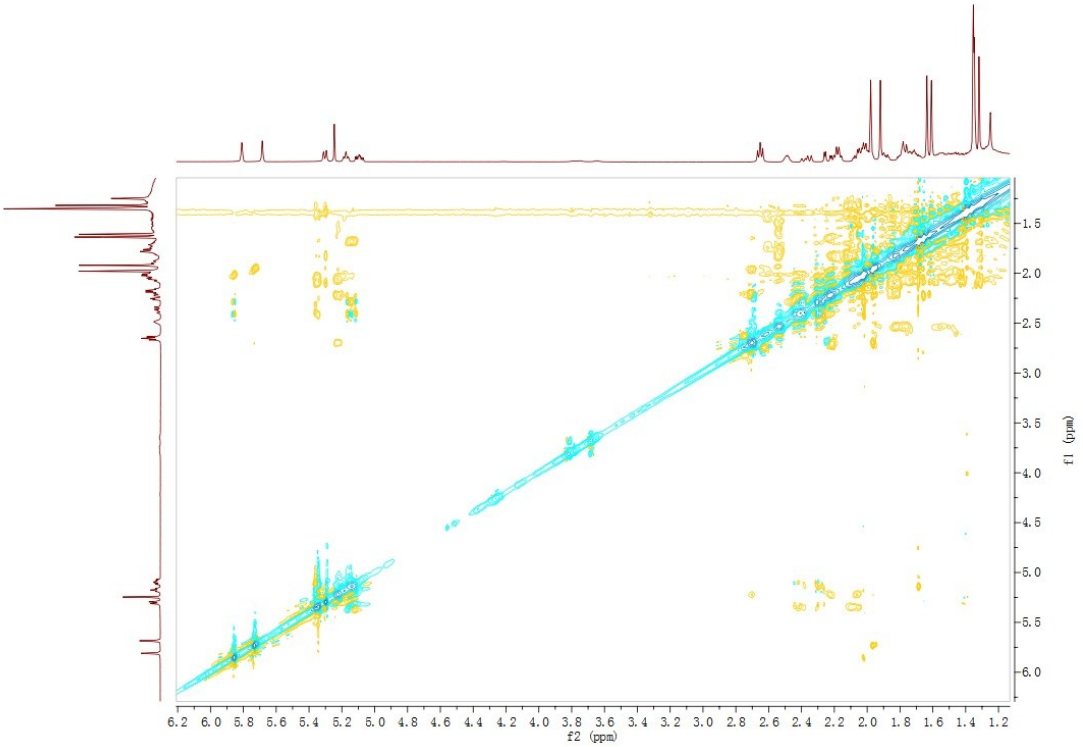
S27. ^{13}C NMR spectrum of compound aphanamene G (**5**) in $CDCl_3$



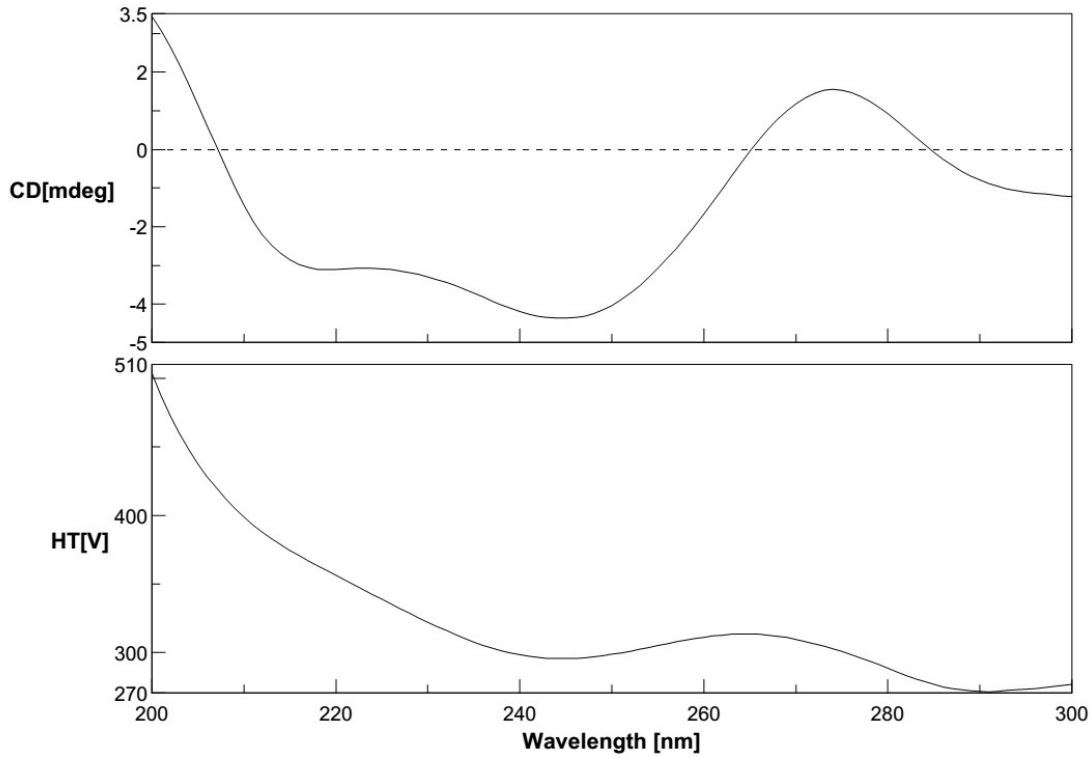
S28. HSQC spectrum of compound aphanamene G (**5**) in $CDCl_3$



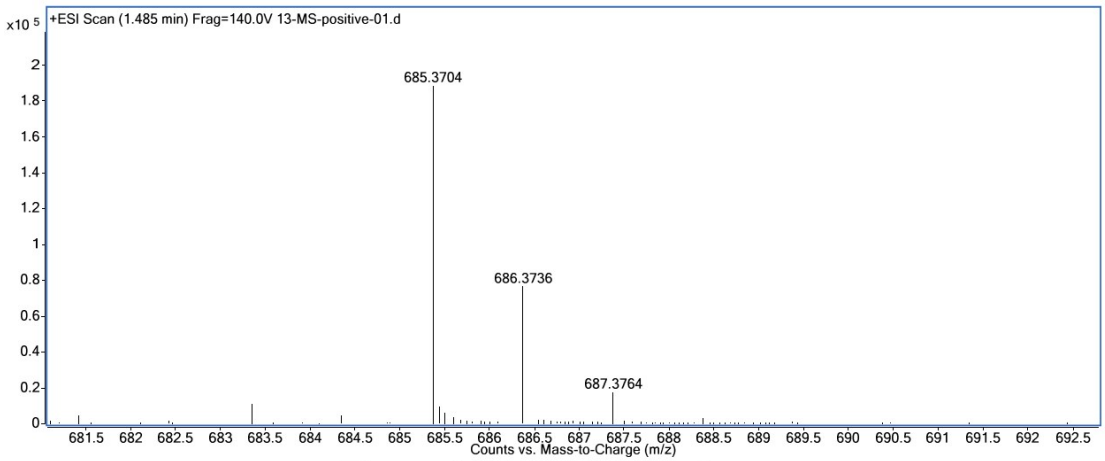
S29. HMBC spectrum of compound aphanamene G (**5**) in $CDCl_3$



S30. ROESY spectrum of compound aphanamene G (**5**) in $CDCl_3$



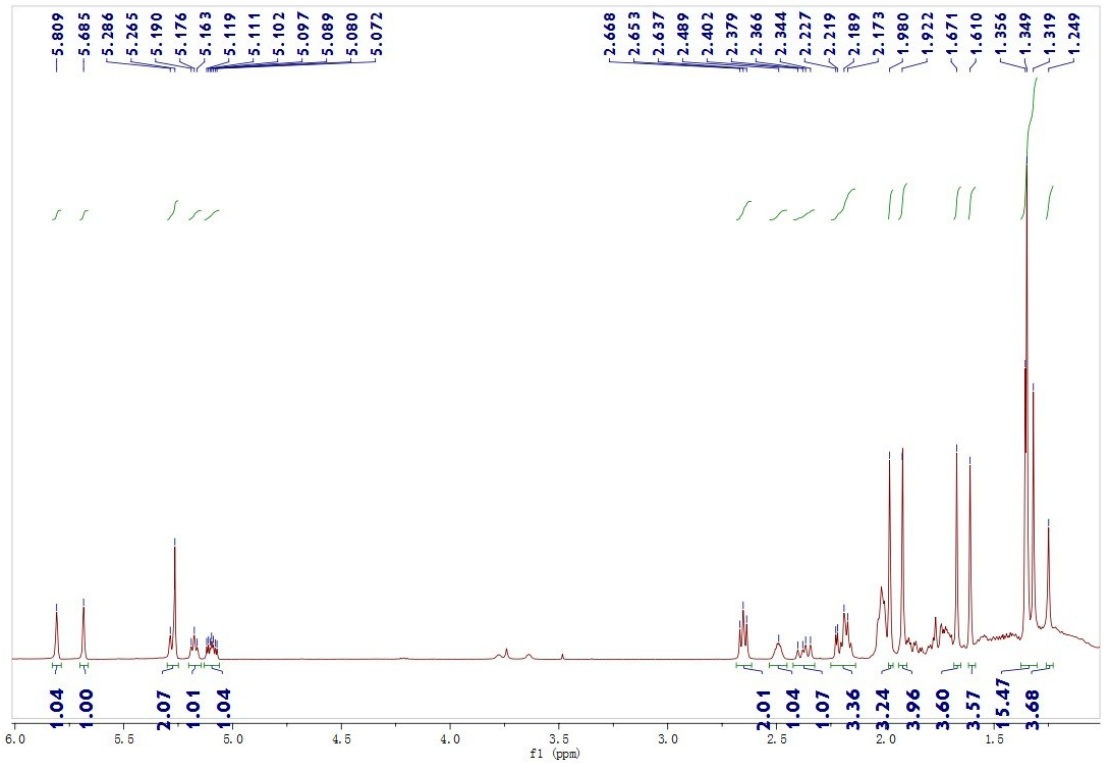
S31. CD spectrum of compound aphanamene G (**5**) in CH_3OH



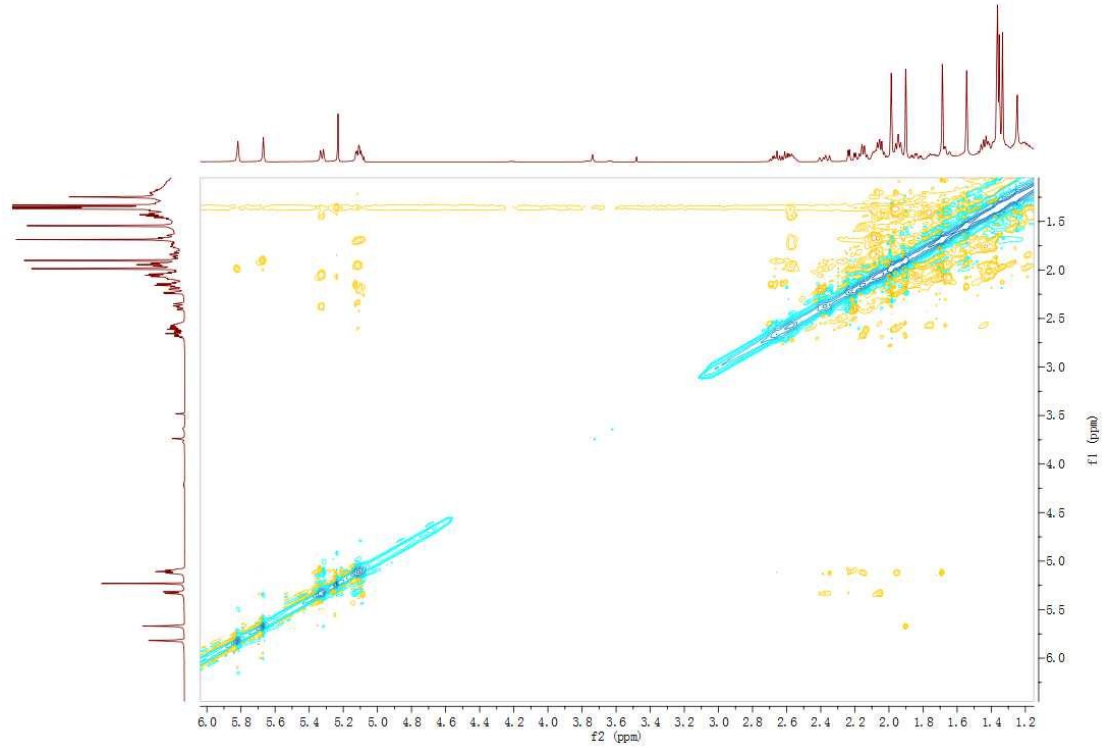
Elemental Composition Calculator

Target m/z:	685.3704	**Result type:**	Positive ions	**Species:**	$[M+Na]^+$
Elements:		C (0-80); H (0-120); O (0-30); N (0-10); Na (0-5); Cl (0-5)			
Ion Formula		**Calculated m/z**		**PPM Error**	
C40H52NaO8		685.3711		0.99	

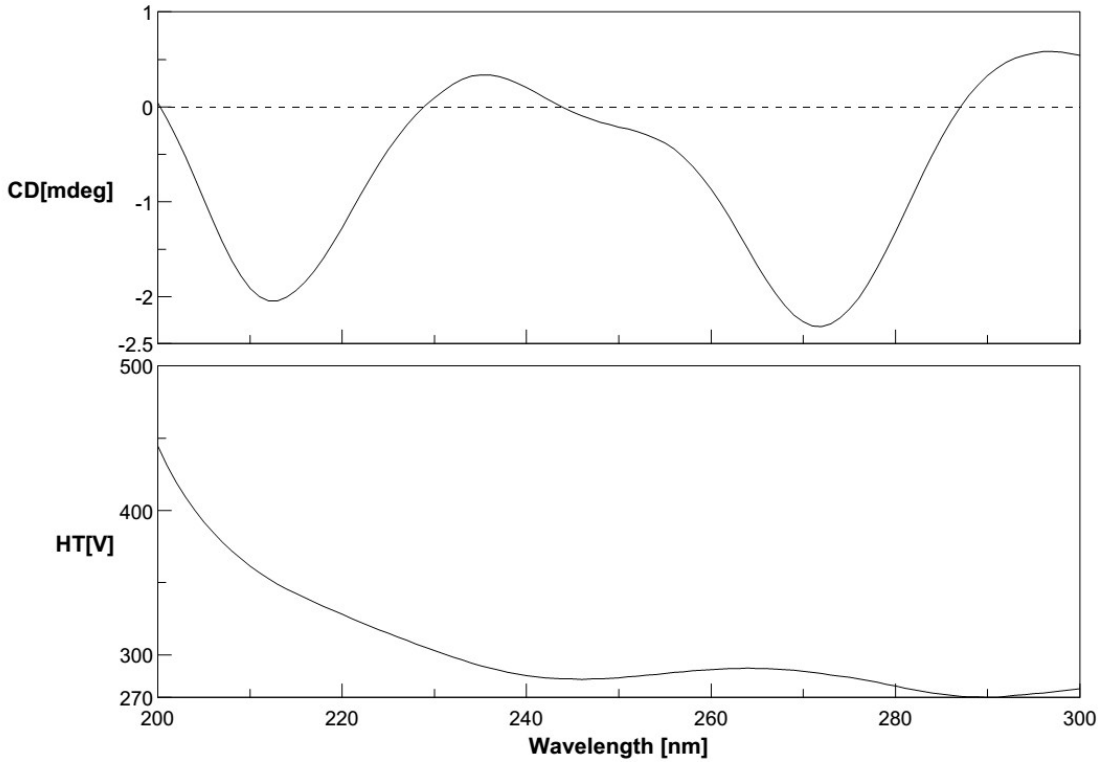
S32. HRESIMS spectrum of compound aphanamene G (**5**) in CH_3OH



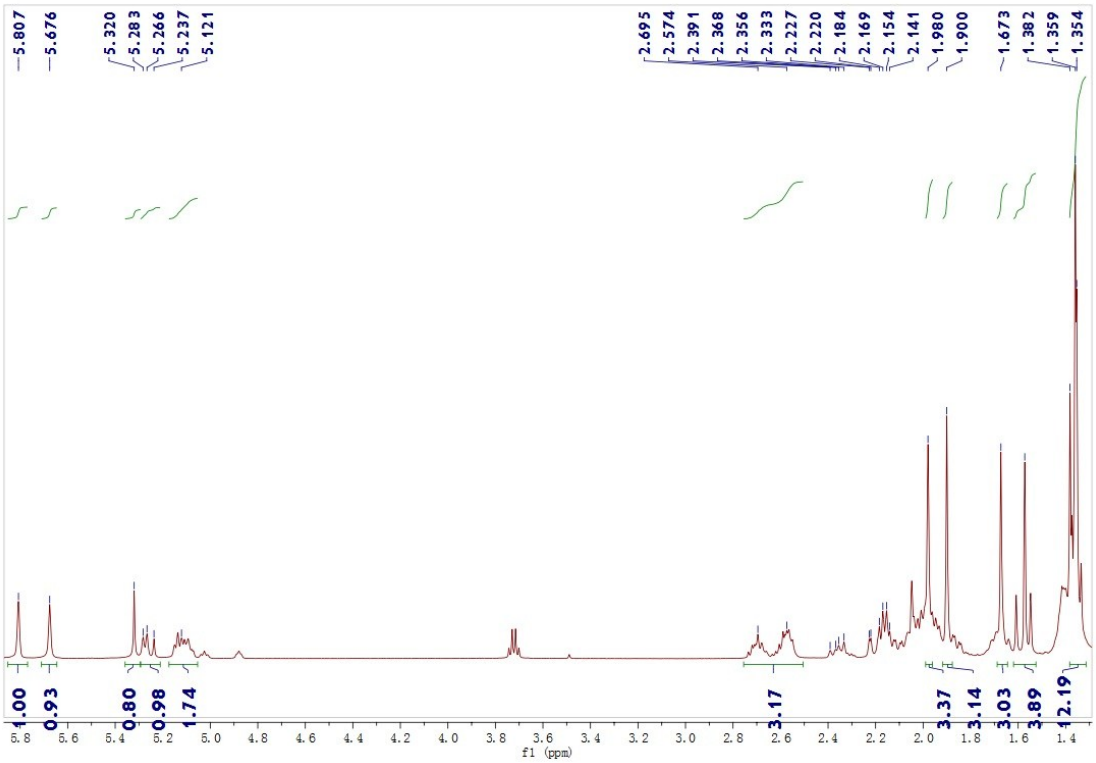
S33. ^{1}H NMR spectrum of compound aphanamene H (**6**) in $CDCl_3$



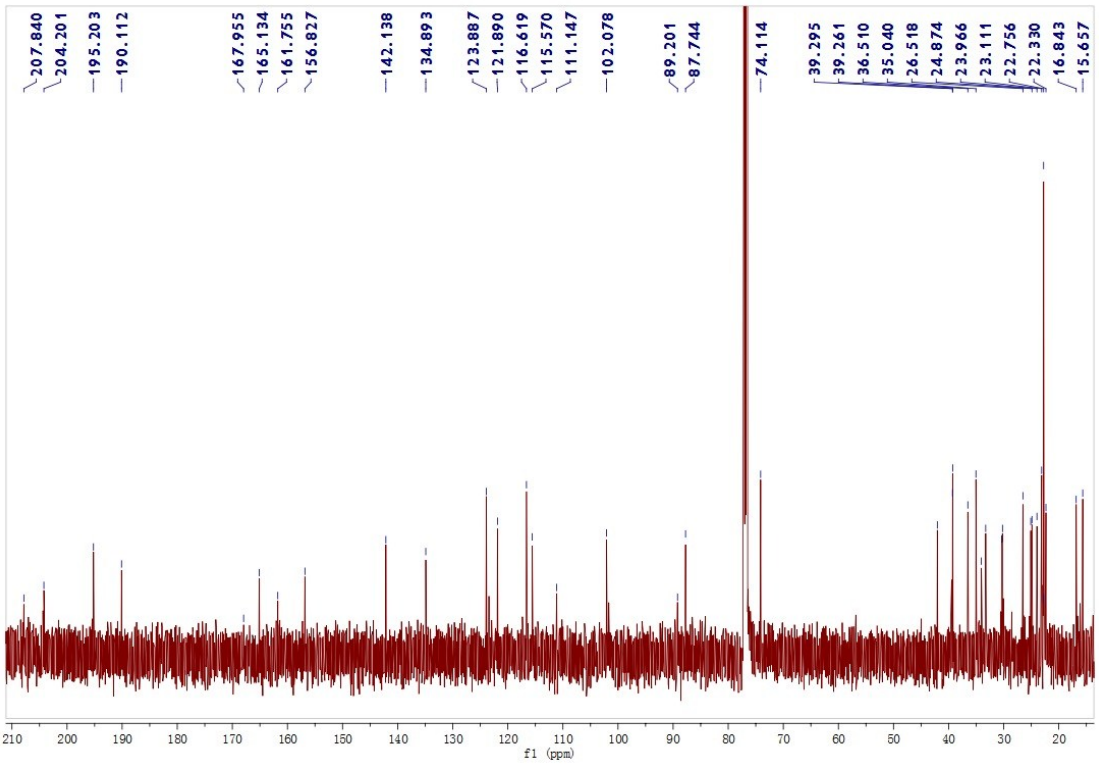
S34. ROESY spectrum of compound aphanamene H (**6**) in $CDCl_3$



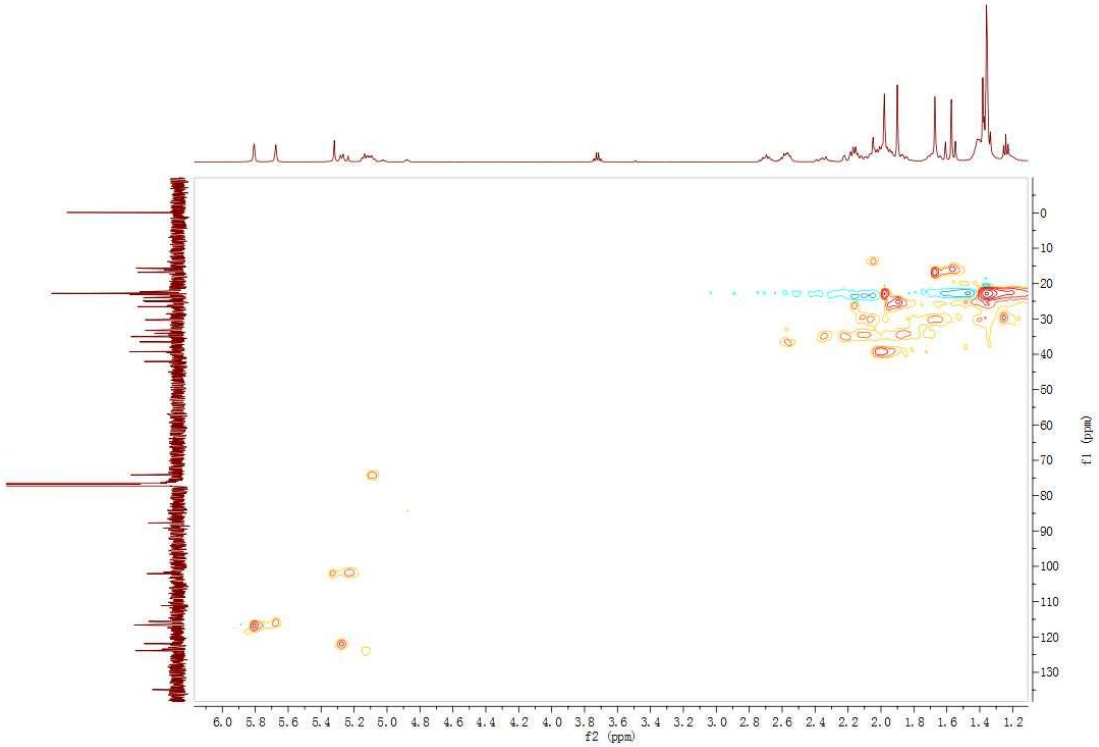
S35. CD spectrum of compound aphanamene H (**6**) in CH_3OH



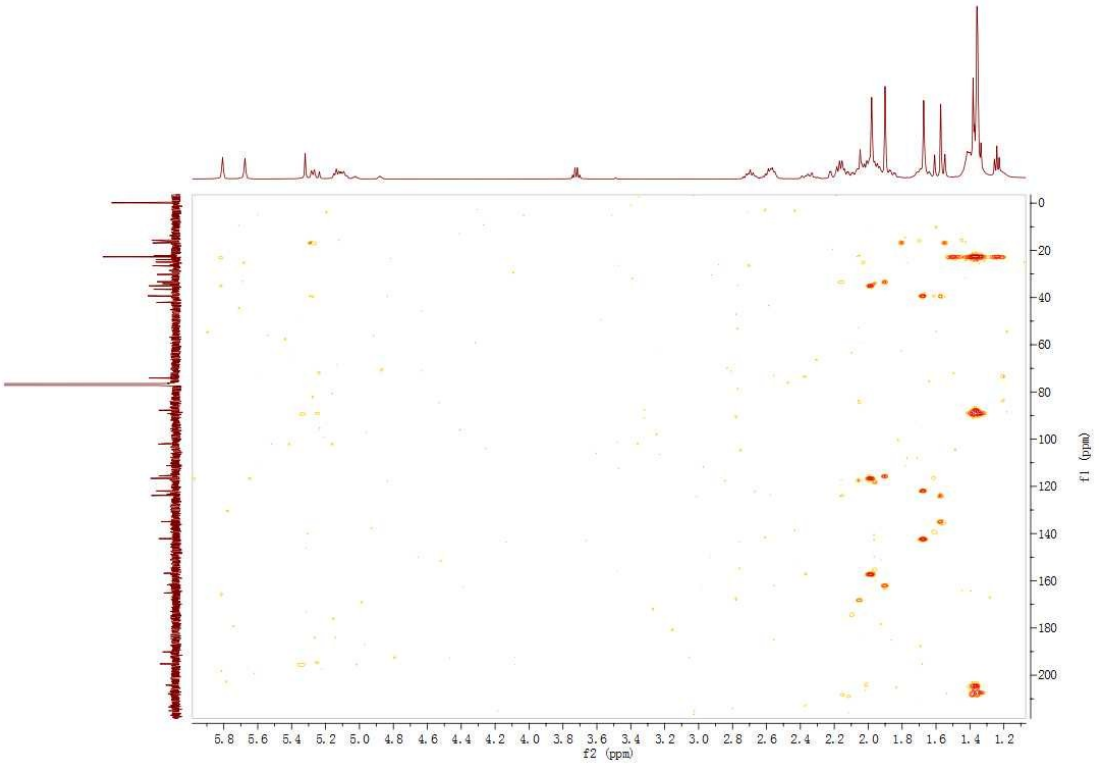
S36. ^1H NMR spectrum of compound aphanamene I (**7**) in $CDCl_3$



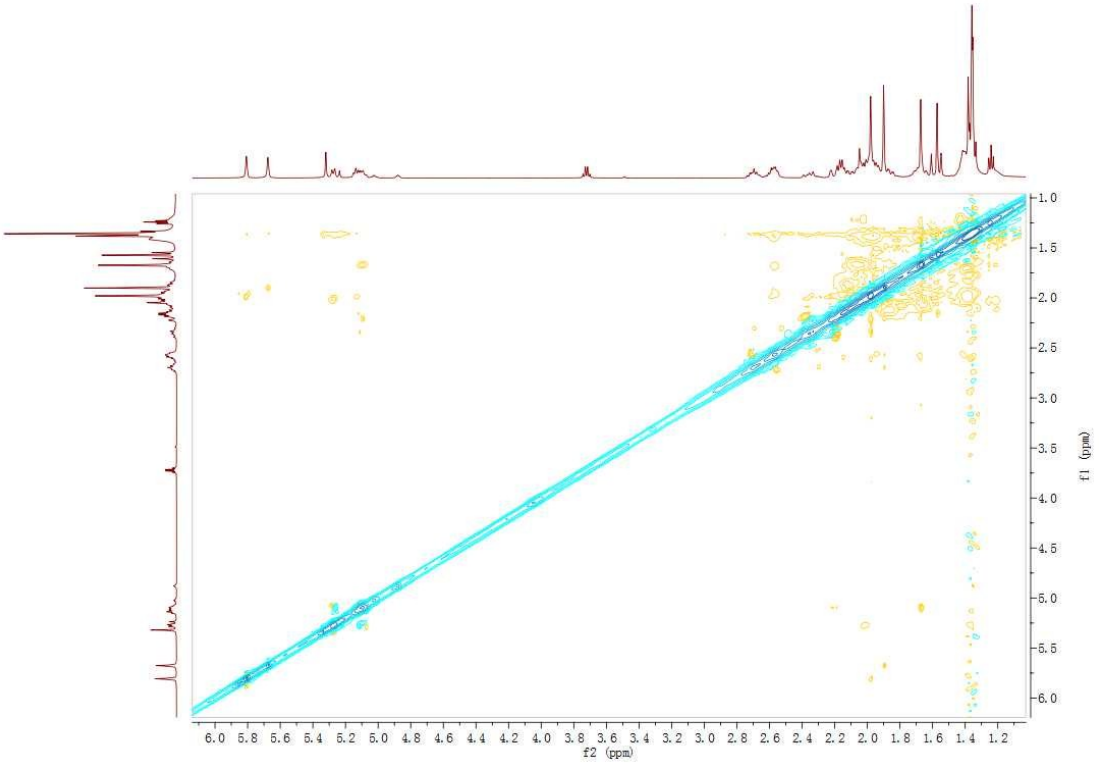
S37. ^{13}C NMR spectrum of compound aphanamene I (**7**) in $CDCl_3$



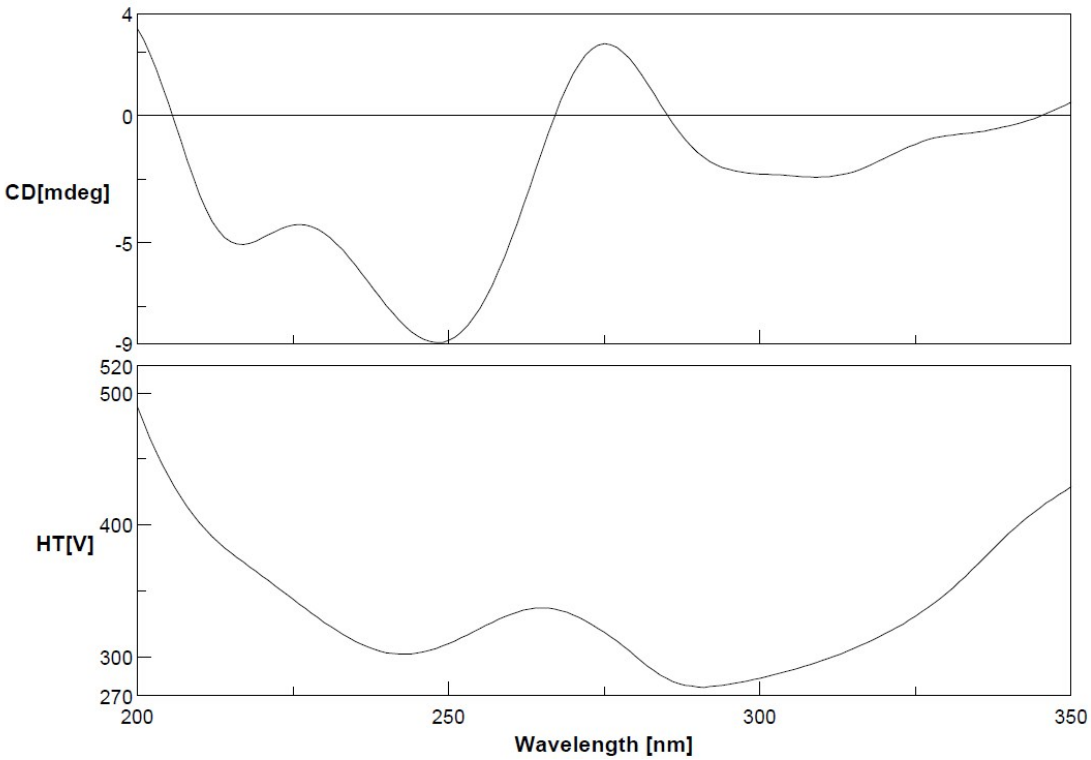
S38. HSQC spectrum of compound aphanamene I (**7**) in $CDCl_3$



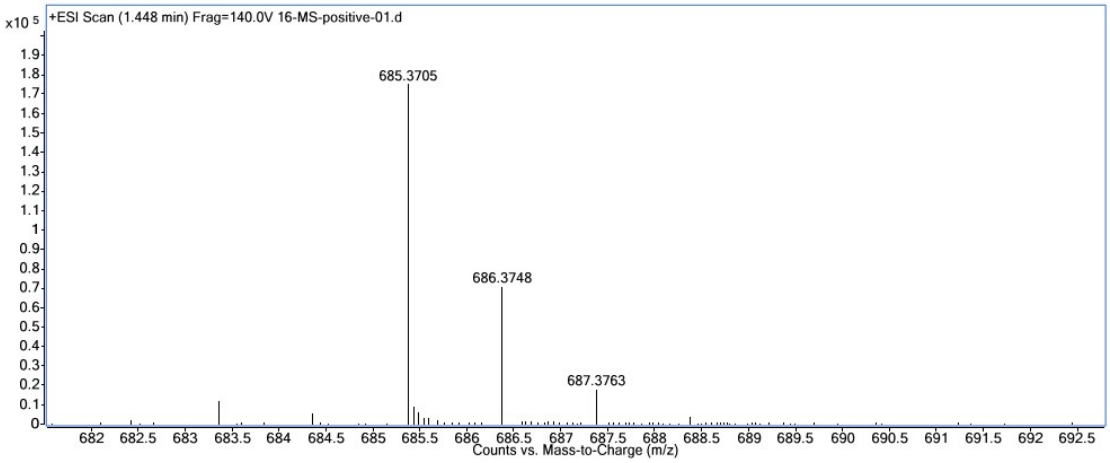
S39. HMBC spectrum of compound aphanamene I (**7**) in $CDCl_3$



S40. ROESY spectrum of compound aphanamene I (**7**) in $CDCl_3$



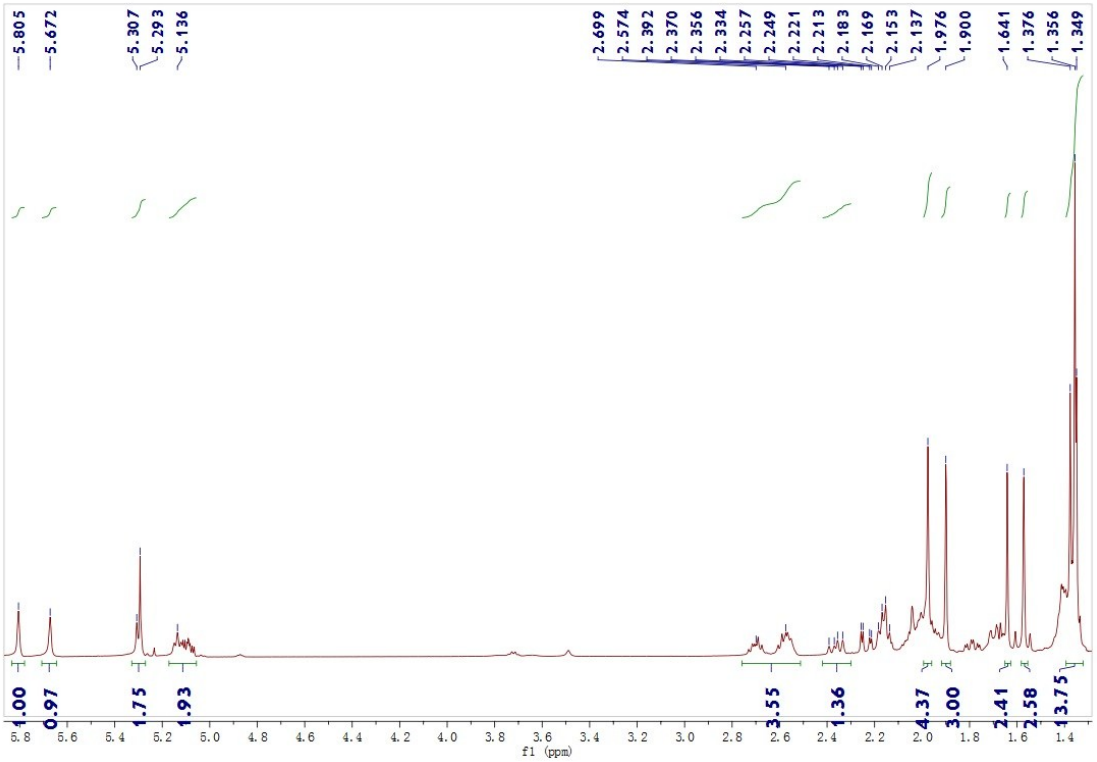
S41. CD spectrum of compound aphanamene I (**7**) in CH_3OH



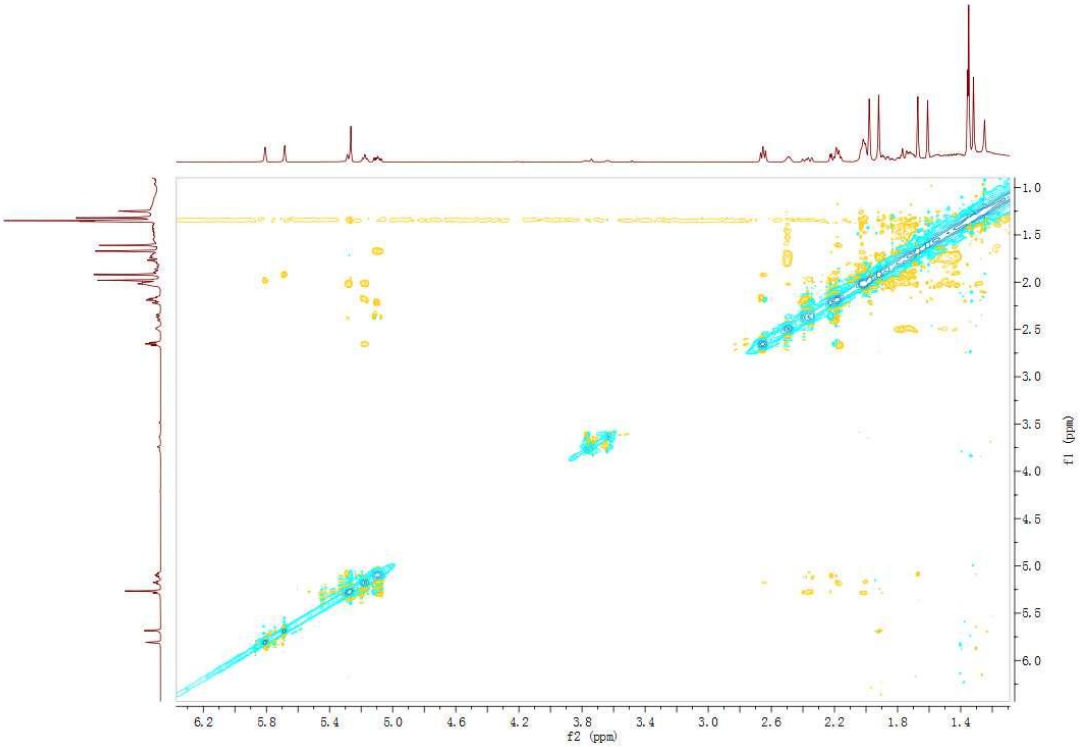
Elemental Composition Calculator

Target m/z:	685.3705	**Result type:**	Positive ions	**Species:**	$[M+Na]^+$
Elements:		C (0-80); H (0-120); O (0-30); N (0-10); Na (0-5); Cl (0-5)			
Ion Formula		**Calculated m/z**		**PPM Error**	
C40H52NaO8		685.3711		0.79	

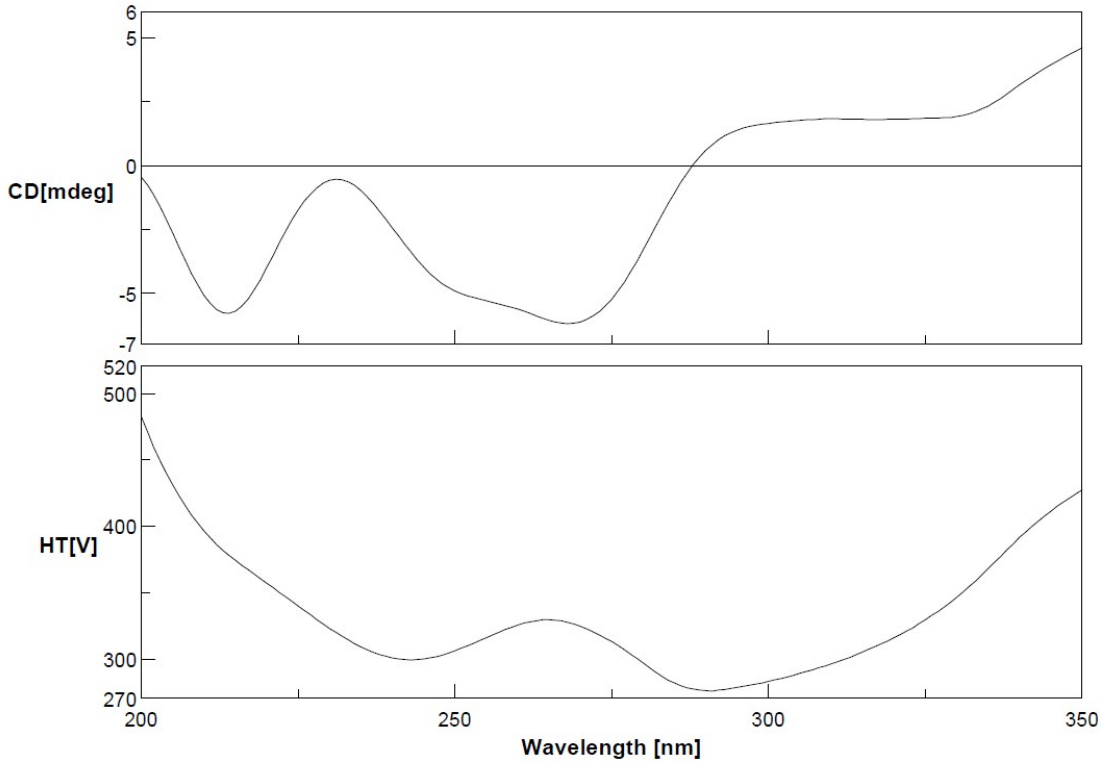
S42. HRESIMS spectrum of compound aphanamene I (**7**) in CH_3OH



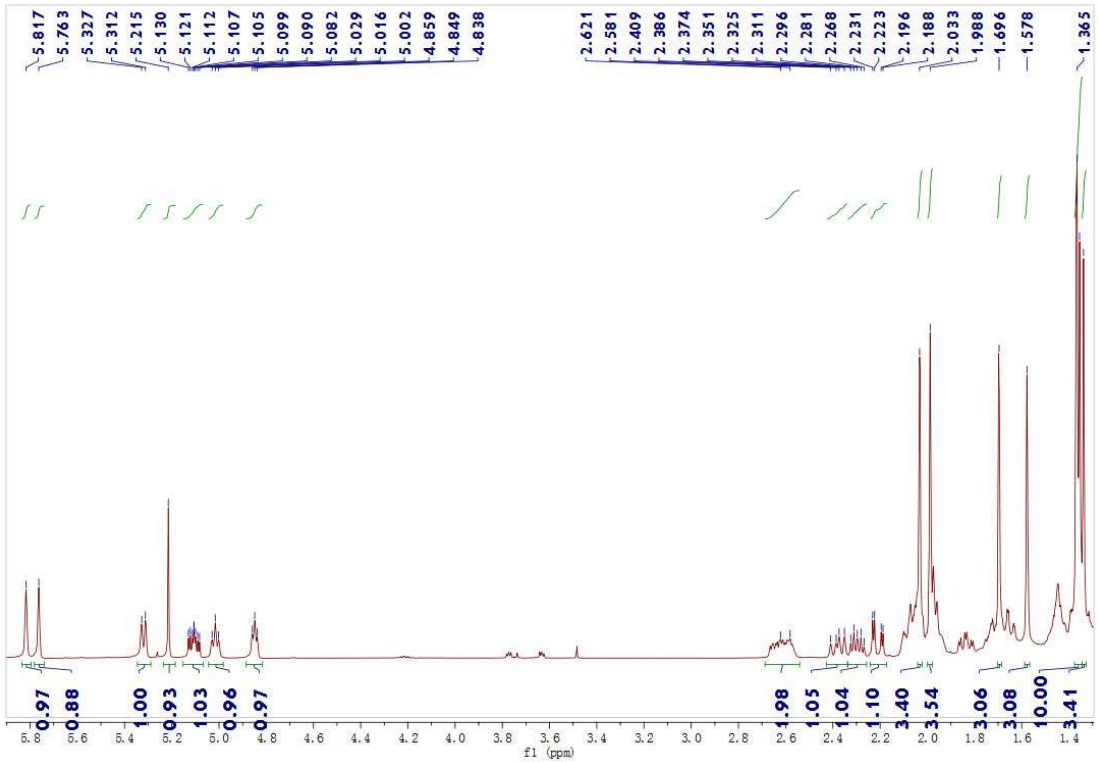
S43. ^{1}H NMR spectrum of compound aphanamene J (**8**) in $CDCl_3$



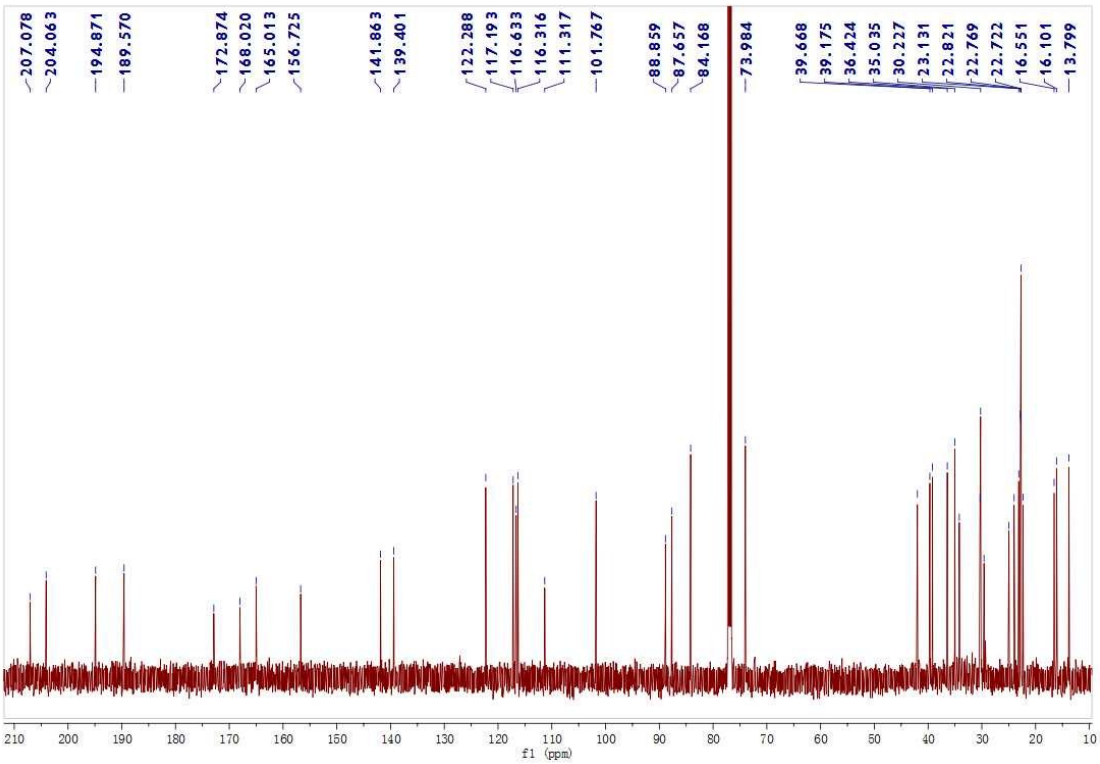
S44. ROESY spectrum of compound aphanamene J (**8**) in $CDCl_3$



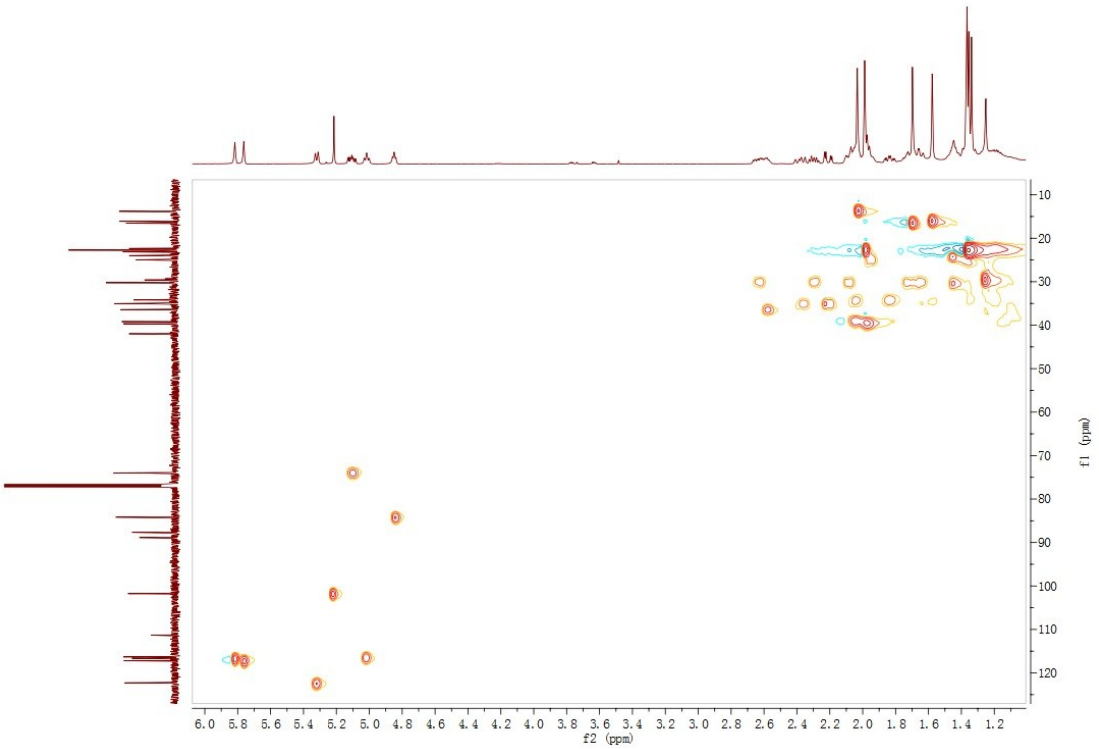
S45. CD spectrum of compound aphanamene J (**8**) in CH_3OH



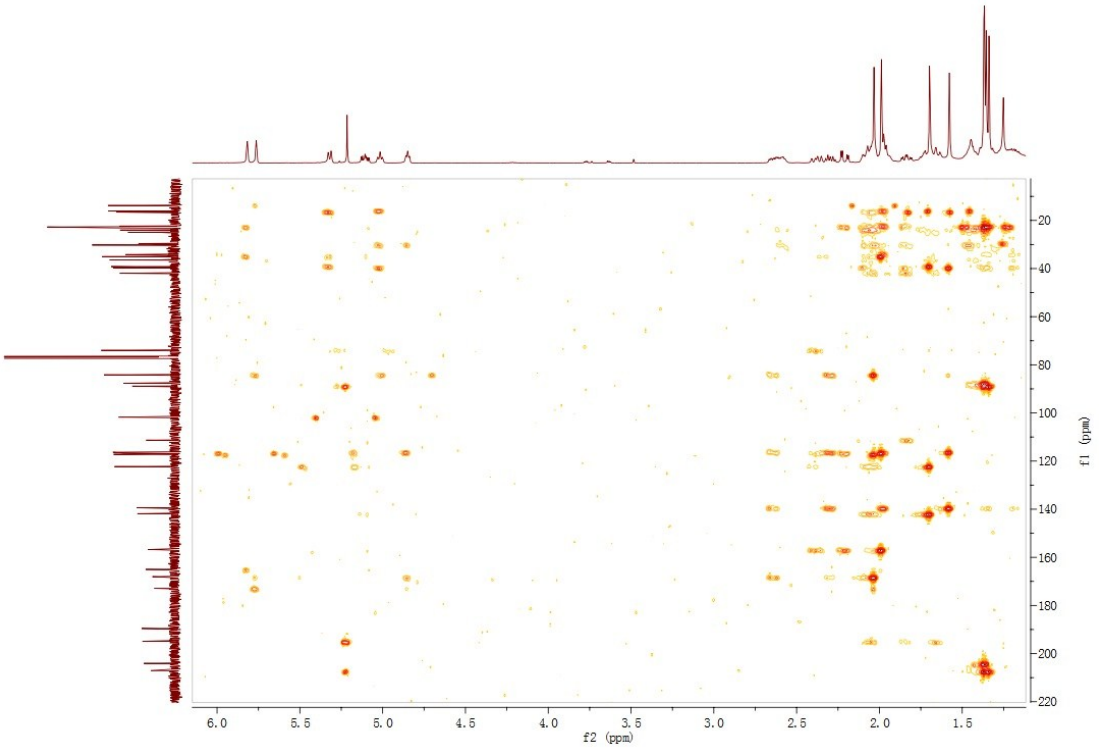
S46. ^1H NMR spectrum of compound aphanamene K (**9**) in $CDCl_3$



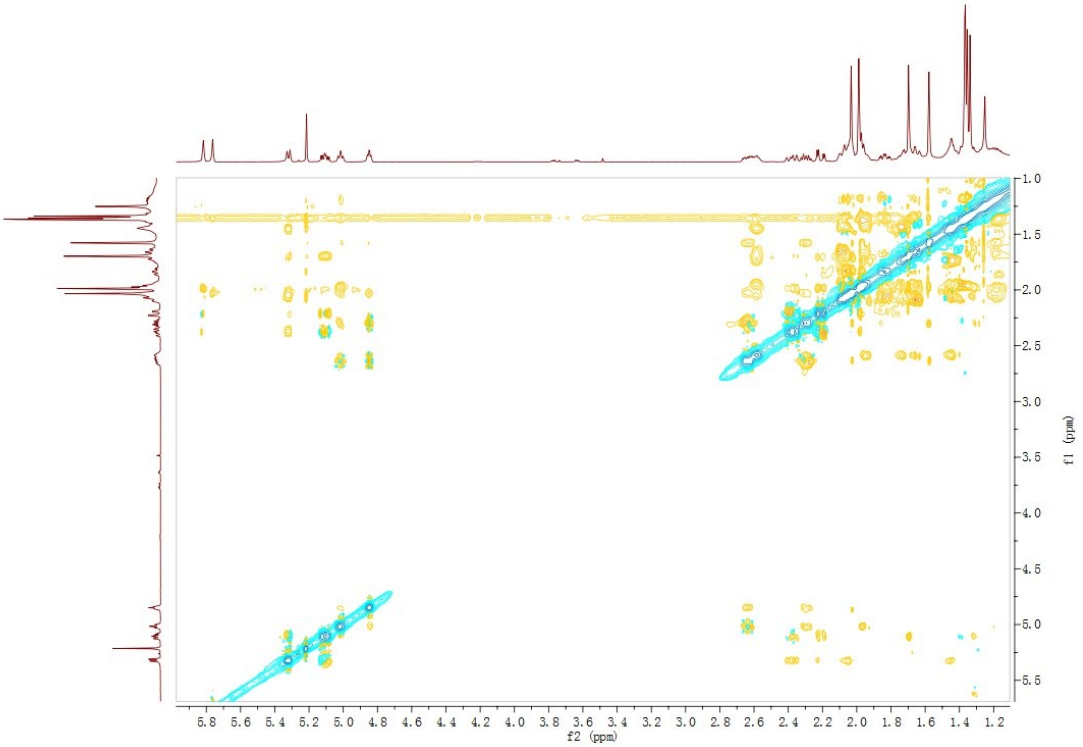
S47. ^{13}C NMR spectrum of compound aphanamene K (**9**) in $CDCl_3$



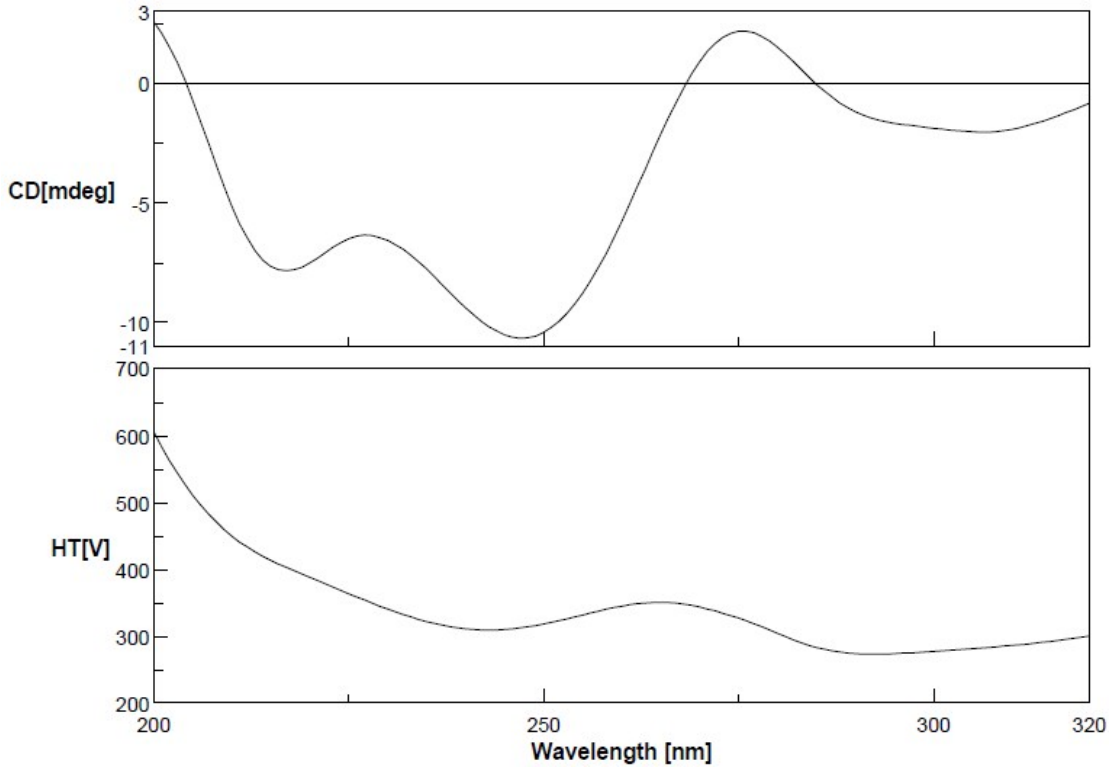
S48. HSQC spectrum of compound aphanamene K (**9**) in $CDCl_3$



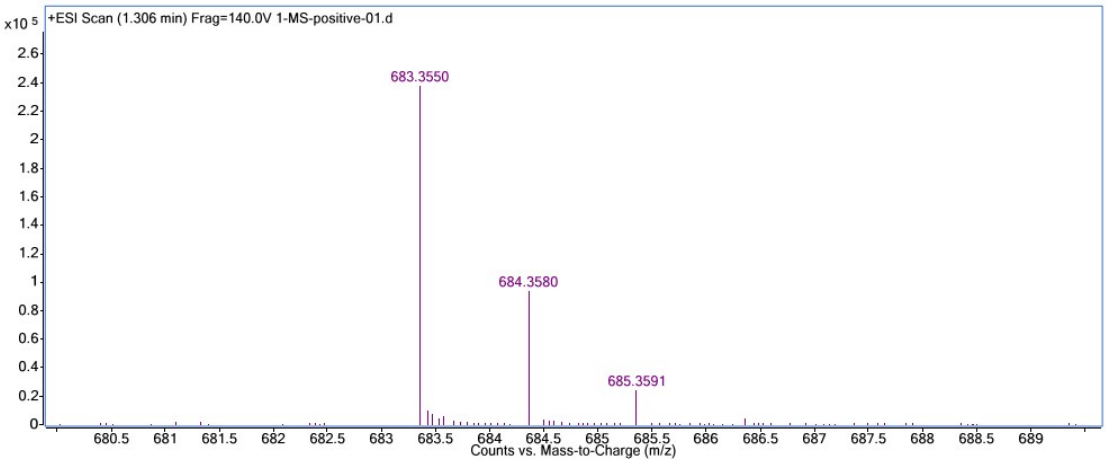
S49. HMBC spectrum of compound aphanamene K (**9**) in $CDCl_3$



S50. ROESY spectrum of compound aphanamene K (**9**) in $CDCl_3$



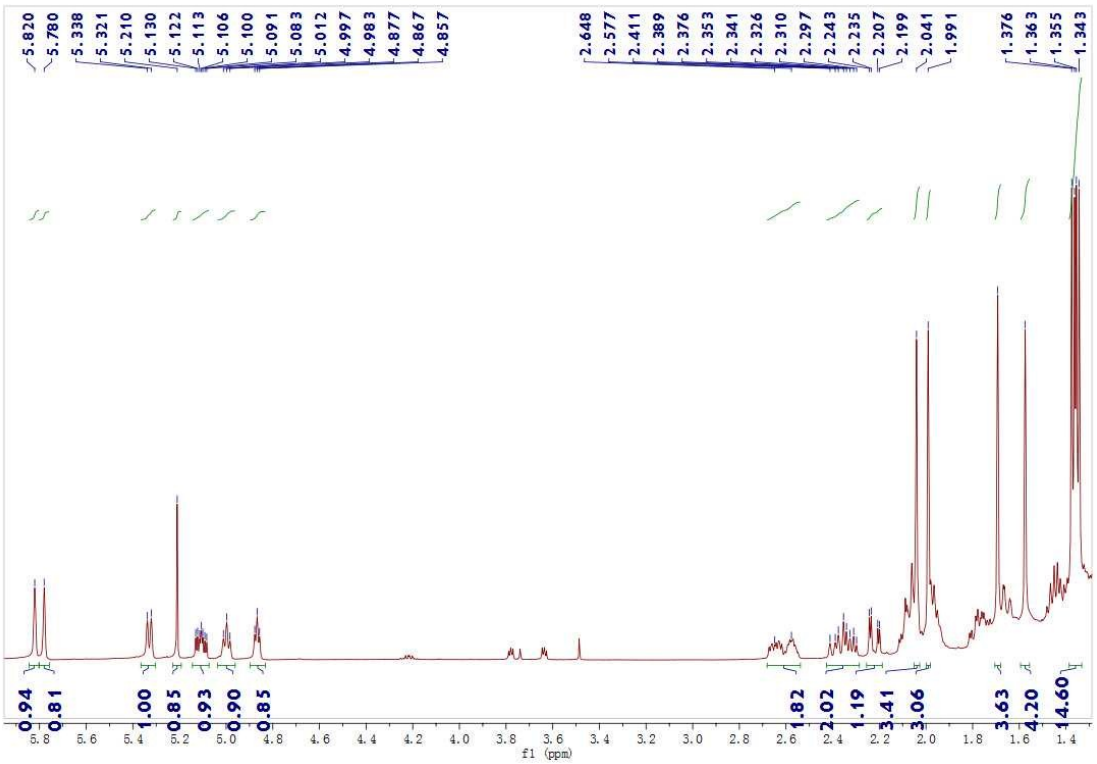
S51. CD spectrum of compound aphanamene K (**9**) in CH_3OH



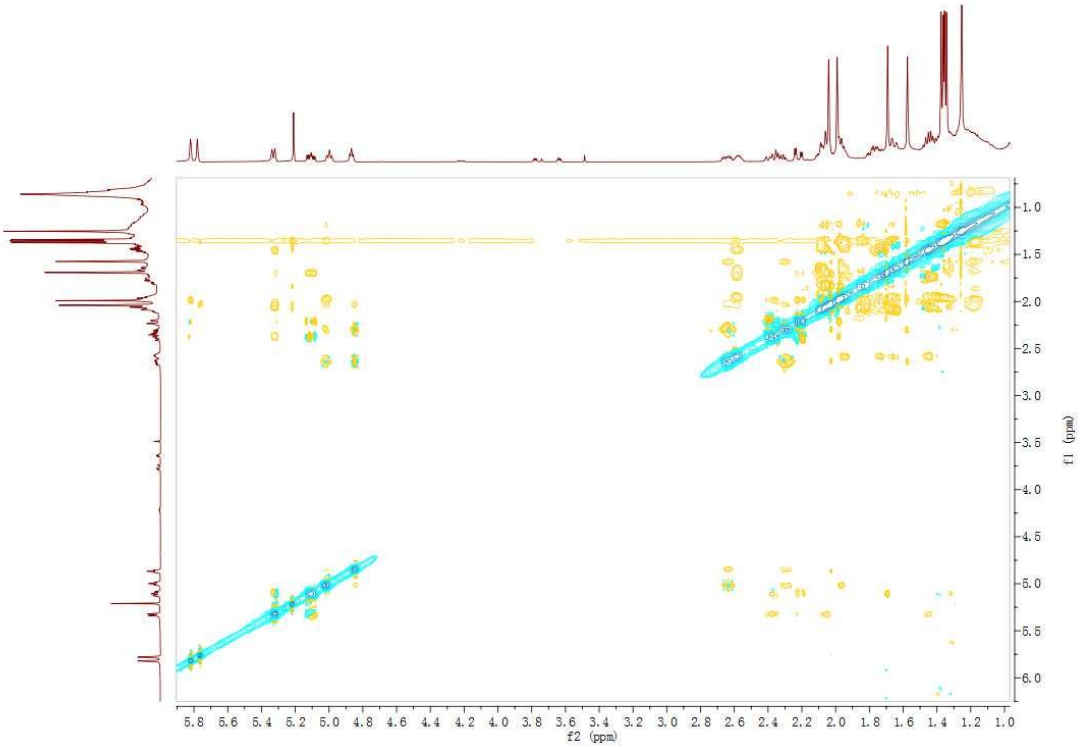
Elemental Composition Calculator

Target m/z:	683.3550	**Result type:**	Positive ions	**Species:**	$[M+Na]^+$
Elements:		C (0-80); H (0-120); O (0-30); N (0-10); Na (0-5); Cl (0-5)			
Ion Formula		**Calculated m/z**		**PPM Error**	
C40H52NaO8		683.3554		0.62	

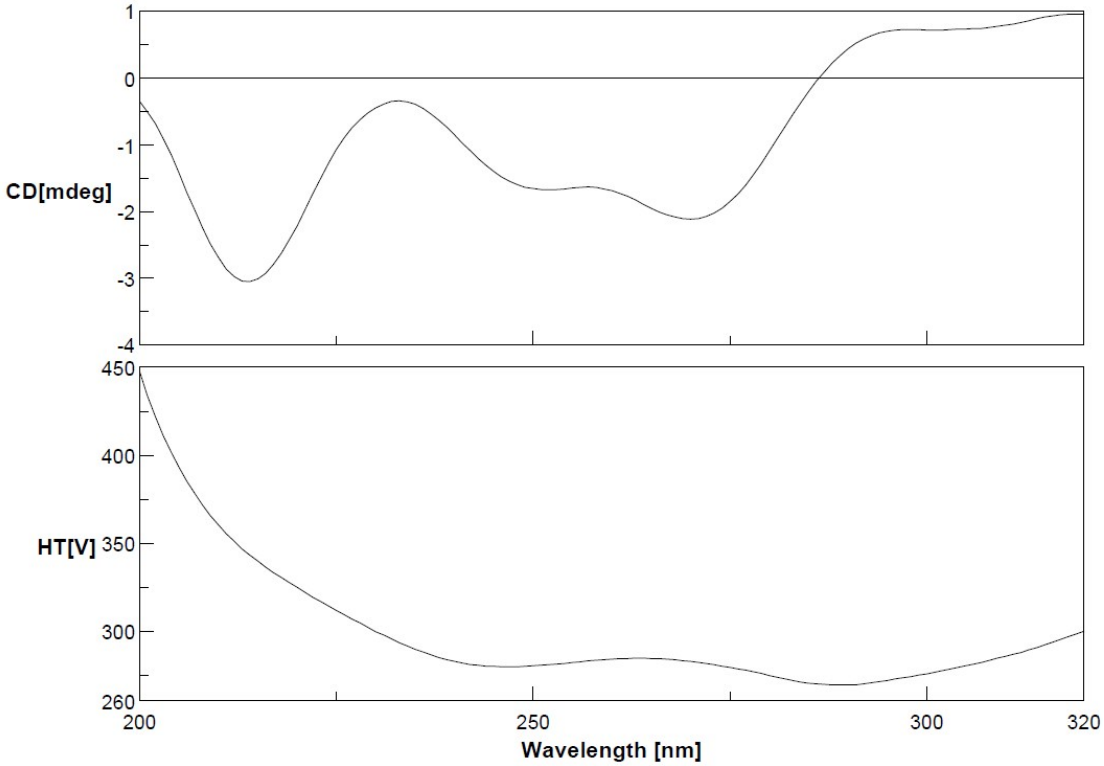
S52. HRESIMS spectrum of compound aphanamene K (**9**) in CH_3OH



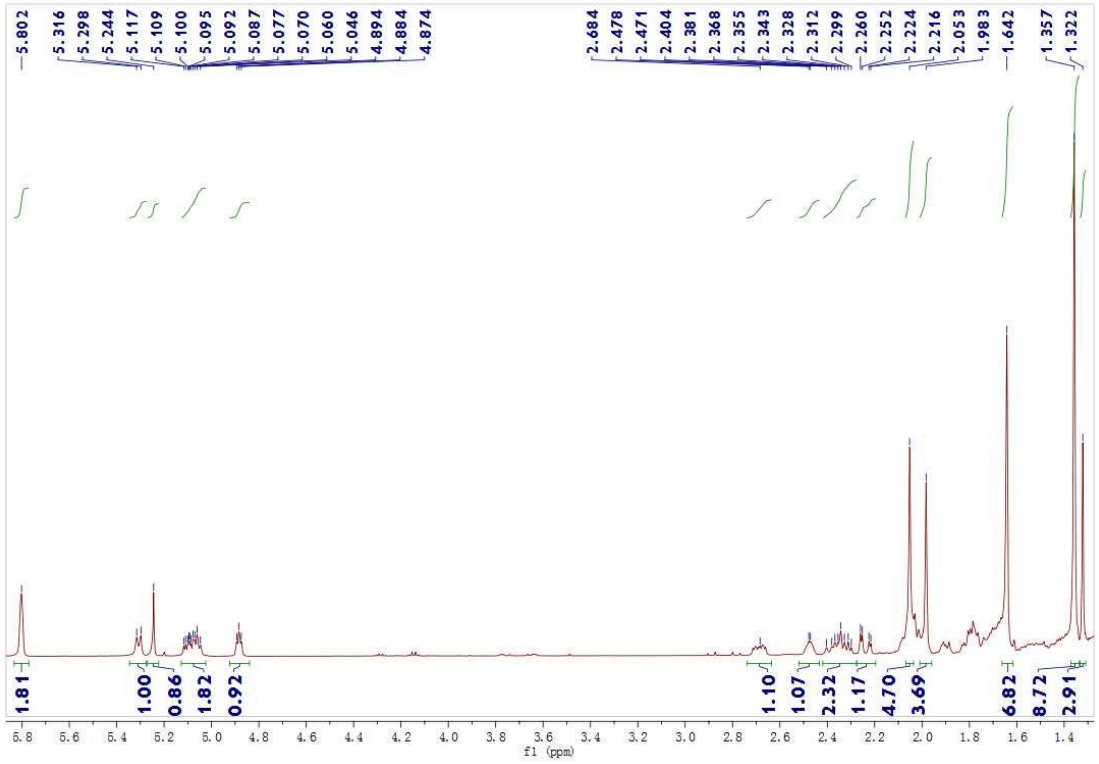
S53. ^{1}H NMR spectrum of compound aphanamene L (**10**) in $CDCl_3$



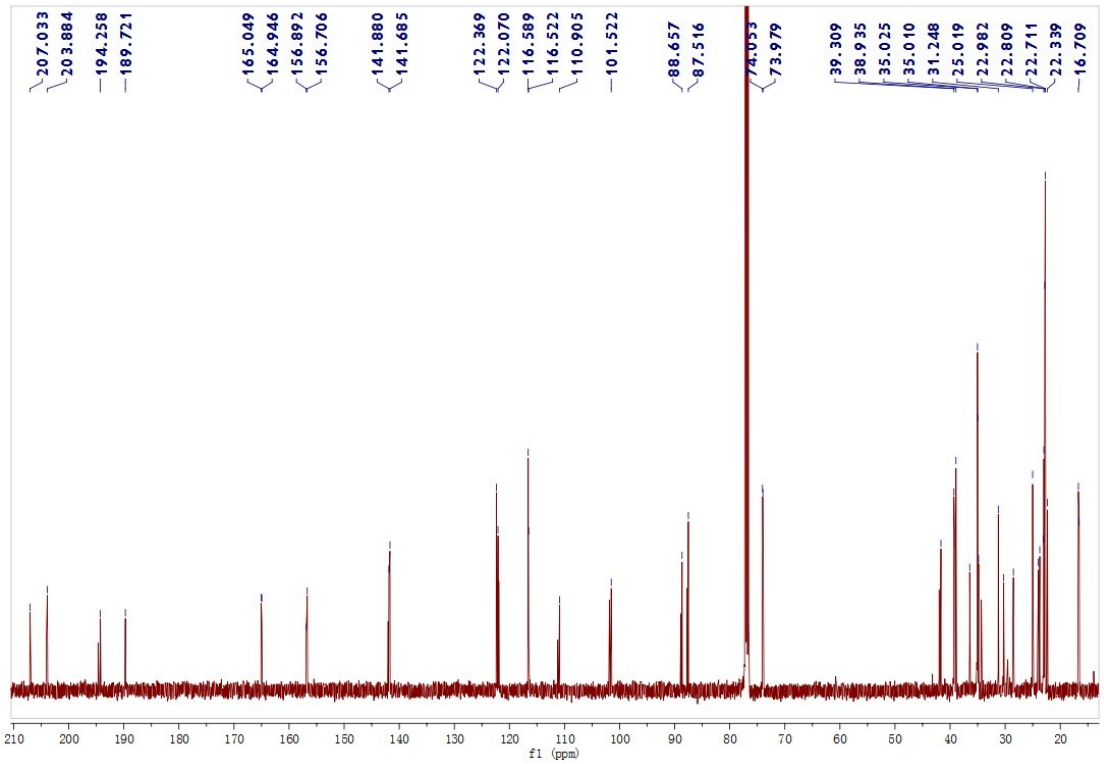
S54. ROESY spectrum of compound aphanamene L (**10**) in $CDCl_3$



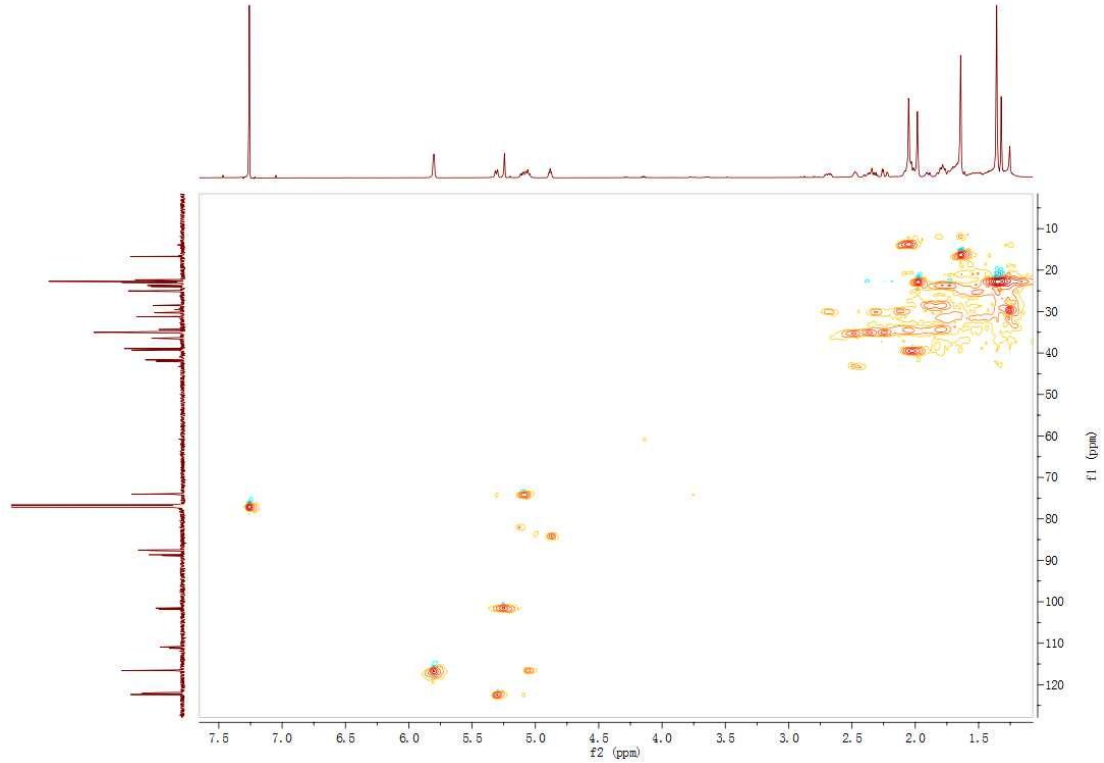
S55. CD spectrum of compound aphanamene L (**10**) in CH_3OH



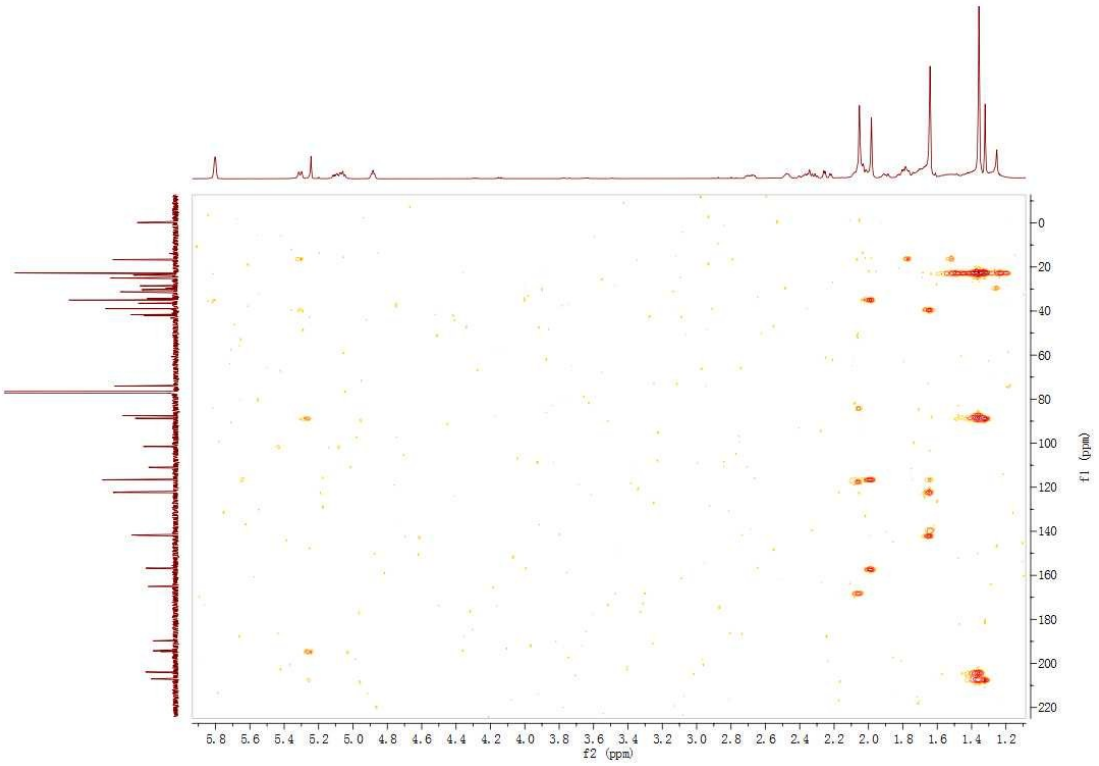
S56. ^1H NMR spectrum of compound aphanamene M (**11**) in $CDCl_3$



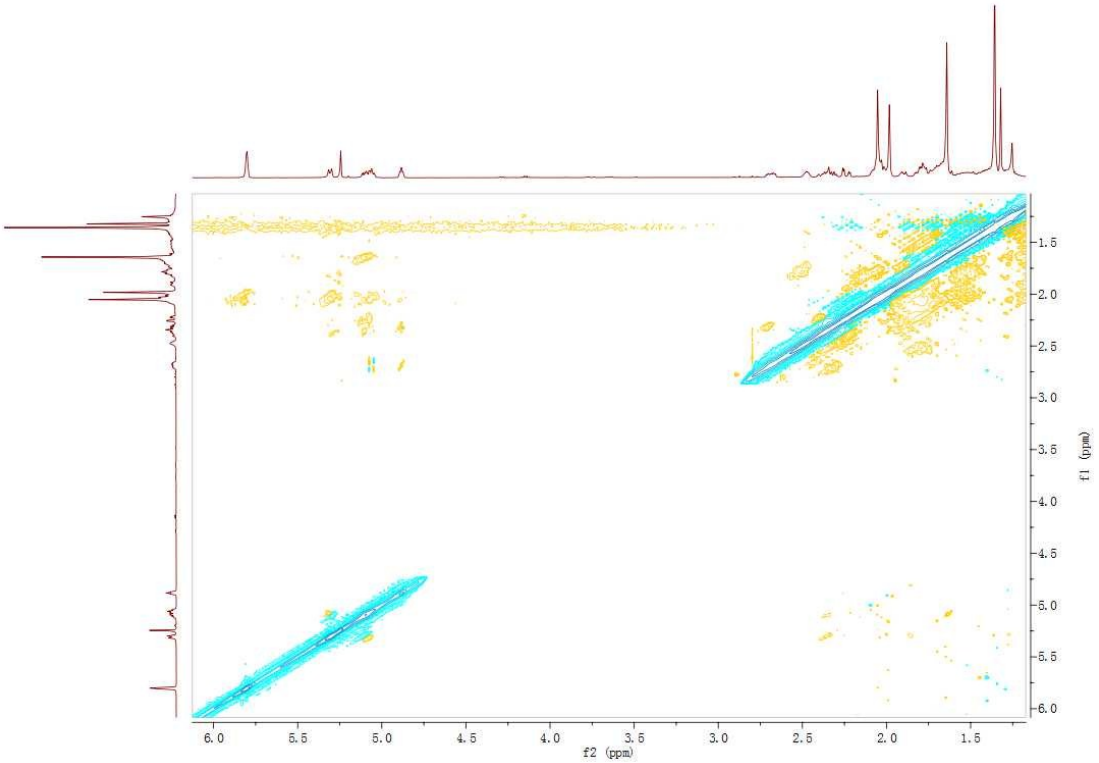
S57. ^{13}C NMR spectrum of compound aphanamene M (**11**) in $CDCl_3$



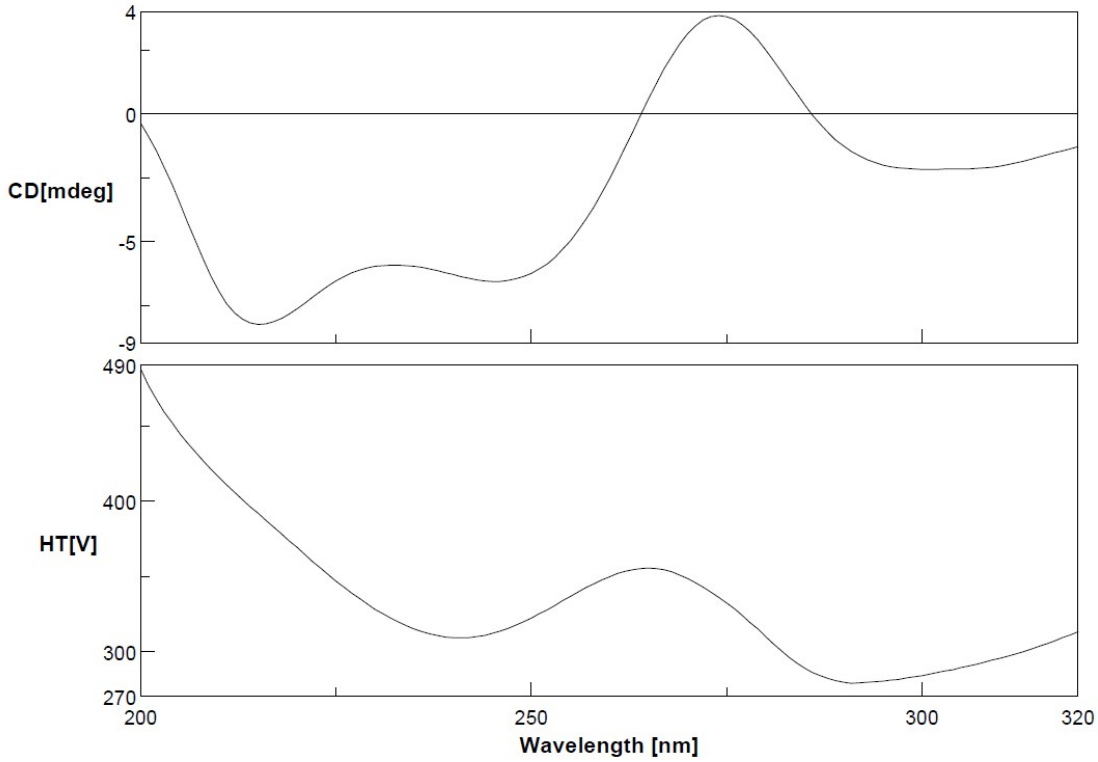
S58. HSQC spectrum of compound aphanamene M (**11**) in $CDCl_3$



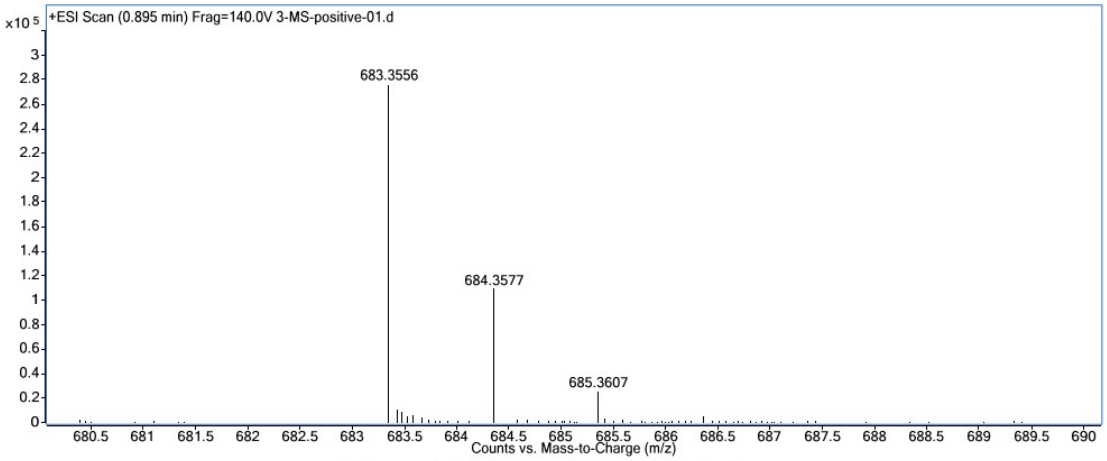
S59. HMBC spectrum of compound aphanamene M (**11**) in $CDCl_3$



S60. ROESY spectrum of compound aphanamene M (**11**) in $CDCl_3$



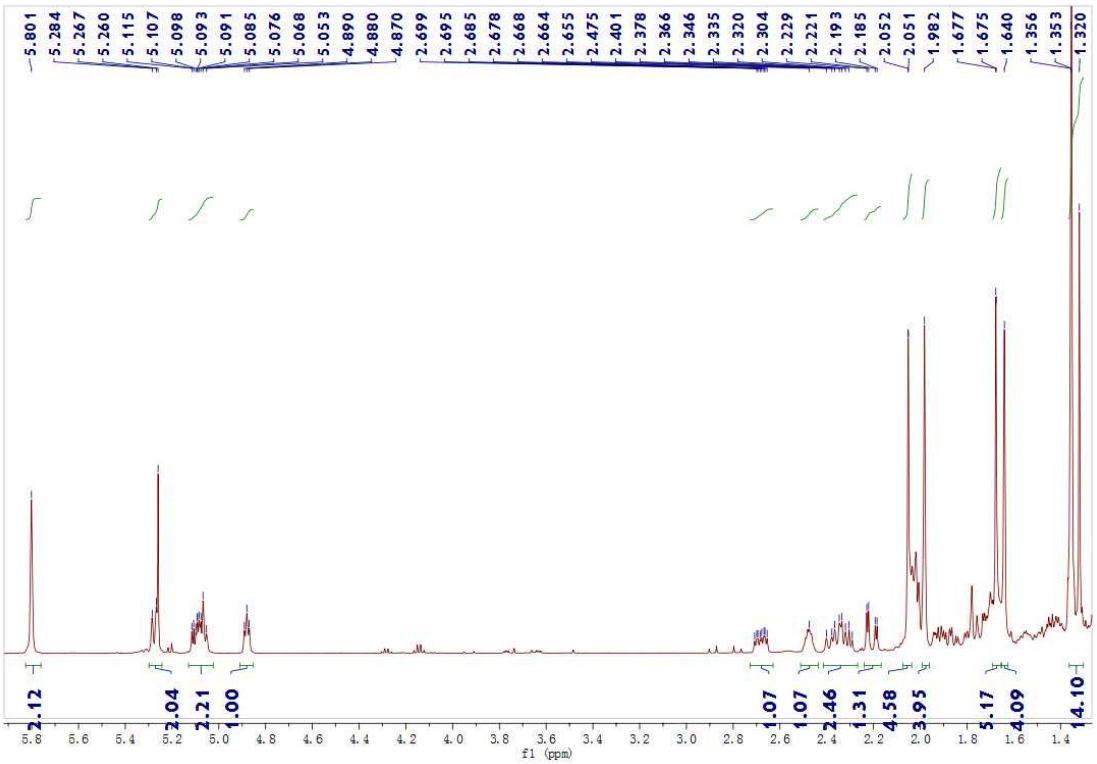
S61. CD spectrum of compound aphanamene M (**11**) in CH_3OH



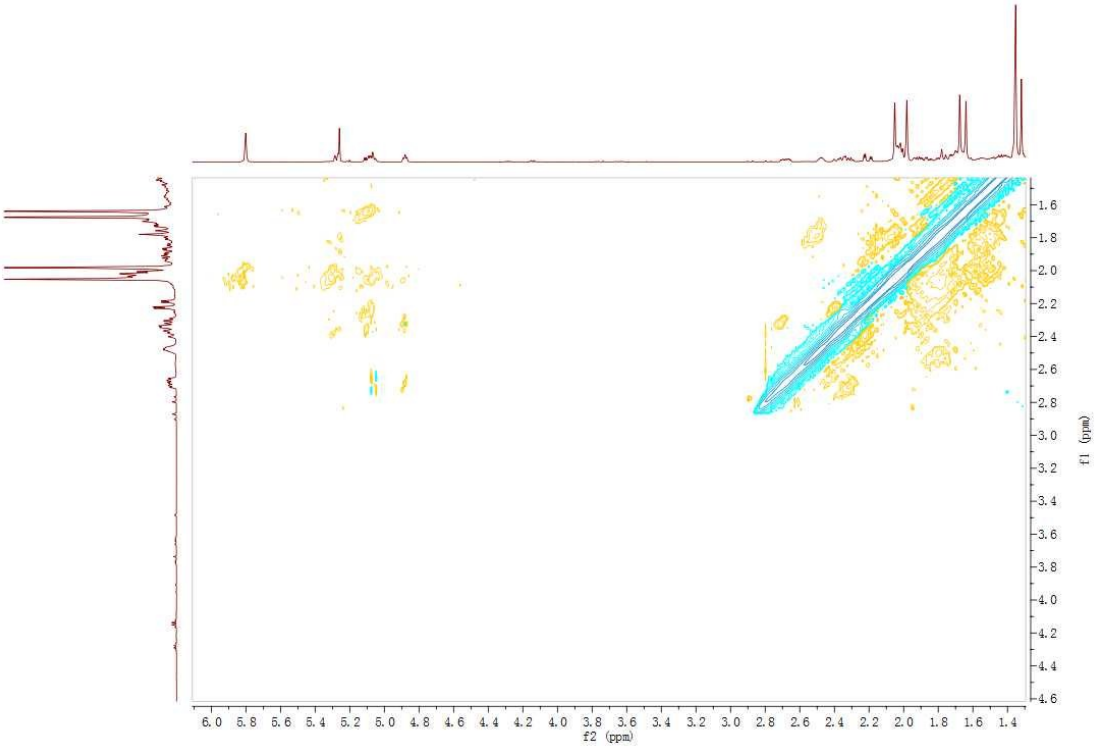
Elemental Composition Calculator

Target m/z:	683.3556	**Result type:**	Positive ions	**Species:**	$[M+Na]^+$
Elements:		C (0-80); H (0-120); O (0-30); N (0-10); Na (0-5); Cl (0-5)			
Ion Formula		**Calculated m/z**		**PPM Error**	
C40H52NaO8		683.3554		-0.21	

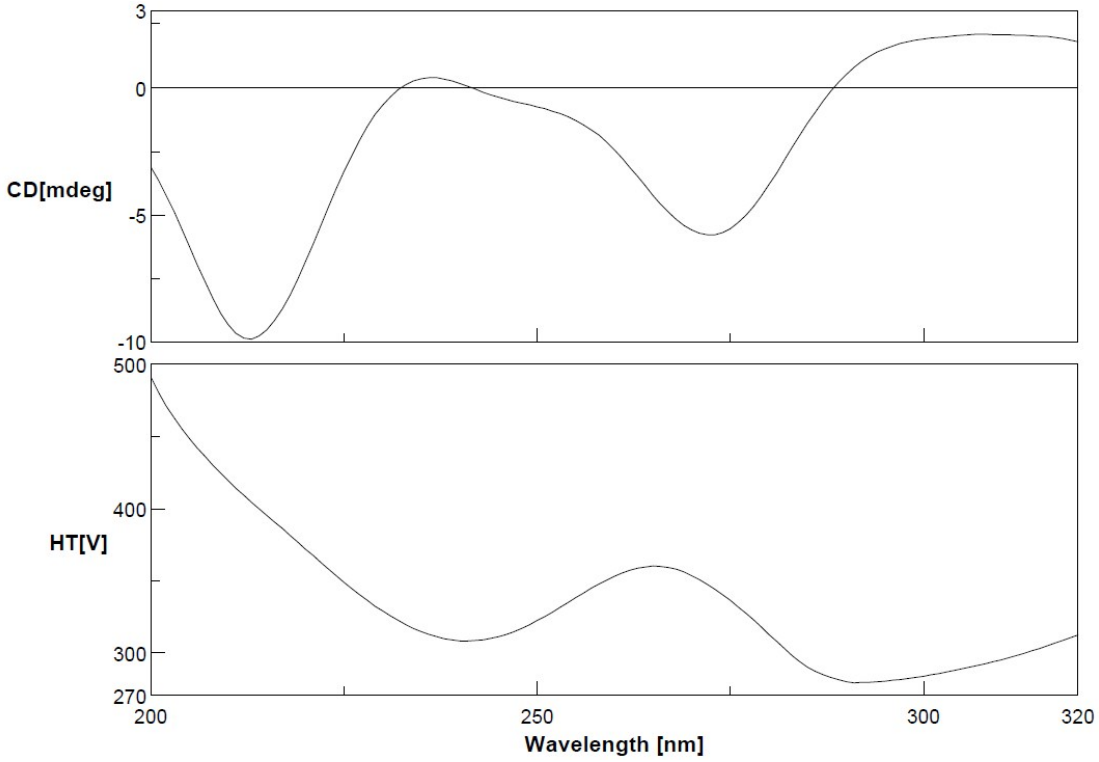
S62. HRESIMS spectrum of compound aphanamene M (**11**) in CH_3OH



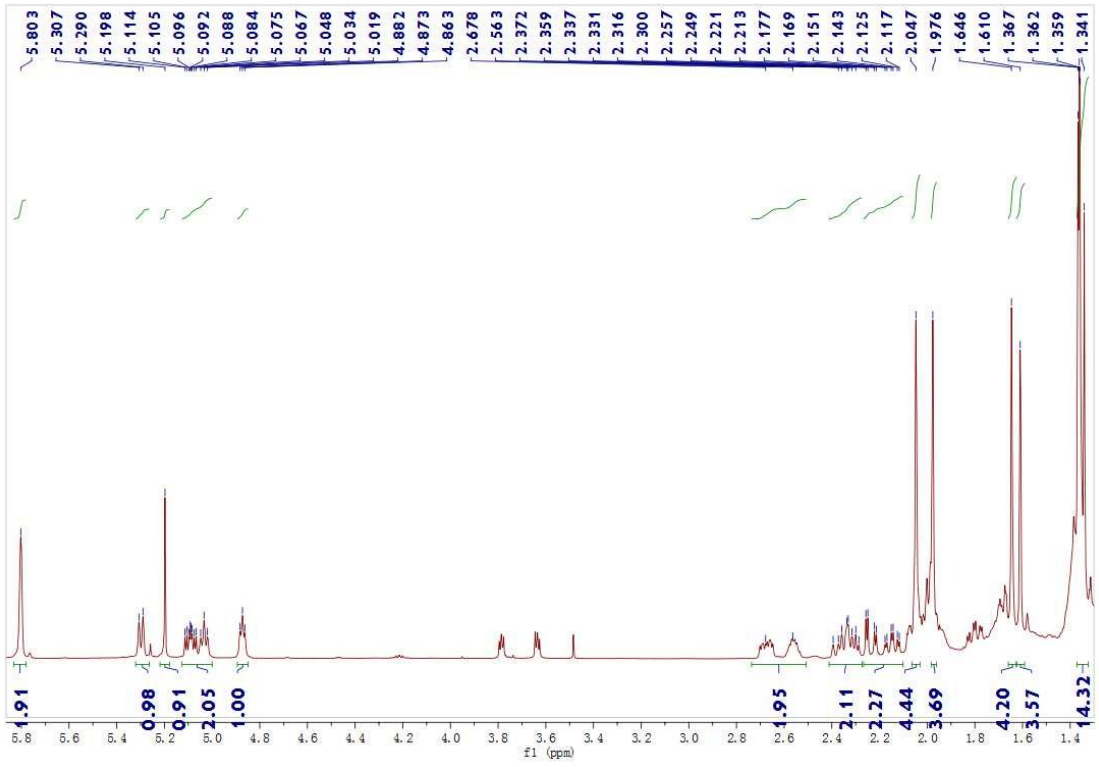
S63. ^{1}H NMR spectrum of compound aphanamene N (**12**) in $CDCl_3$



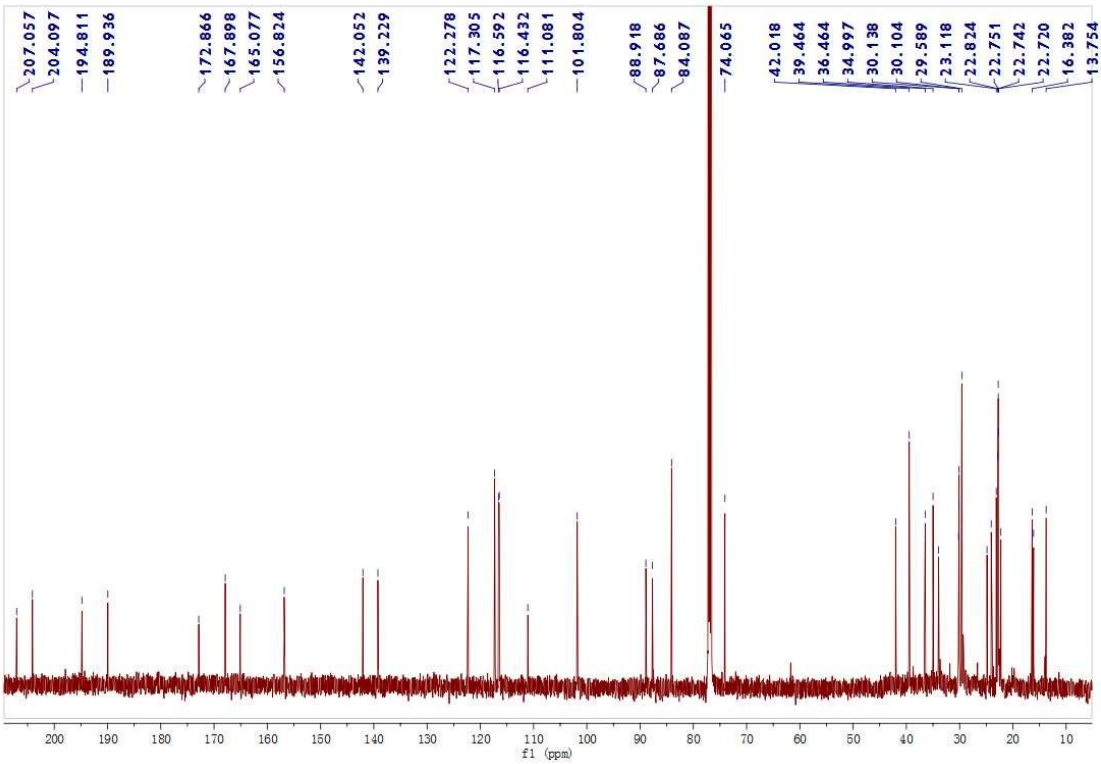
S64. ROESY spectrum of compound aphanamene N (**12**) in $CDCl_3$



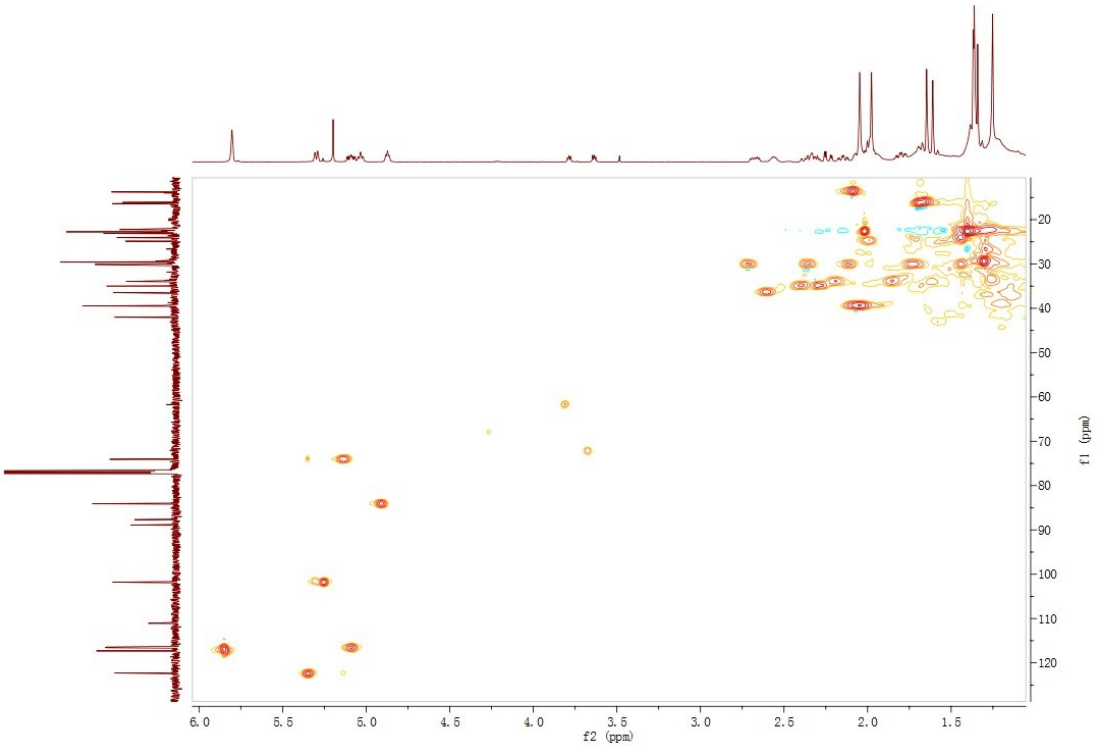
S65. CD spectrum of compound aphanamene N (**12**) in CH_3OH



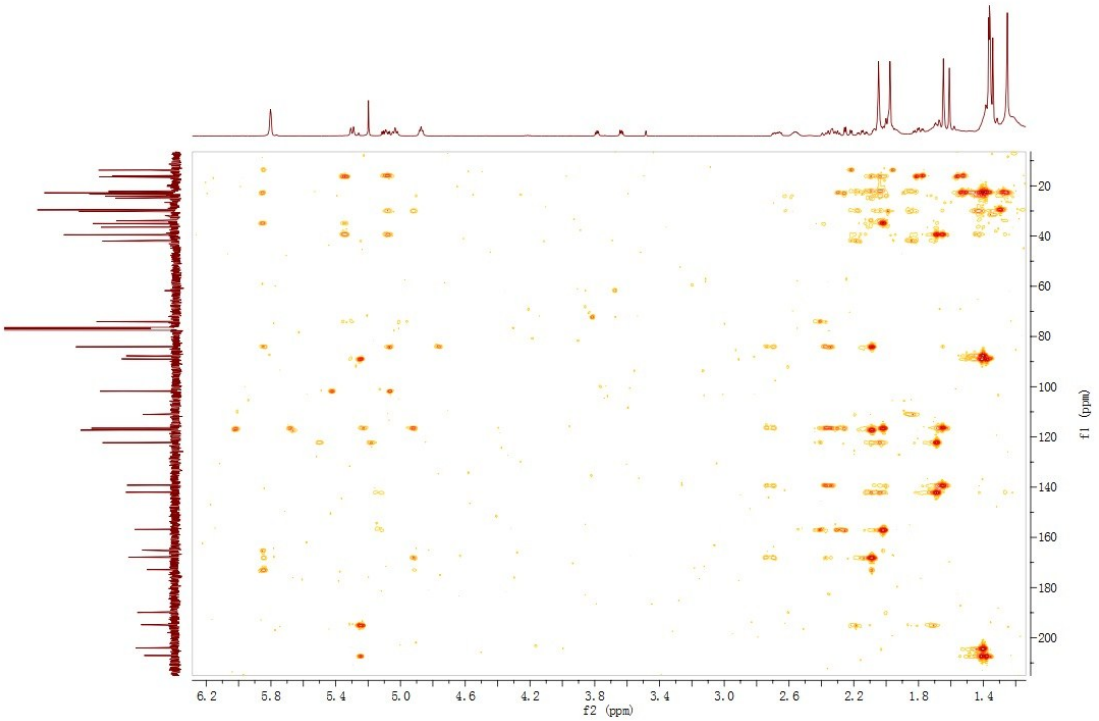
S66. ^{1}H NMR spectrum of compound aphanamene O (**13**) in $CDCl_3$



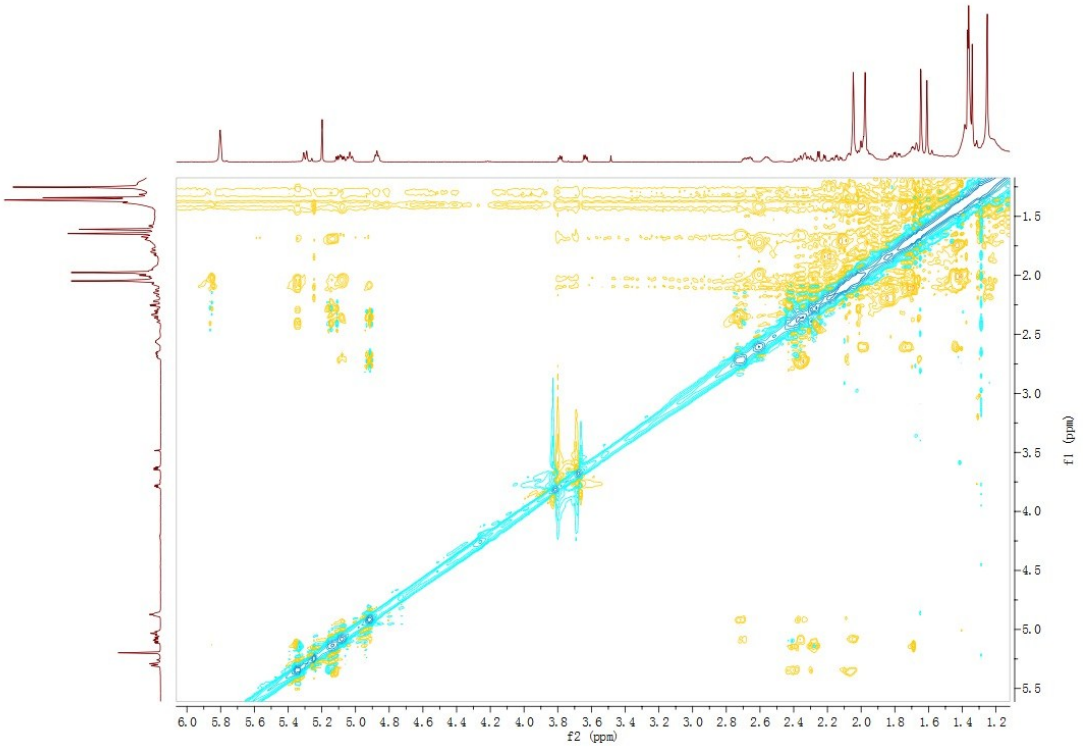
S67. ^{13}C NMR spectrum of compound aphanamene O (**13**) in $CDCl_3$



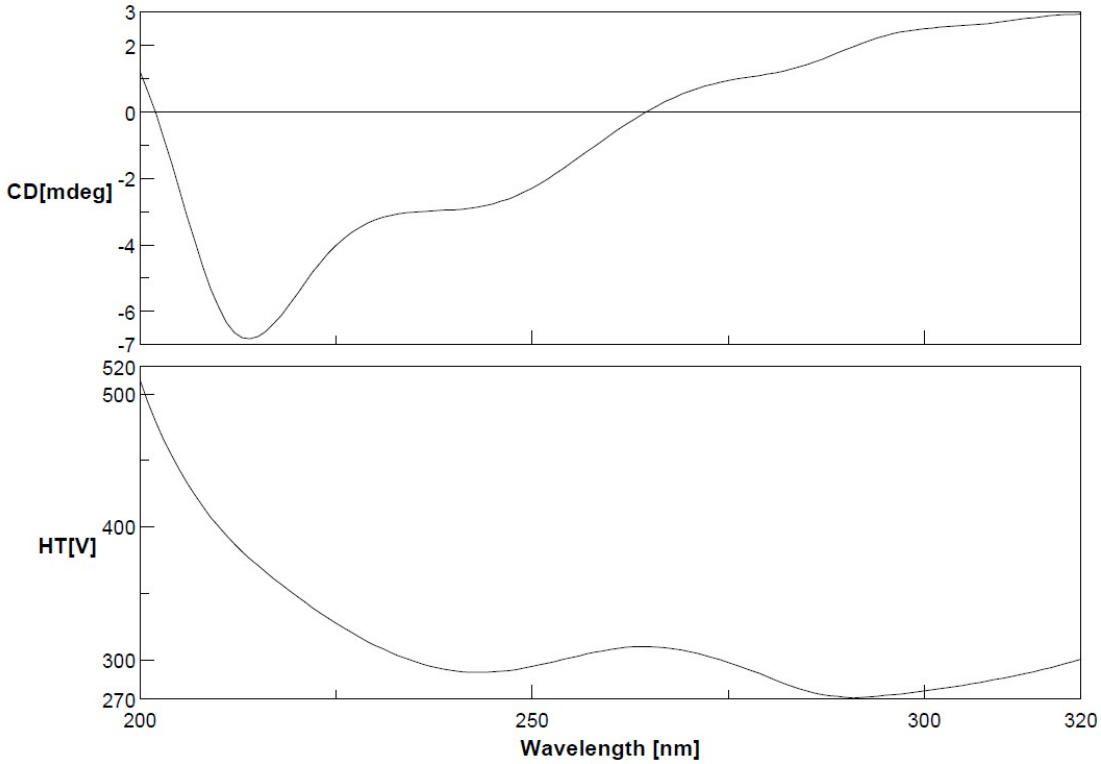
S68. HSQC spectrum of compound aphanamene O (**13**) in $CDCl_3$



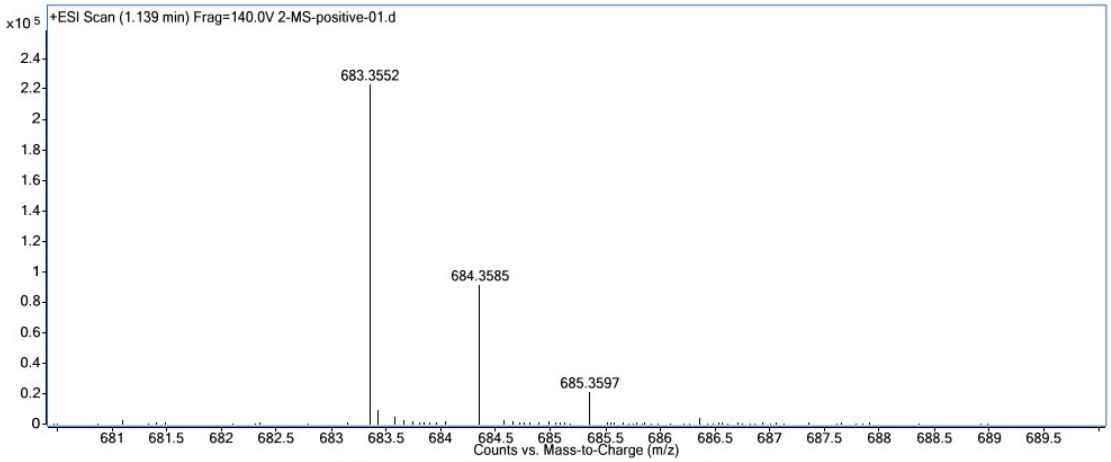
S69. HMBC spectrum of compound aphanamene O (**13**) in $CDCl_3$



S70. ROESY spectrum of compound aphanamene O (**13**) in $CDCl_3$



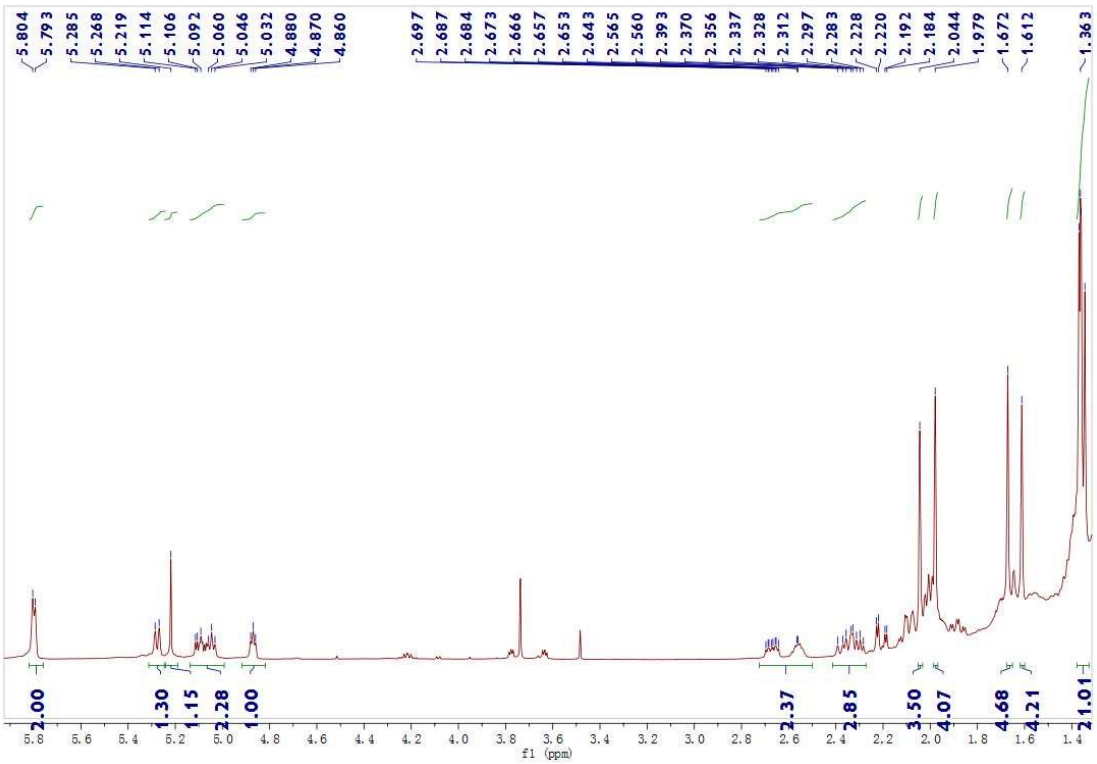
S71. CD spectrum of compound aphanamene O (**13**) in CH_3OH



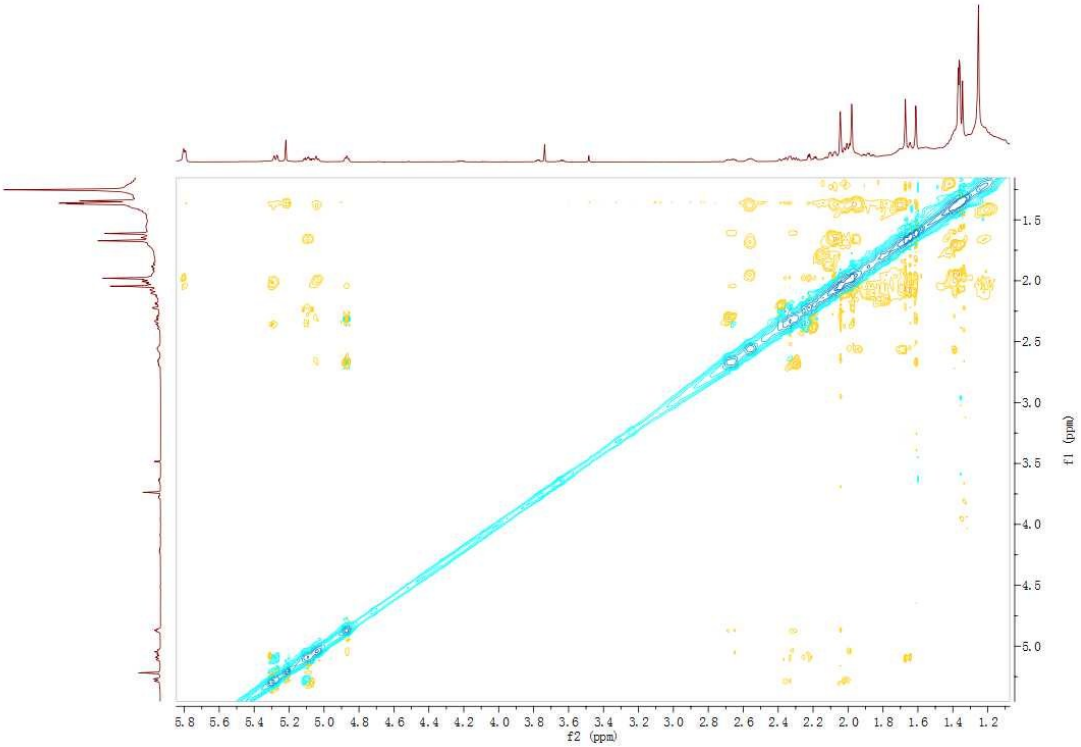
Elemental Composition Calculator

Target m/z:	683.3552	**Result type:**	Positive ions	**Species:**	$[M+Na]^+$
Elements:		C (0-80); H (0-120); O (0-30); N (0-10); Na (0-5); Cl (0-5)			
Ion Formula		**Calculated m/z**		**PPM Error**	
C40H52NaO8		683.3554		0.41	

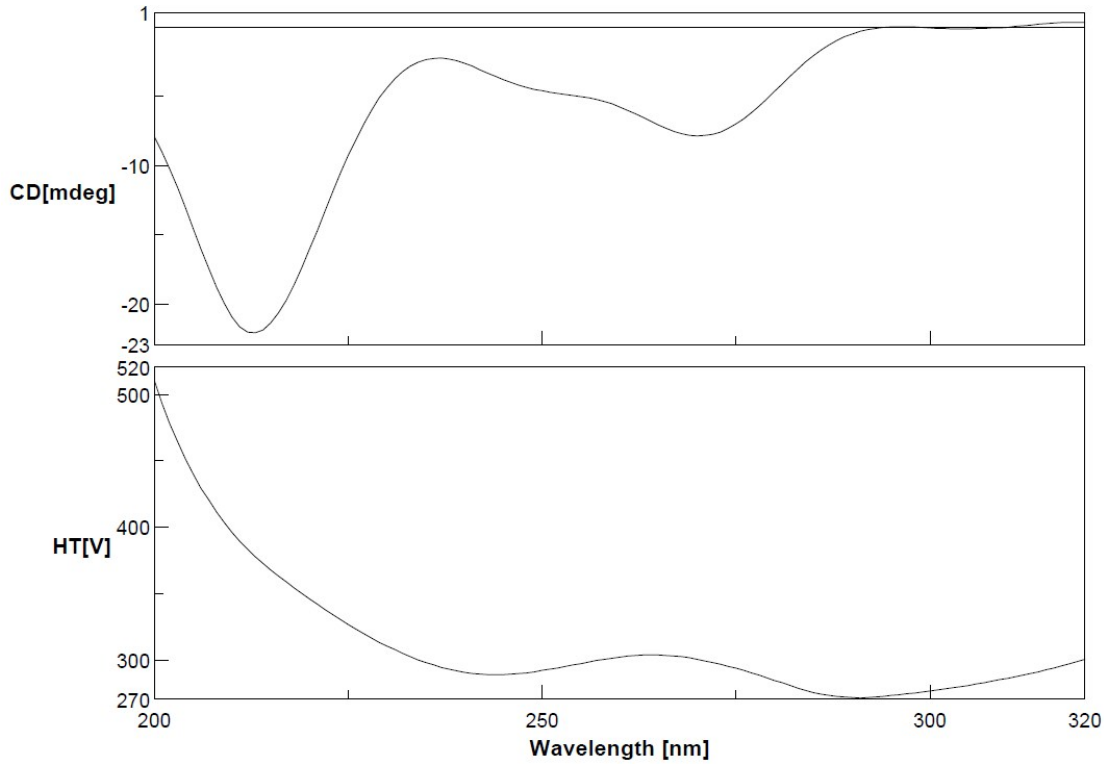
S72. HRESIMS spectrum of compound aphanamene O (**13**) in CH_3OH



S73. ^{1}H NMR spectrum of compound aphanamene P (**14**) in $CDCl_3$



S74. ROESY spectrum of compound aphanamene P (**14**) in $CDCl_3$



S75. CD spectrum of compound aphanamene P (**14**) in CH_3OH