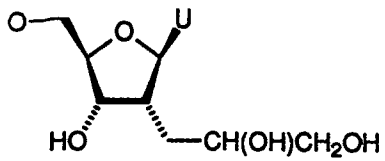
JOC The Journal of Organic Chemistry

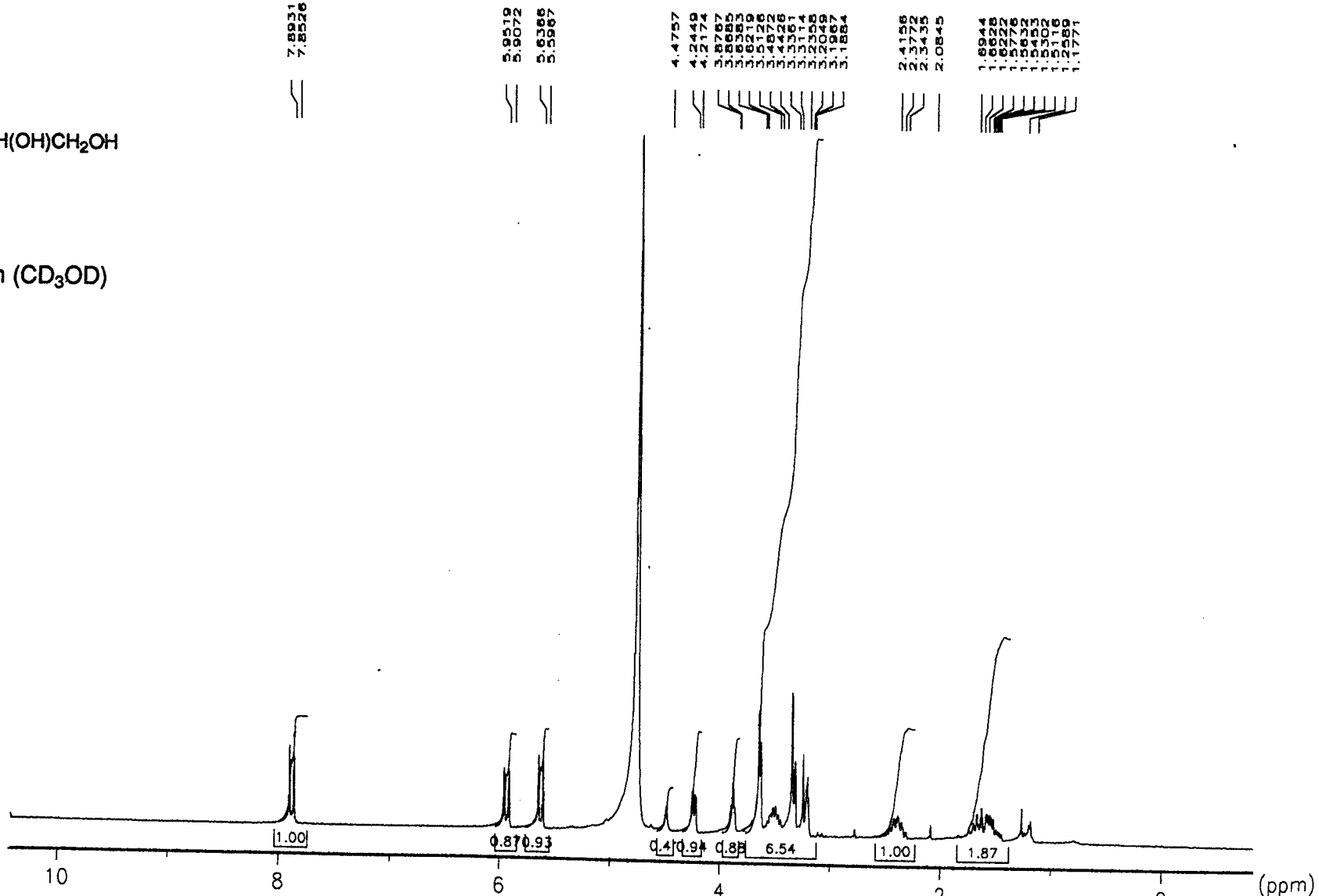
J. Org. Chem., 1996, 61(26), 9213-9222, DOI:10.1021/jo9614607

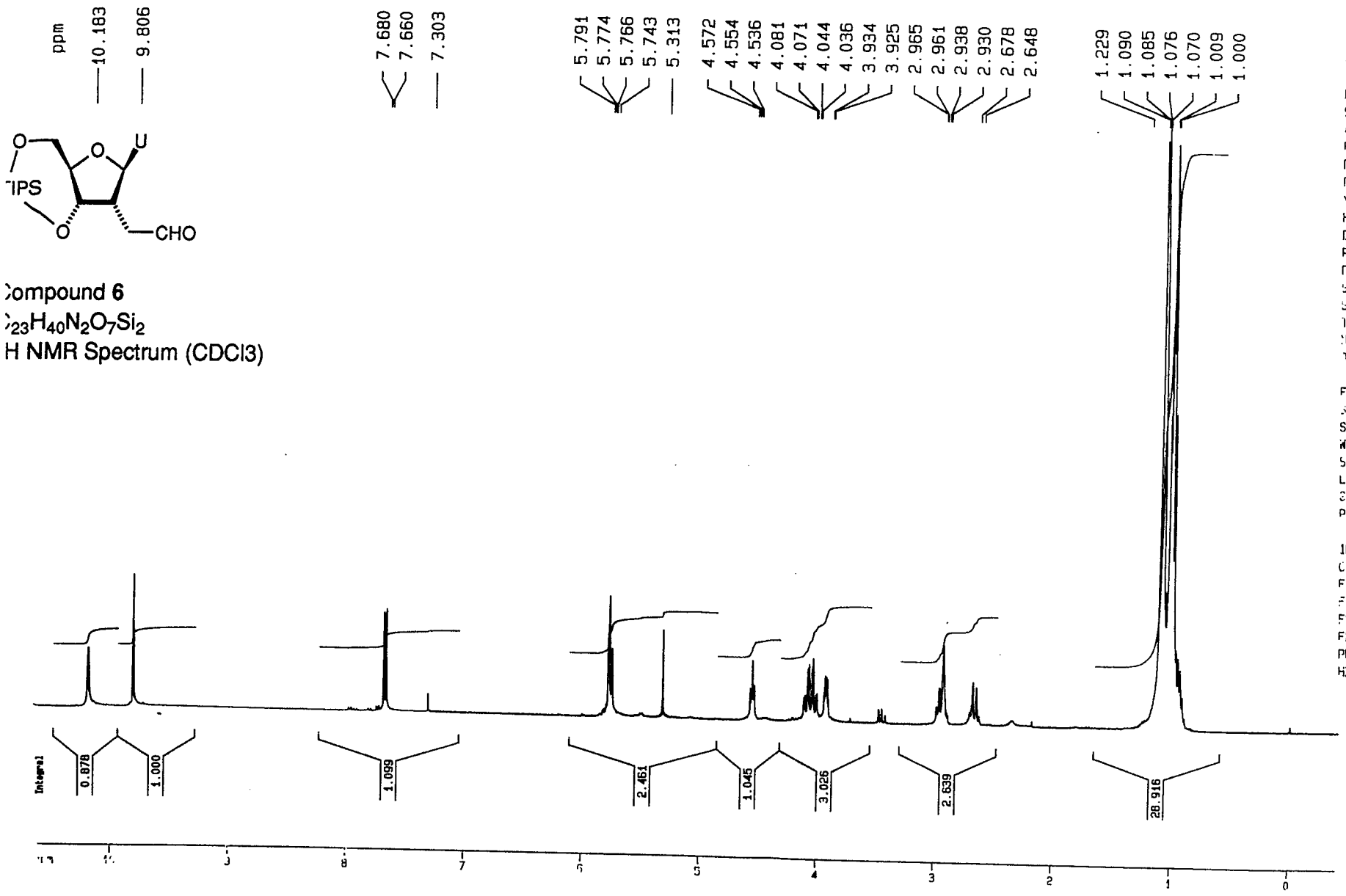


ompound **4**

$_{12}H_{18}N_2O_7$

1 NMR Spectrum (CD_3OD)

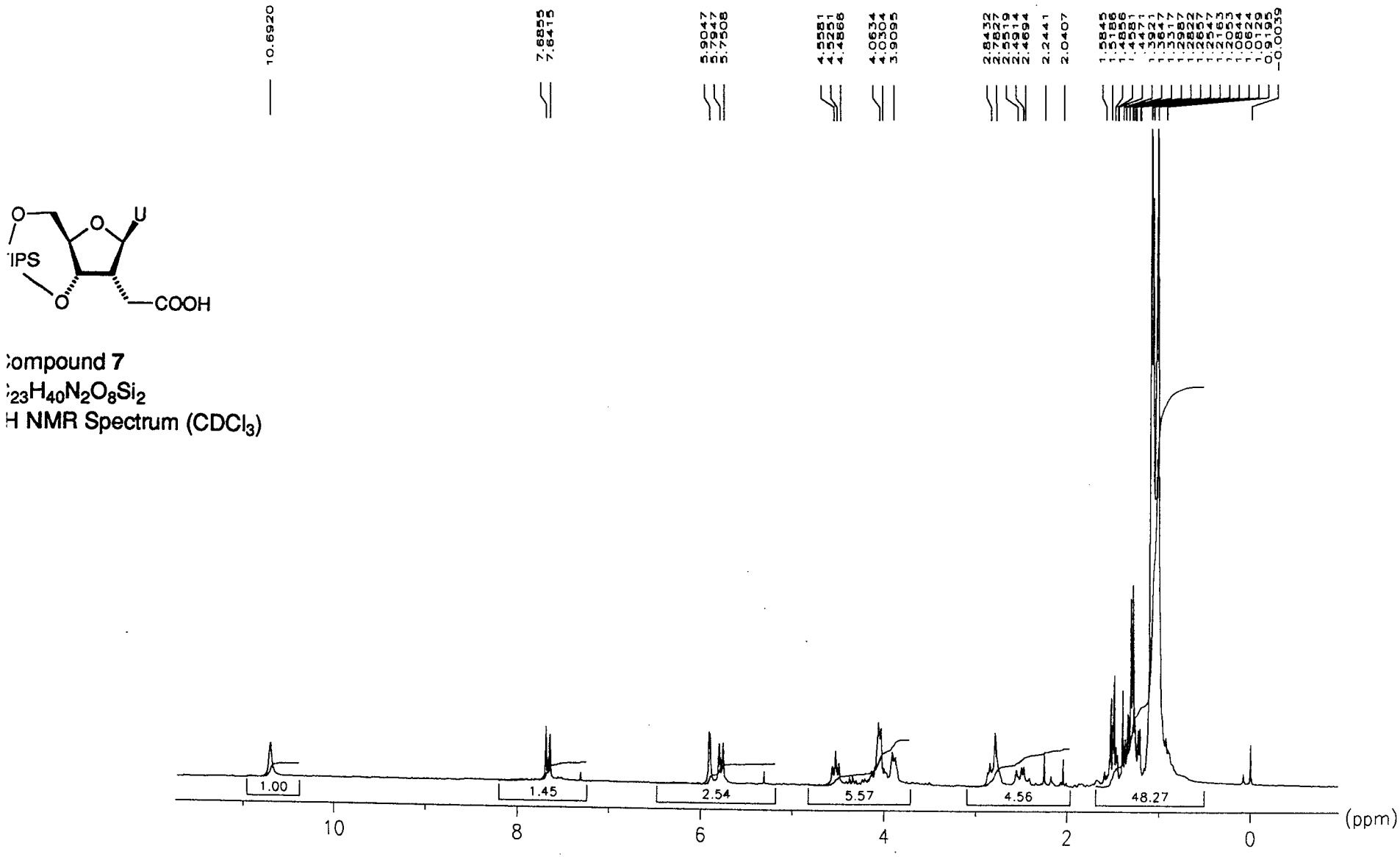




Compound 6

$C_{23}H_{40}N_2O_7Si_2$

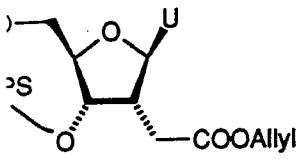
H NMR Spectrum (CDCl3)



Compound **7**

$C_{23}H_{40}N_2O_8Si_2$

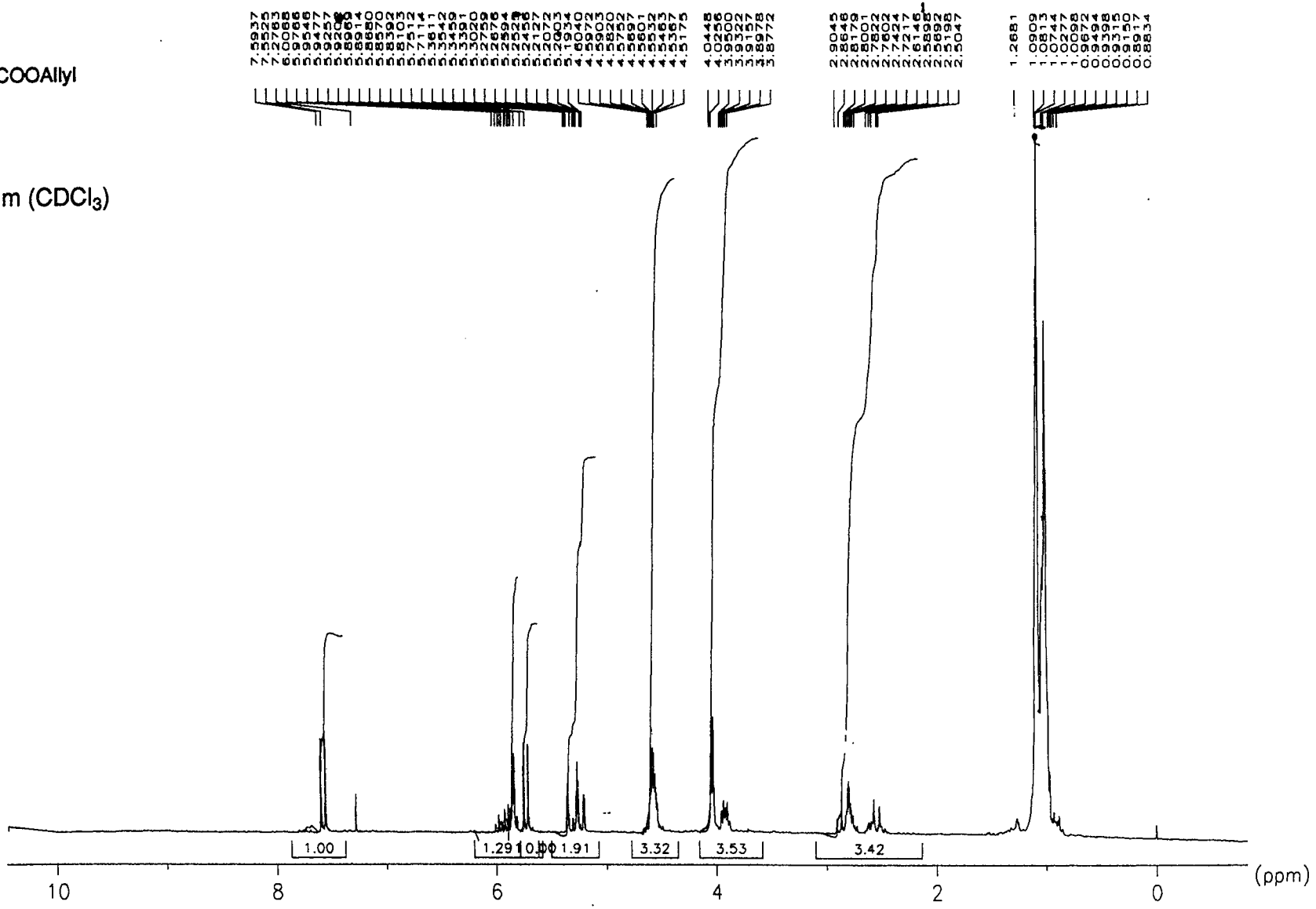
H NMR Spectrum ($CDCl_3$)

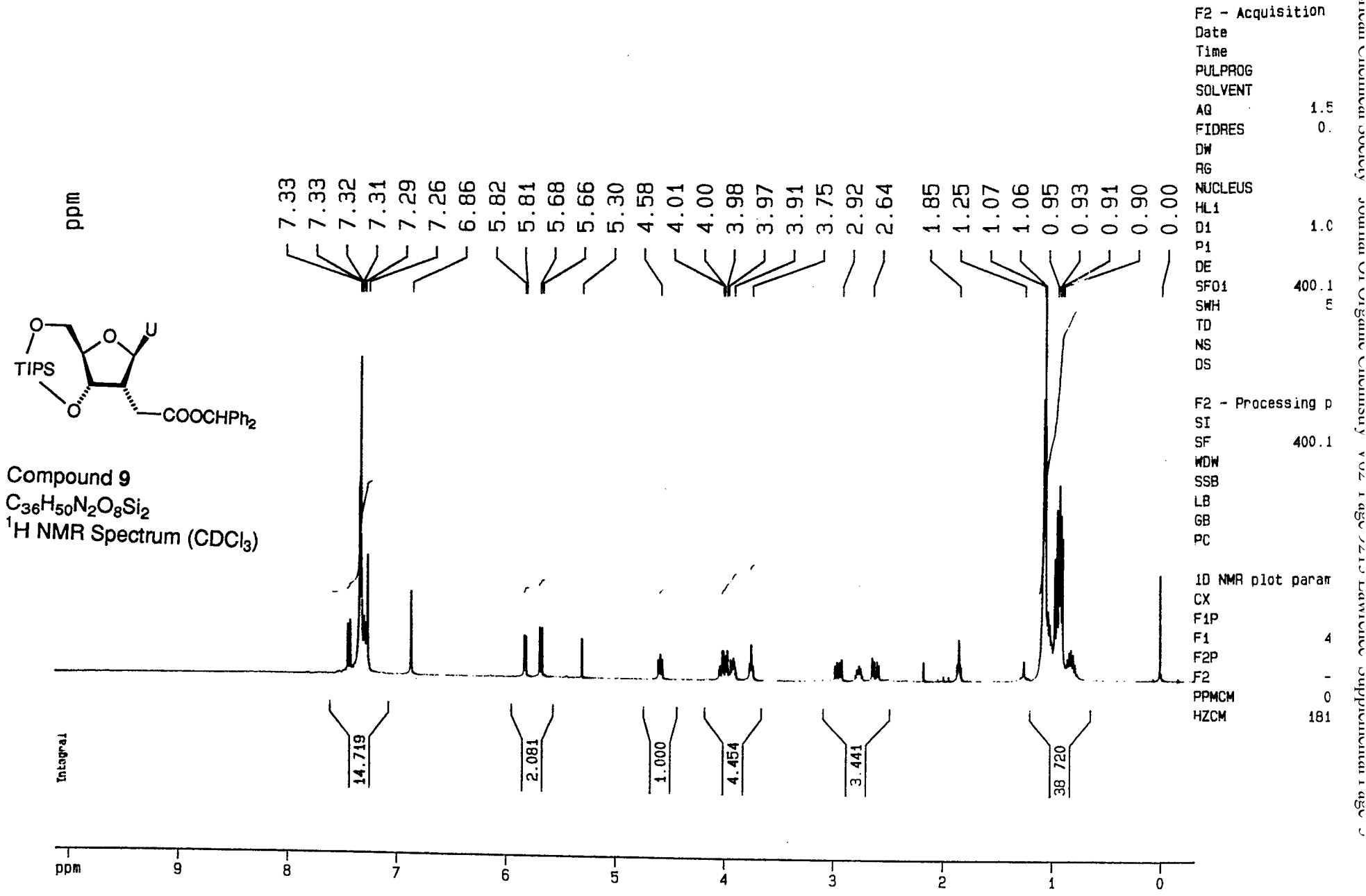


mpound **8**

$_{6}H_{45}N_2O_8Si_2$

NMR Spectrum ($CDCl_3$)

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7.33
7.32
7.31
7.29
7.26
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5.68
5.66
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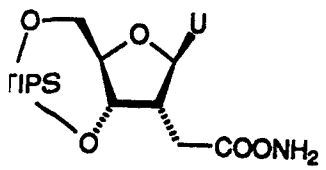
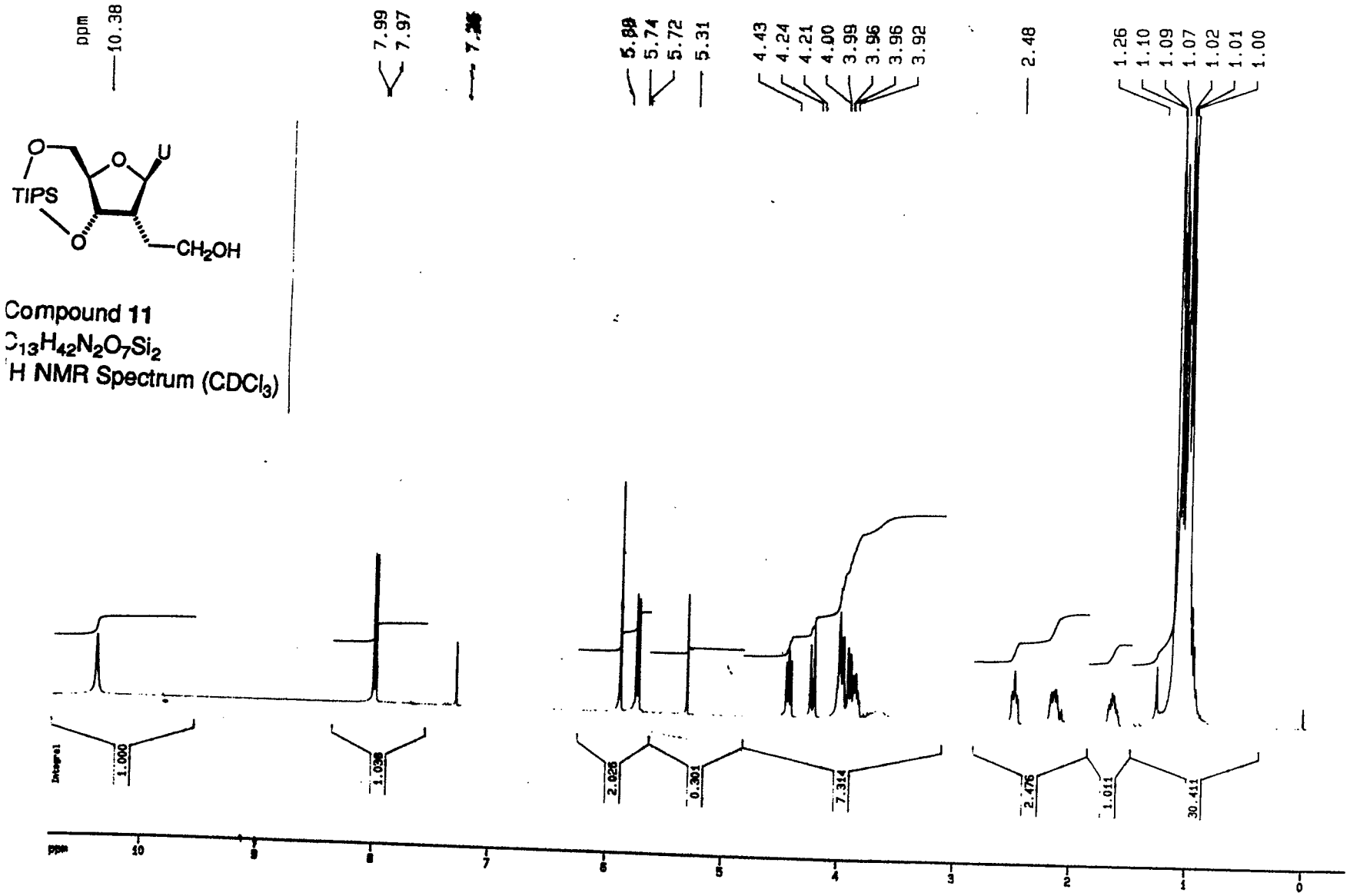
4.00
3.98
3.97
3.91
3.75
2.92
2.64
1.85
1.25
1.07
1.06
0.95
0.93
0.91
0.90
0.00
TIPS
COOCHPh2
Compound 9
C36H50N2O8Si2
1H NMR Spectrum (CDCl3)
Integral
14.719
2.081
1.000
4.454
3.441
38.720
ppm
9
8
7
6
5
4
3
2
1
0
F2 - Acquisition
Date
Time
PULPROG
SOLVENT
AQ 1.5
FIDRES 0.
DW
RG
NUCLEUS
HL1
D1 1.0
P1
DE
SFO1 400.1
SWH 5
TD
NS
DS
F2 - Processing p
SI
SF 400.1
WDW
SSB
LB
GB
PC
1D NMR plot param
CX
F1P
F1 4
F2P
F2 -
PPMCM 0
HZCM 181



Compound 10
$C_{23}H_{41}N_3O_7Si_2$
^{1}H NMR Spectrum ($CDCl_3$)



Compound **11**

$C_{13}H_{42}N_2O_7Si_2$

^1H NMR Spectrum ($CDCl_3$)

Current Data Parame
NAME A.LAMF
EXPNO
PROCNO

F2 - Acquisition Pa
Date 94
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SOLVENT C
AQ 1.499
FIDRES 0.33
DW
RG
NUCLEUS
HL1
D1 1.000
P1
DE 1
SFO1 400.136
SWH 526.
TD 1
NS
DS

F2 - Processing para
SI 1
SF 400.134
NDW
SSB
LB 0
GB
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1D NMR plot parame
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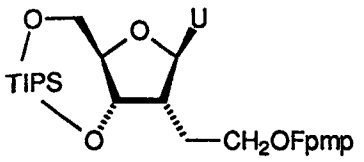
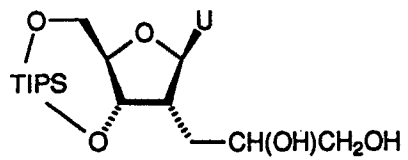


Compound **12**

$C_{35}H_{56}N_3O_8FSi_2$

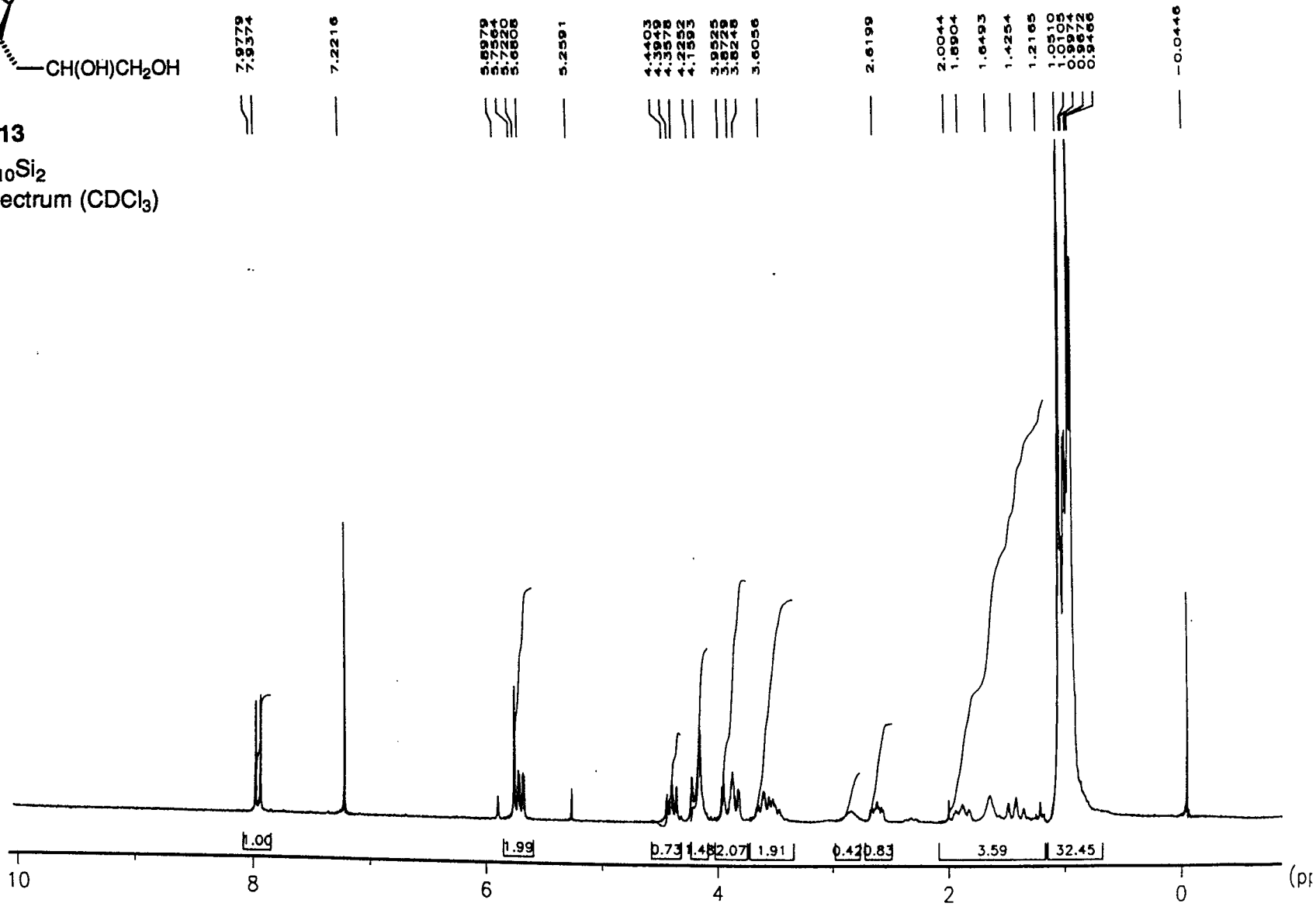
^1H NMR Spectrum ($CDCl_3$)

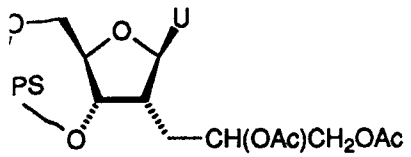


Compound **13**

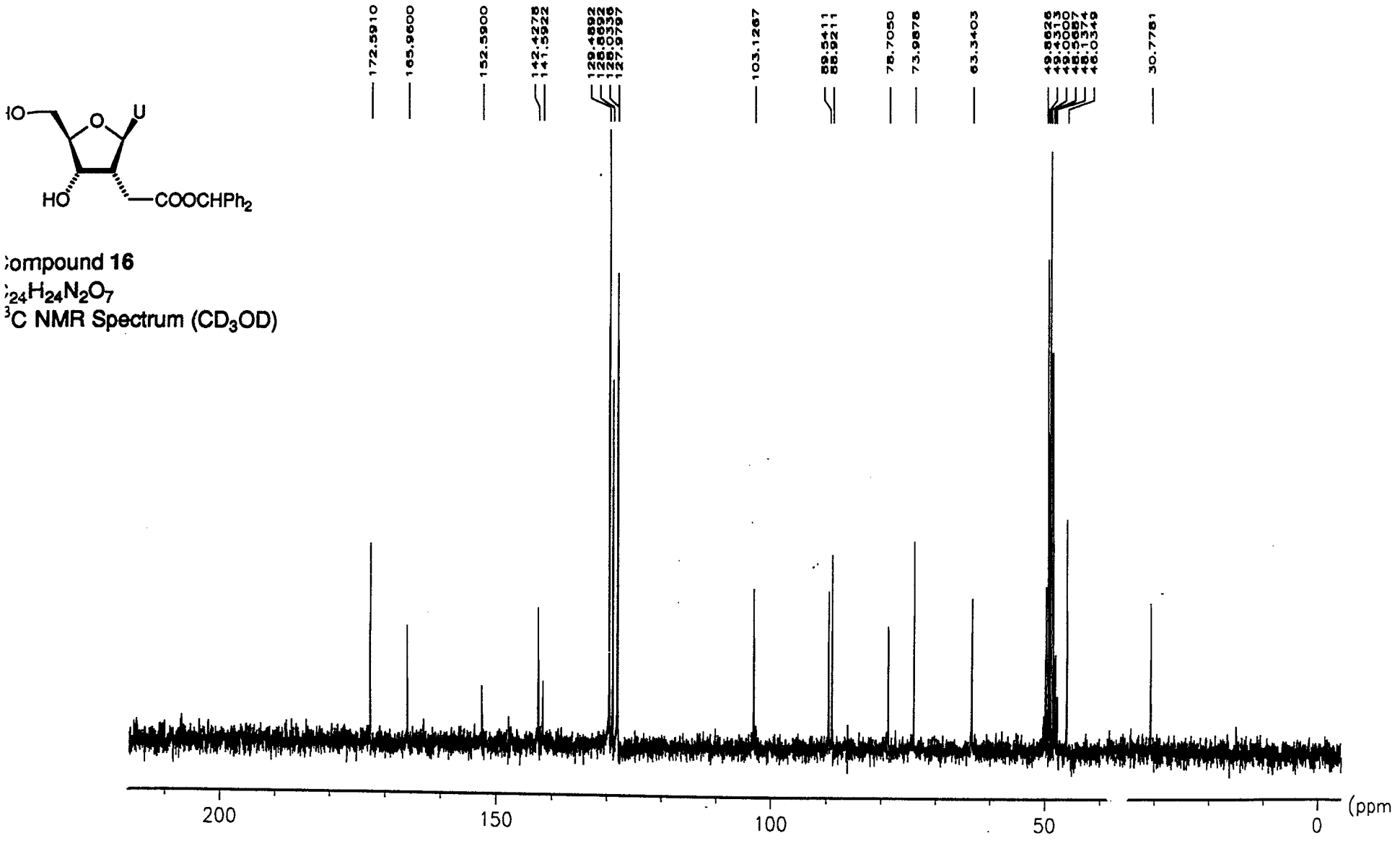
$C_{28}H_{48}N_2O_{10}Si_2$

^{1}H NMR Spectrum ($CDCl_3$)





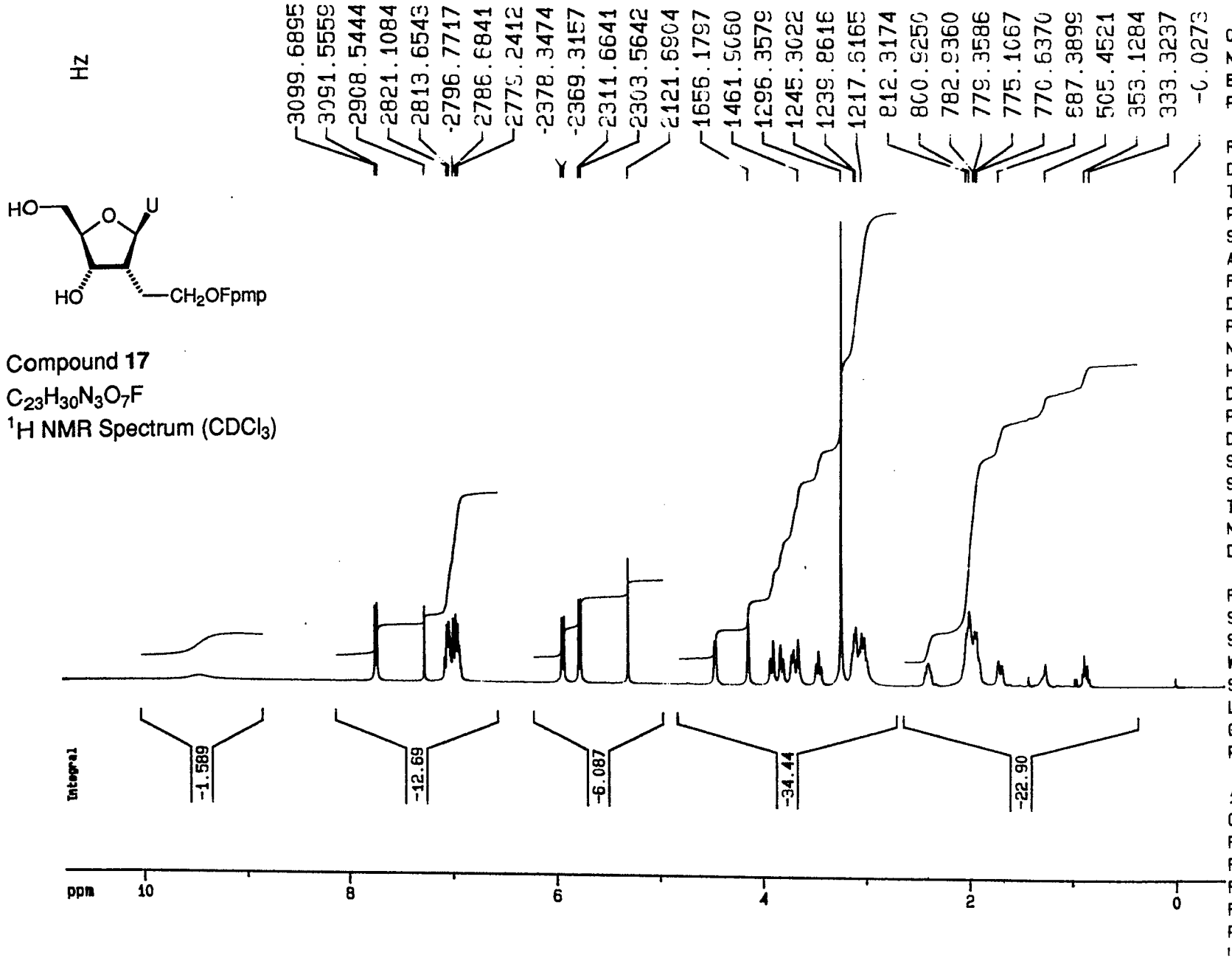
ompound **14**
$_{28}H_{48}N_2O_{10}Si_2$
H NMR Spectrum ($CDCl_3$)

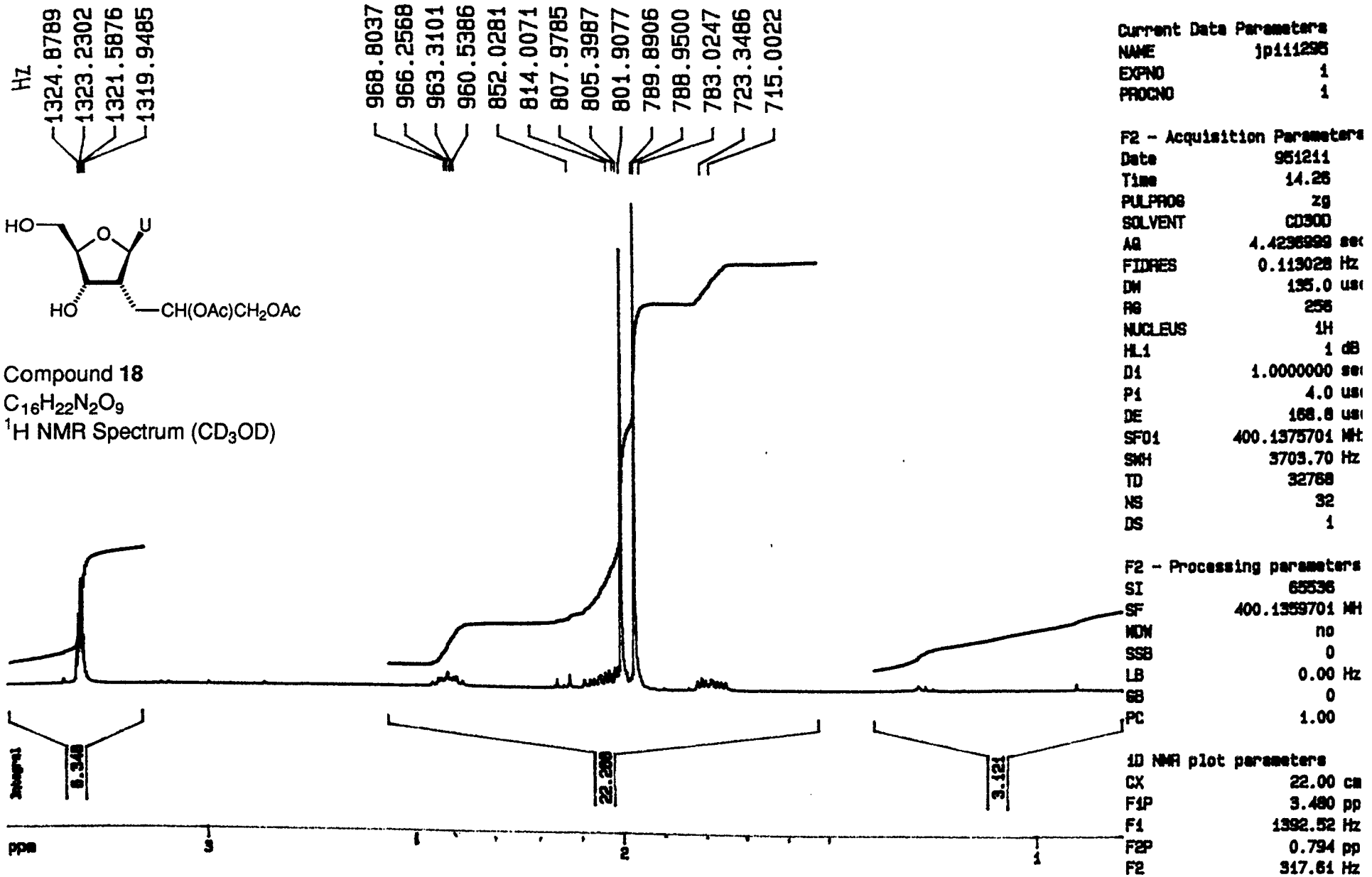


ompound **16**

$C_{24}H_{24}N_2O_7$

^{13}C NMR Spectrum (CD_3OD)





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EXPNO	1
PROCNO	1

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NUCLEUS	1H
HL1	1 dB
D1	1.0000000 sec
P1	4.0 usec
DE	168.8 usec
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SWH	3703.70 Hz
TD	32768
NS	32
DS	1

F2 - Processing parameters

SI	65536
SF	400.1359701 MHz
WDW	no
SSB	0
LB	0.00 Hz
GB	0
PC	1.00

1D NMR plot parameters

CX	22.00 cm
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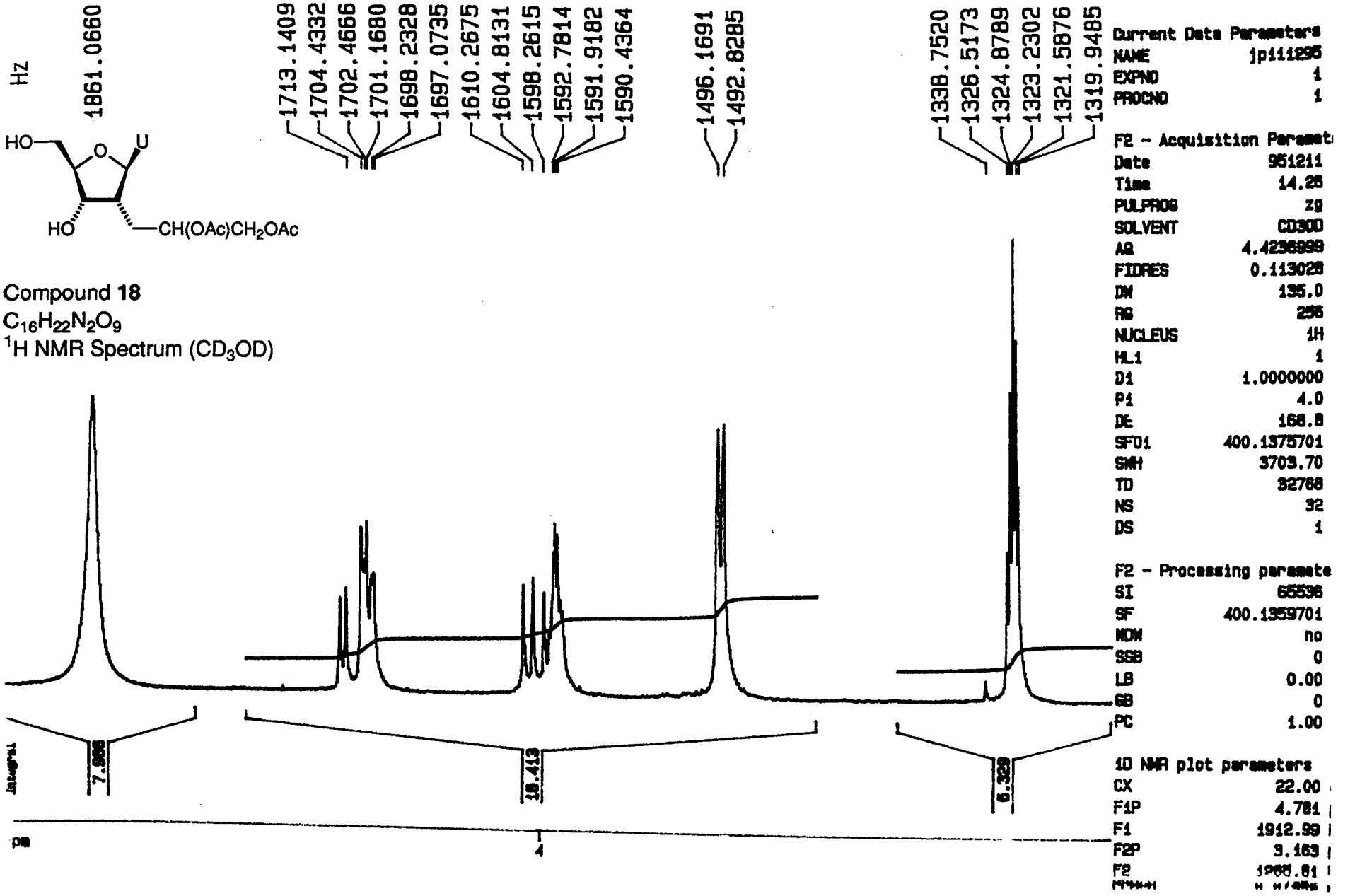
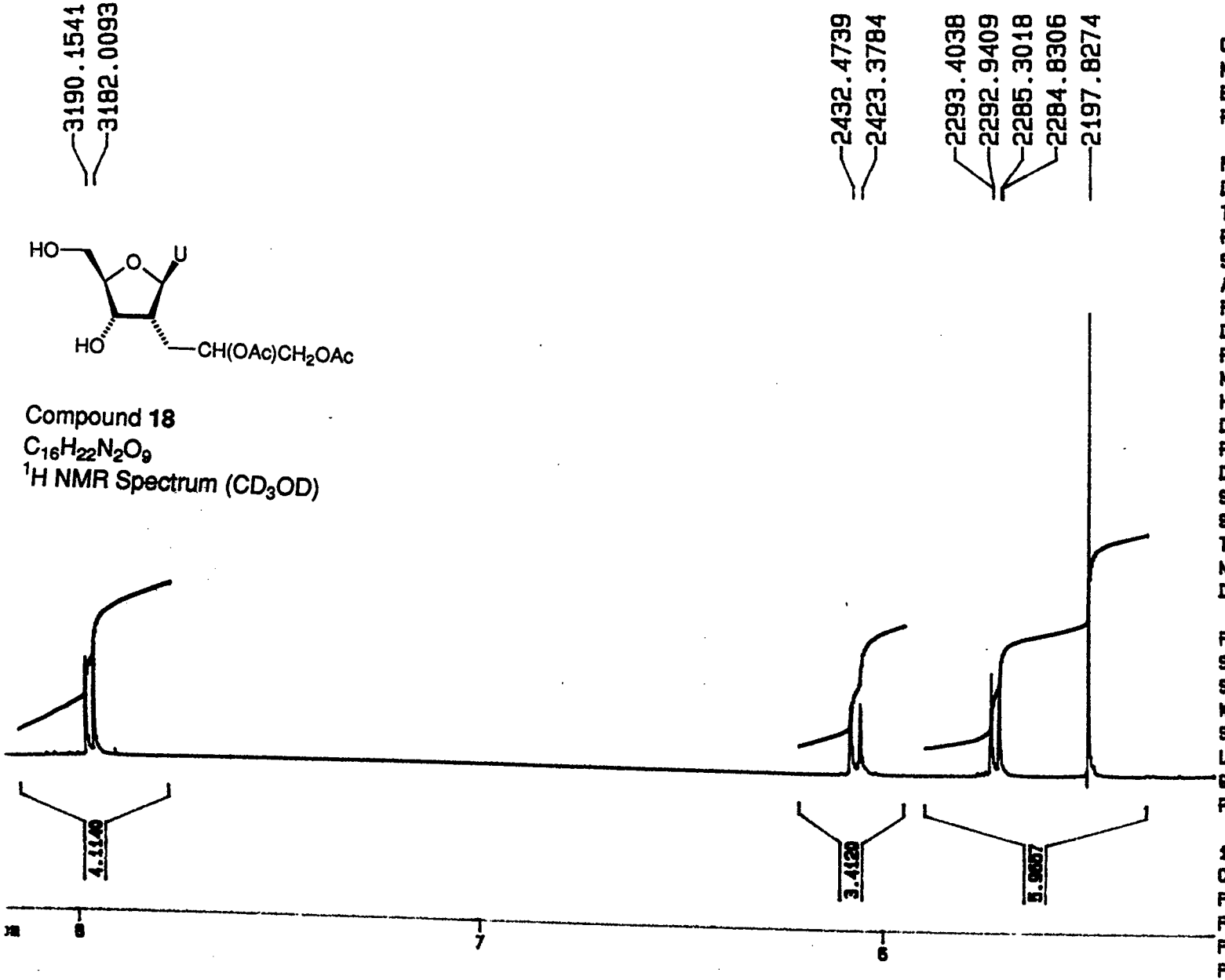
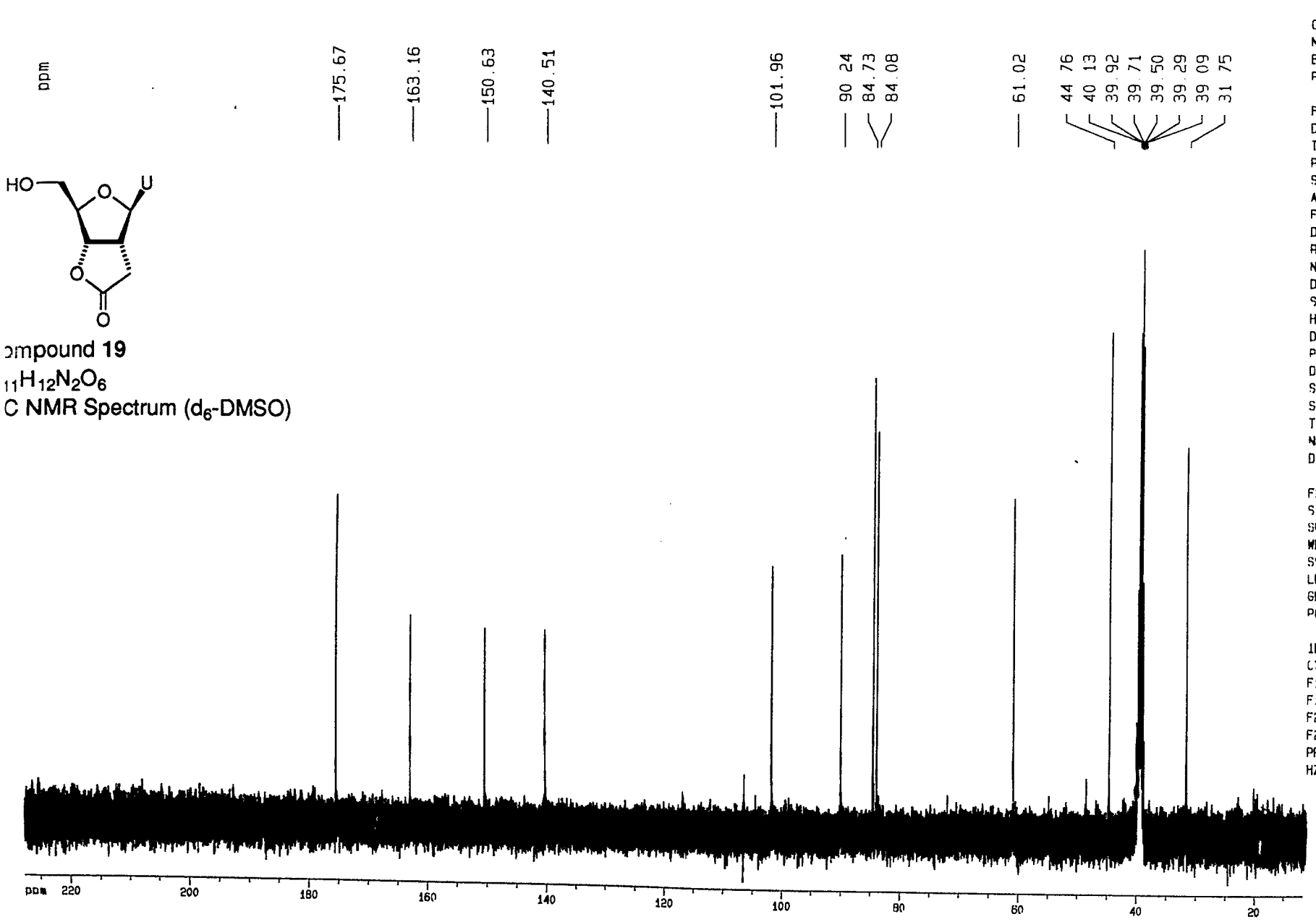
F2P	0.794 ppm
F2	317.61 Hz
PPMCM	0.12211 ppm/cm
HZCM	48.86941 Hz/cm



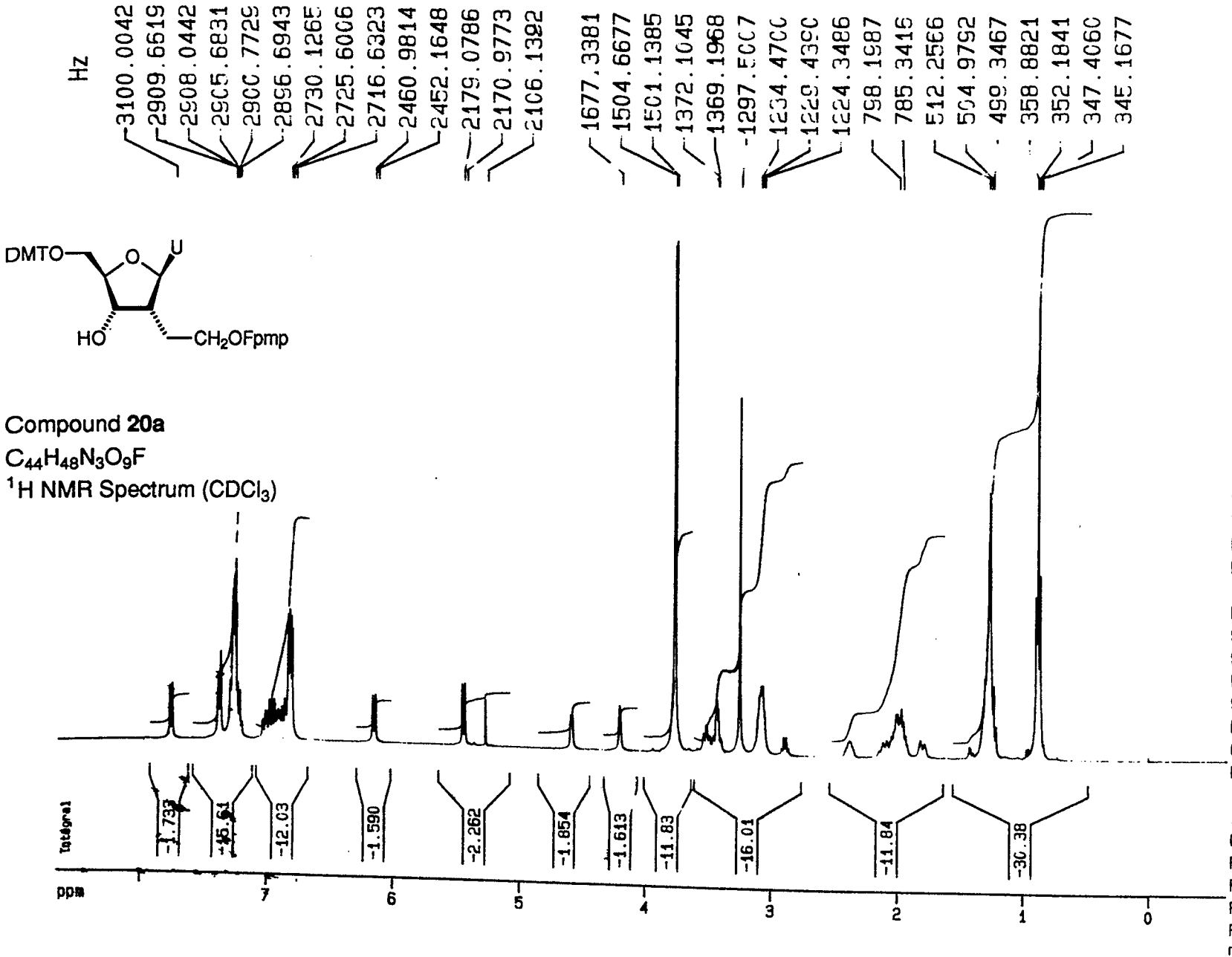


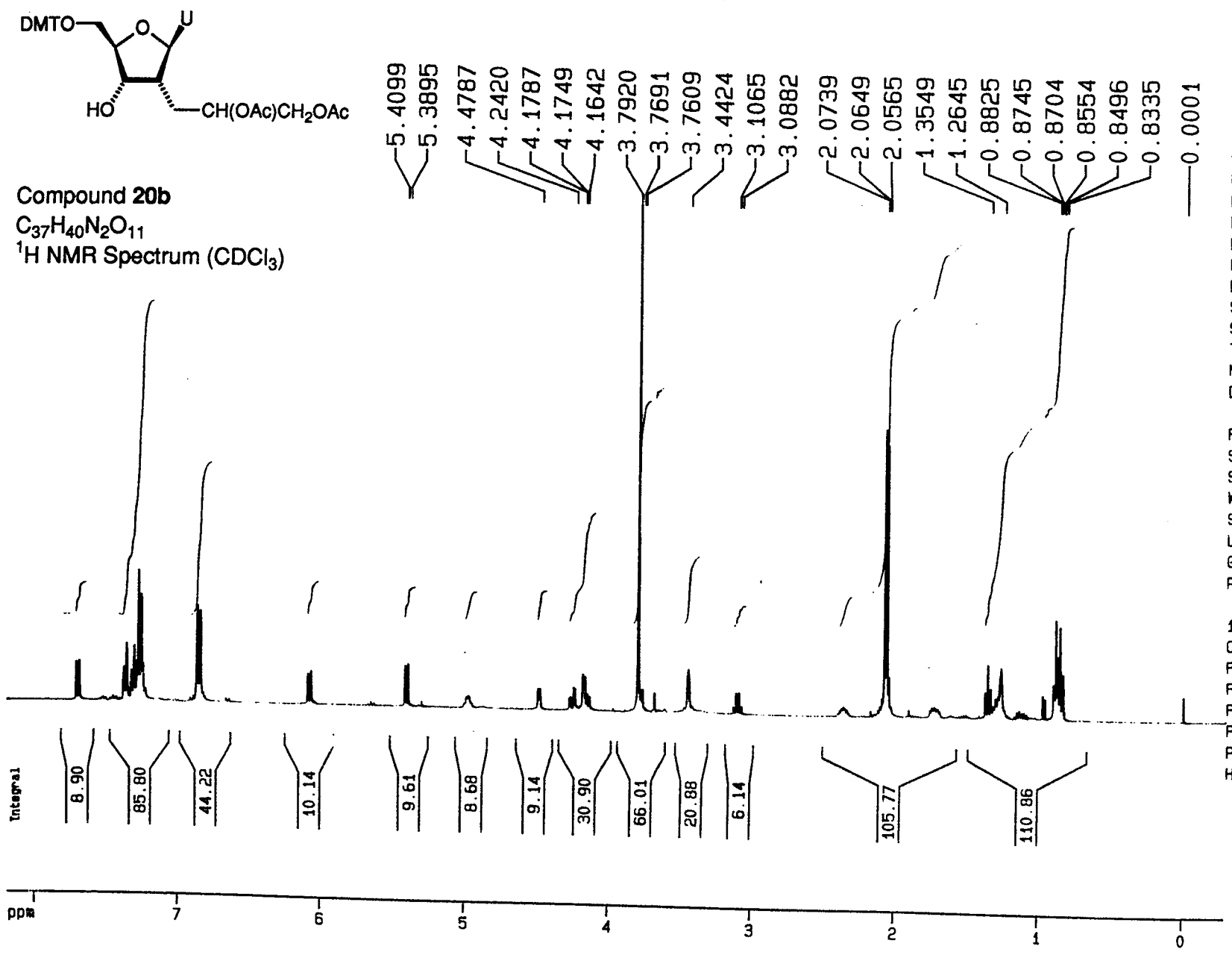


ompound **19**

$_{11}H_{12}N_2O_6$

C NMR Spectrum (d_6-DMSO)




DMTO
O
U
HO
CH(OAc)CH2OAc
Compound 20b
C37H40N2O11
1H NMR Spectrum (CDCl3)
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5.3895
4.4787
4.2420
4.1787
4.1749
4.1642
3.7920
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3.7609
3.4424
3.1065
3.0882
2.0739
2.0649
2.0565
1.3549
1.2645
0.8825
0.8745
0.8704
0.8554
0.8496
0.8335
0.0001
Integral
8.90
85.80
44.22
10.14
9.61
8.68
9.14
30.90
66.01
20.88
6.14
105.77
110.86
ppm
7
6
5
4
3
2
1
0
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Date 96(
Time 1
PULPROG
SOLVENT CI
AQ 3.112
FIDRES 0.16(
DW
RG
NUCLEUS
HL1
D1 1.000(
P1
DE 1
SFO1 400.136
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DS
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SF 400.134
WDW
SSB
LB
GB
PC
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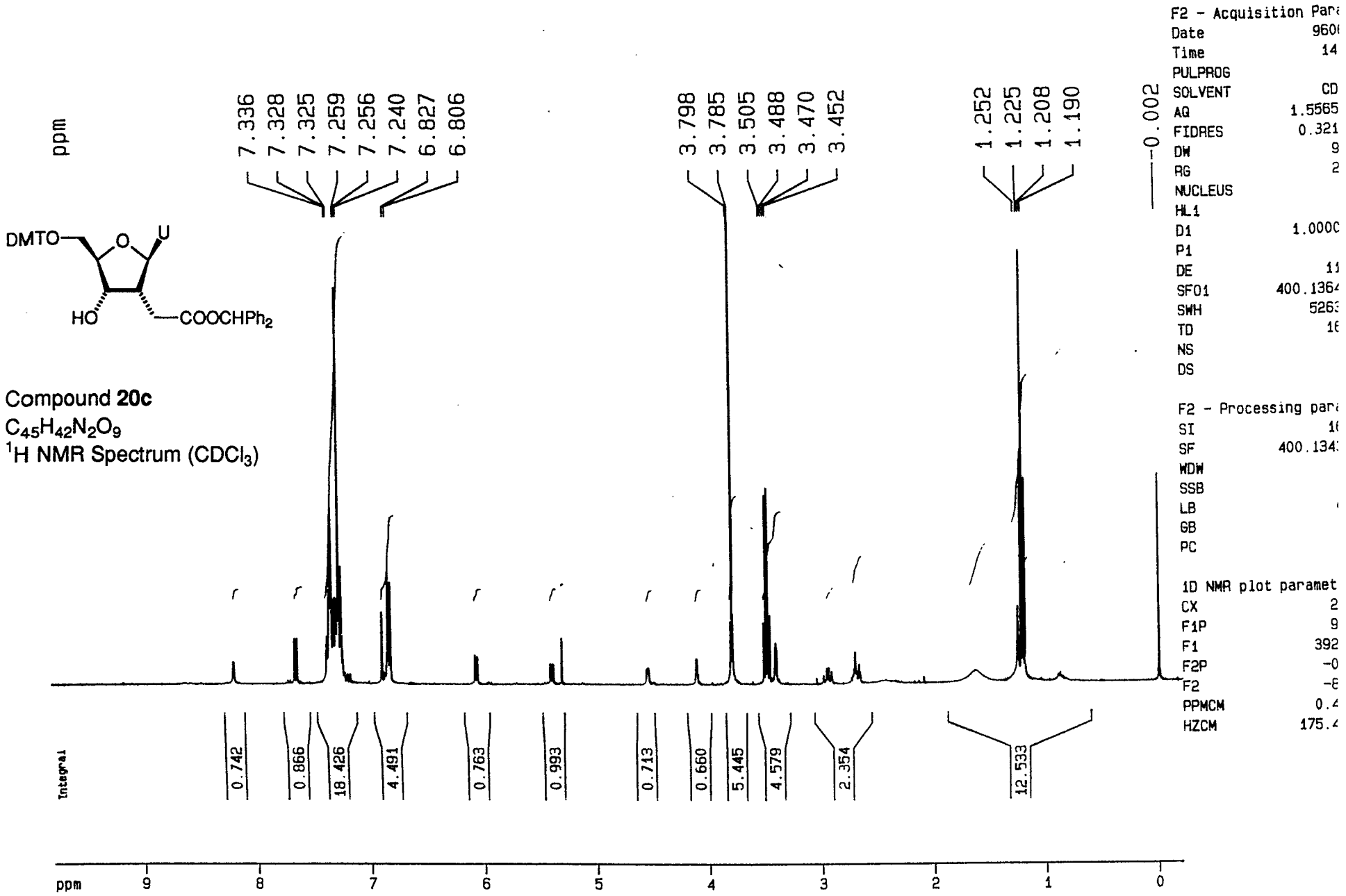

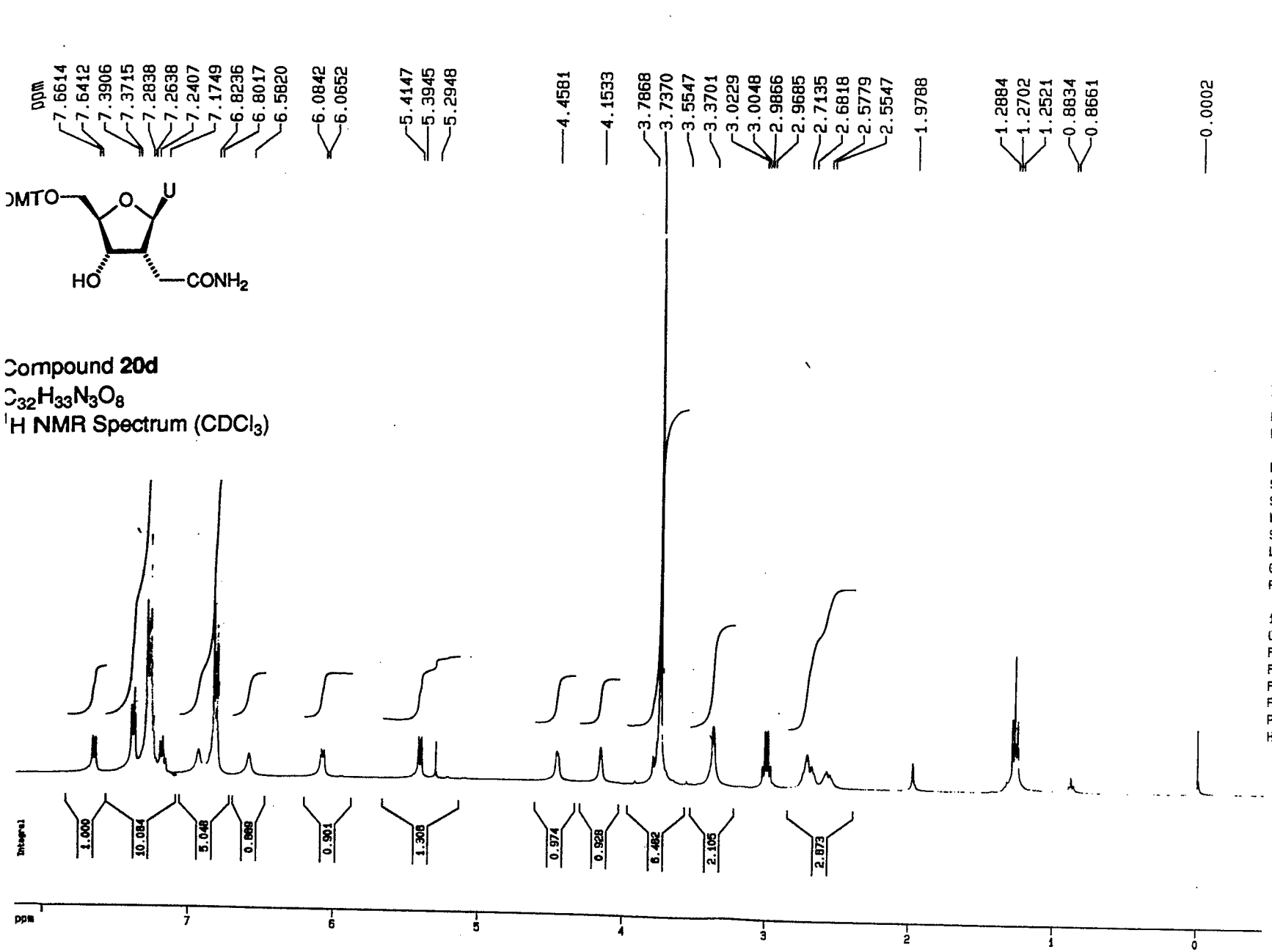


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EXPNO
PROCNO

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F2 - Processing param
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SSB
LB                   0.
GB
PC                   4.

1D NMR plot parameter
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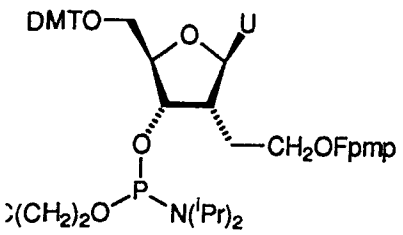
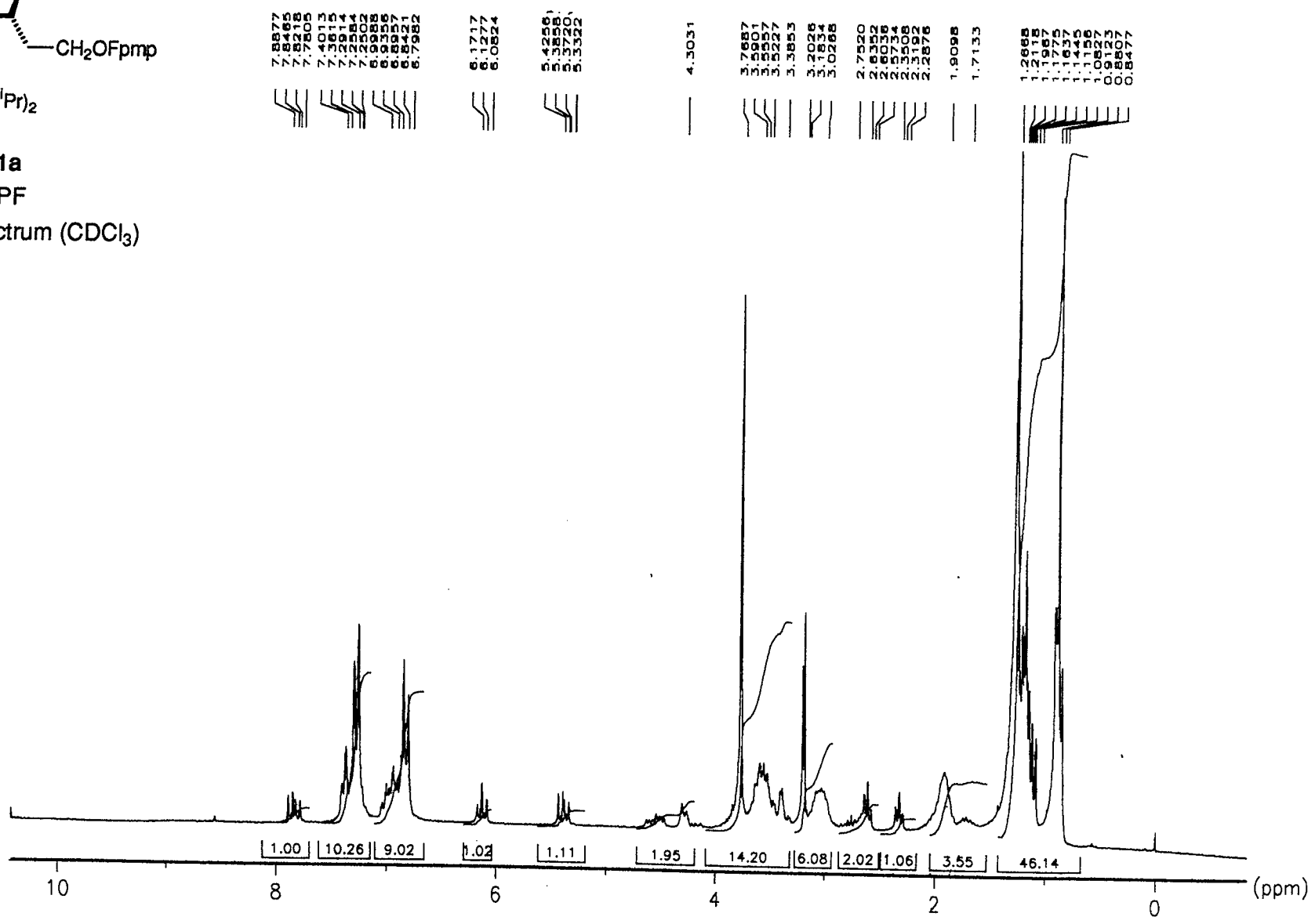
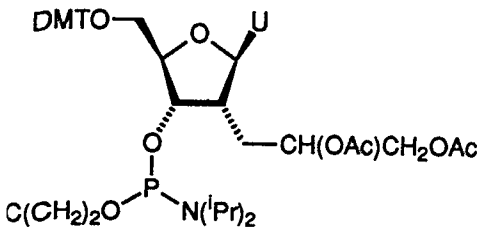


Compound **21a**

$C_{53}H_{65}N_5O_{10}PF$

^1H NMR Spectrum ($CDCl_3$)

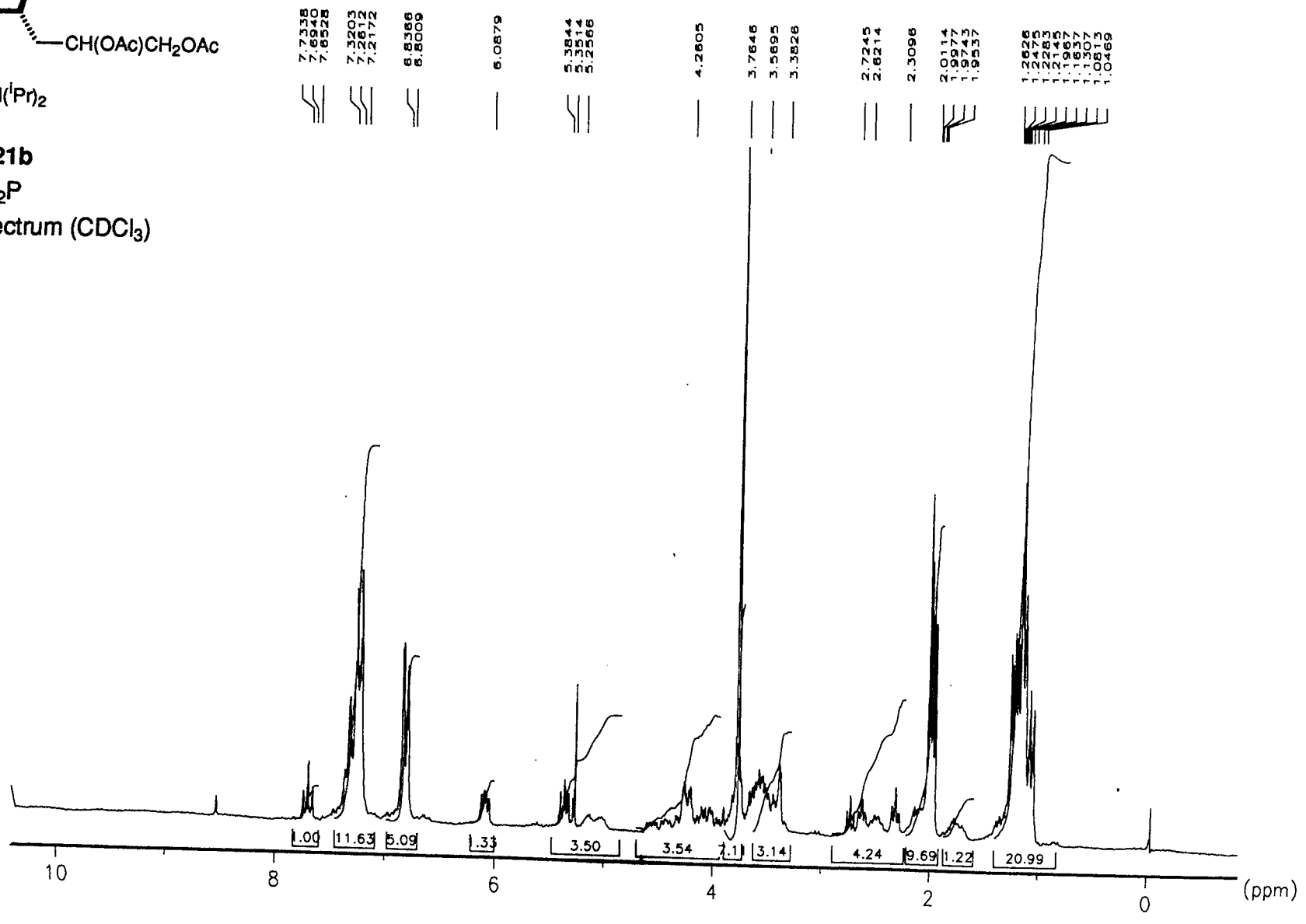


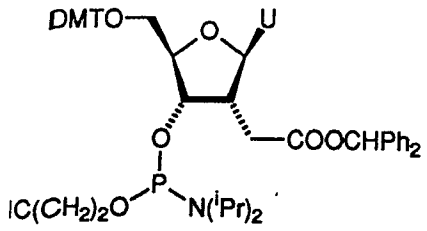


Compound 21b

$C_{46}H_{57}N_4O_{12}P$

^1H NMR Spectrum ($CDCl_3$)

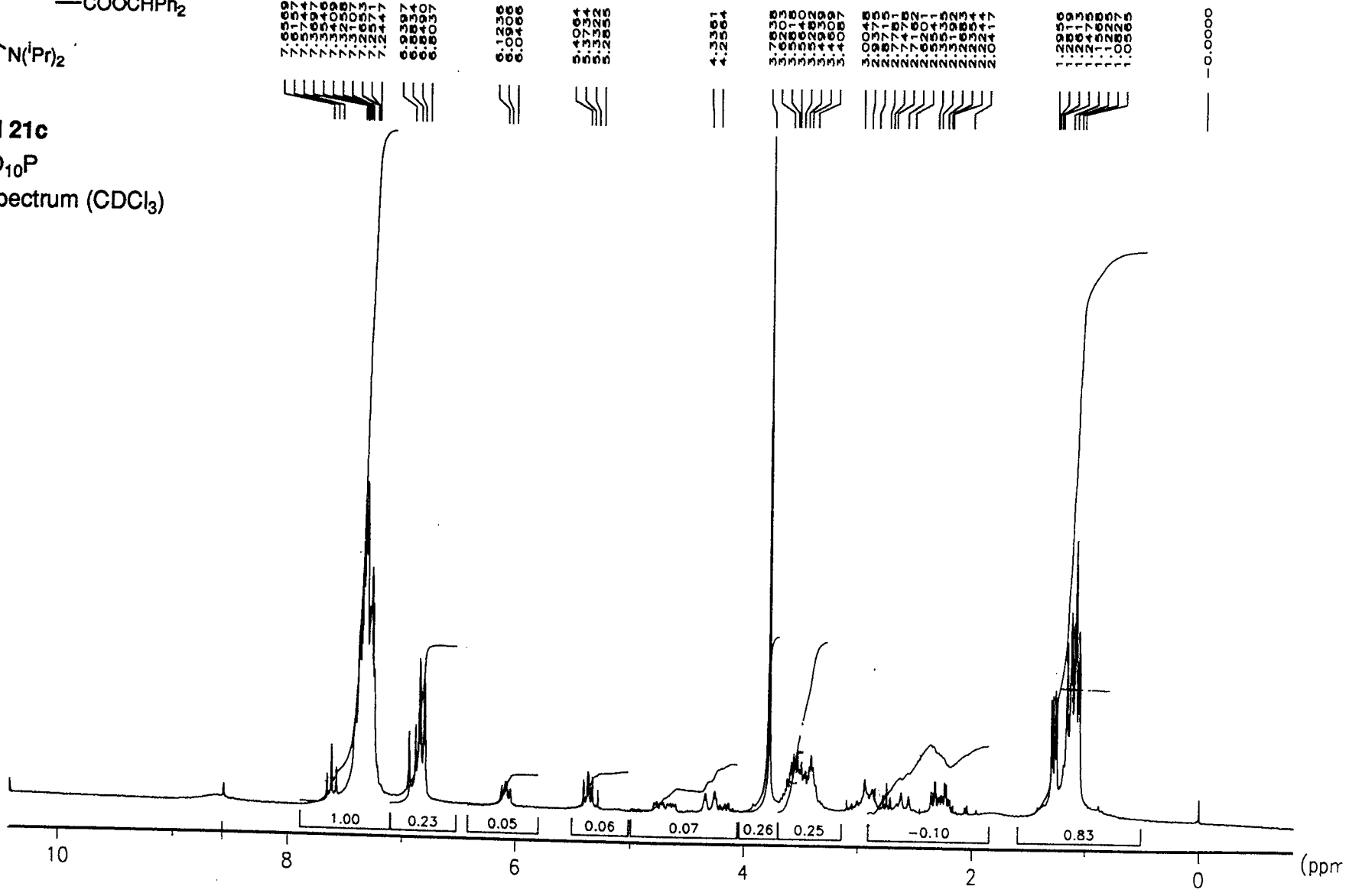


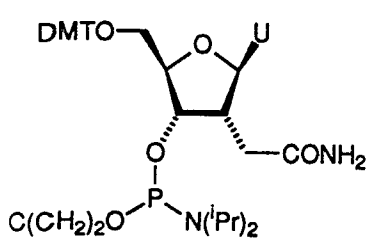


Compound **21c**

$C_{50}H_{60}N_4O_{10}P$

^1H NMR Spectrum ($CDCl_3$)

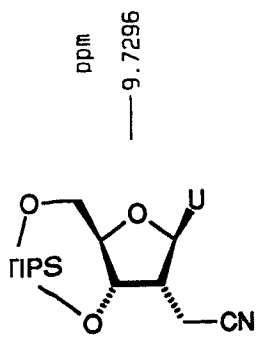




Compound 21d

$C_{41}H_{50}N_5O_9P$

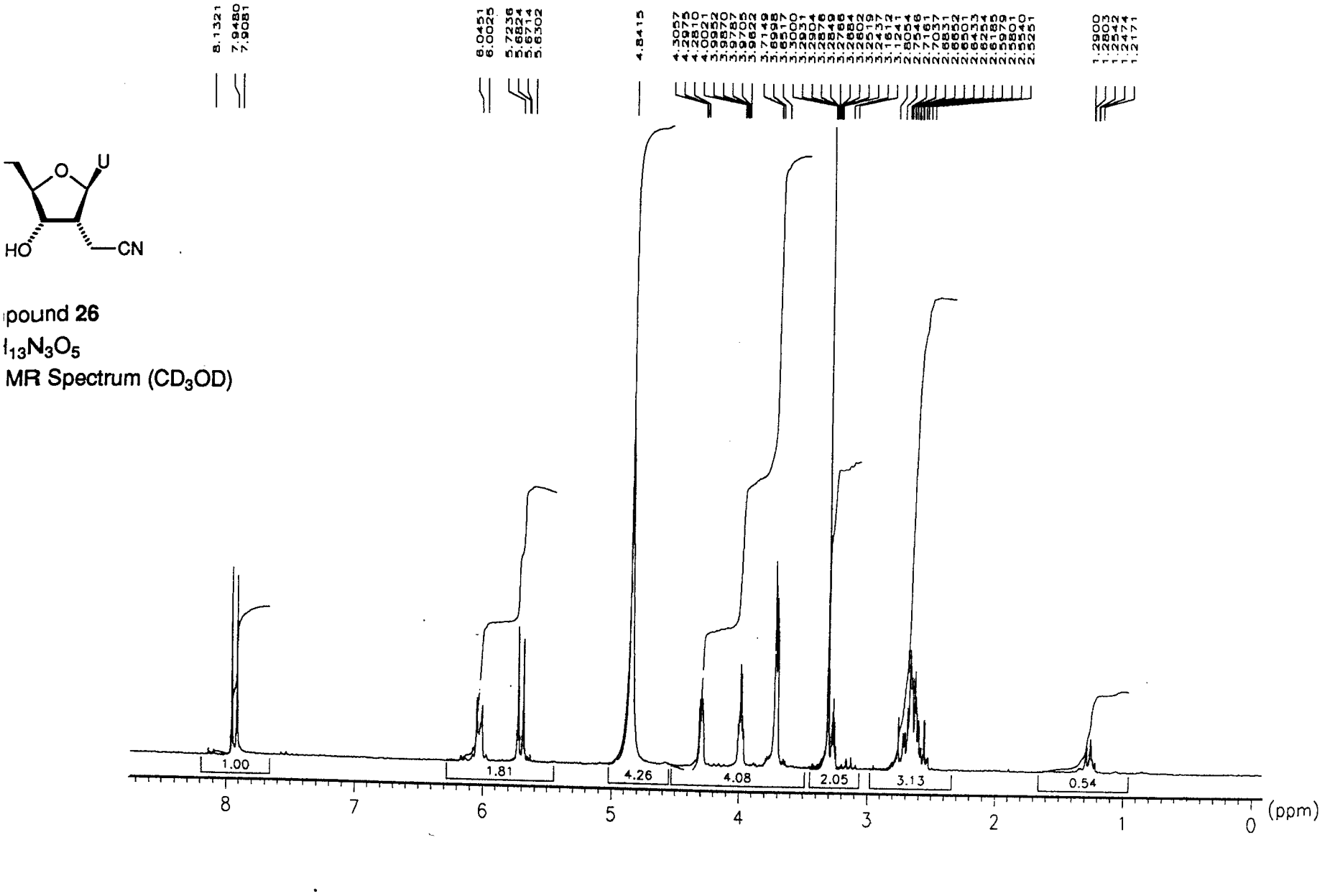
^1H NMR Spectrum ($CDCl_3$)



Compound **25**

$C_{23}H_{39}N_3O_6Si_2$

^1H NMR Spectrum ($CDCl_3$)



pound **26**

$_{13}N_3O_5$

MR Spectrum (CD_3OD)