

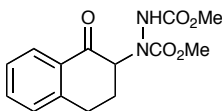
Supplementary Information

Asymmetric α -Amination Reaction of Alkenyl Trifluoroacetates Catalyzed by Chiral
Phosphine-Silver Complex

Akira Yanagisawa*, Ryoji Miyake, and Kazuhiro Yoshida

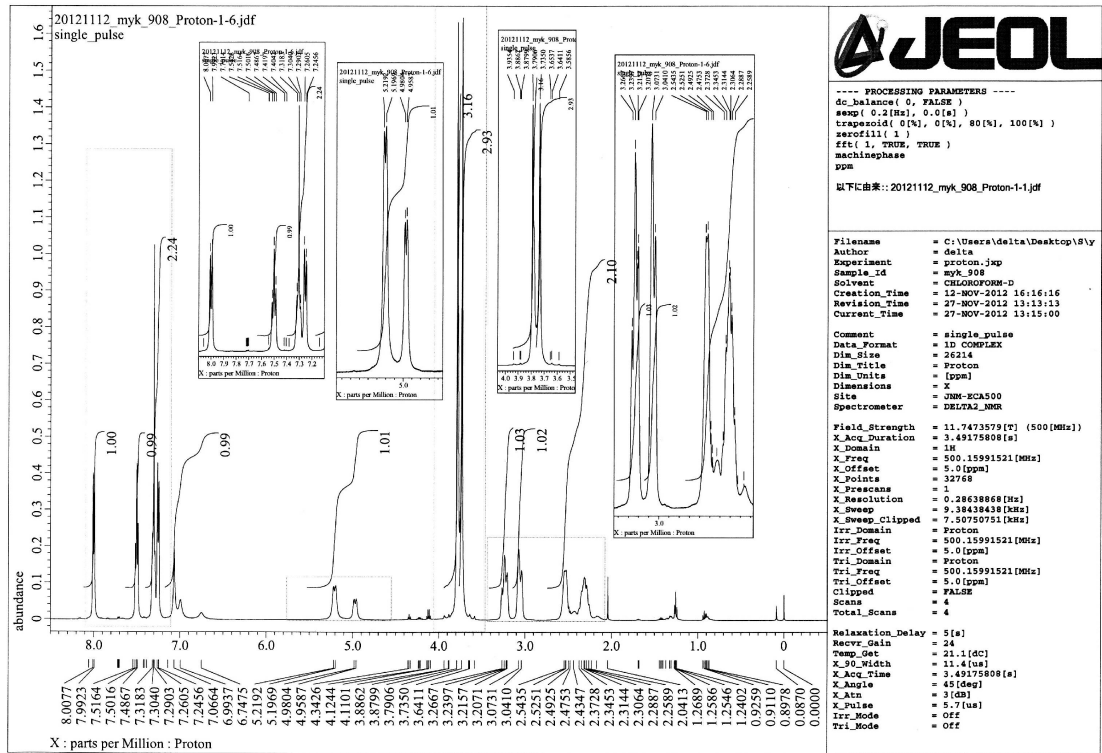
Department of Chemistry, Graduate School of Science, Chiba University, Chiba 263-8522, Japan
ayanagi@faculty.chiba-u.jp

NMR (^1H and ^{13}C) and HPLC Charts for Products in Tables 2 and 3

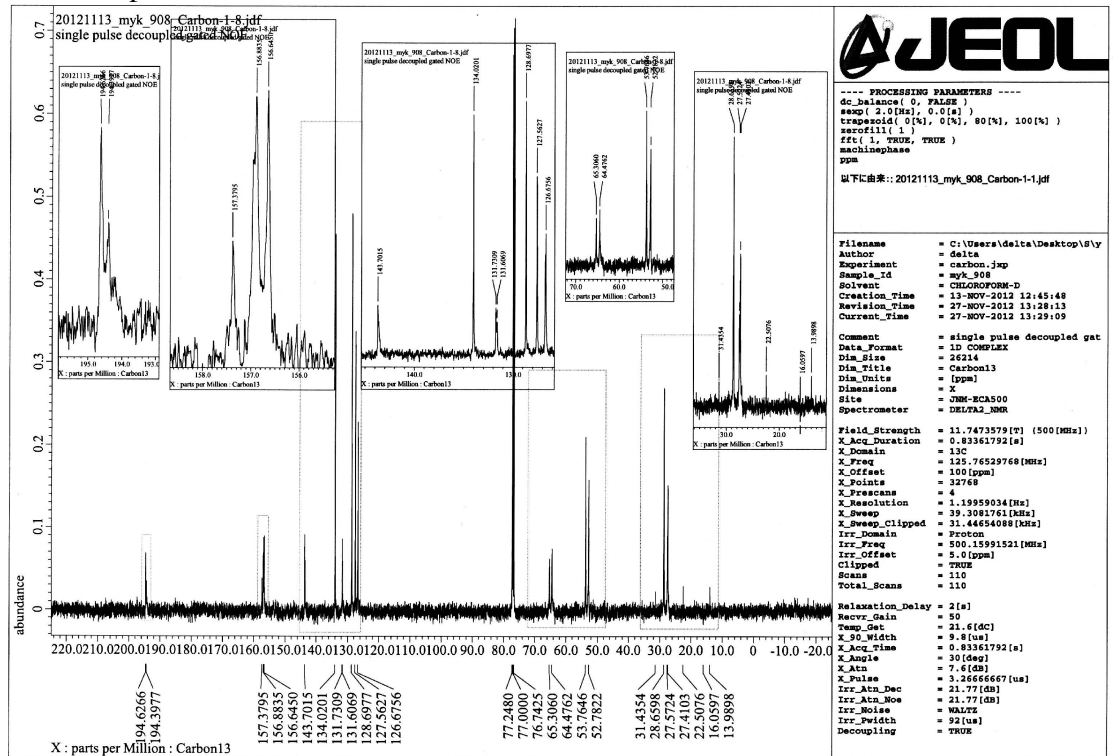


Dimethyl 1-(1-oxo-1,2,3,4-tetrahydronaphthalen-2-yl)hydrazine-1,2-dicarboxylate (3aa, Table 2, entry 1; Table 3, entry 2).

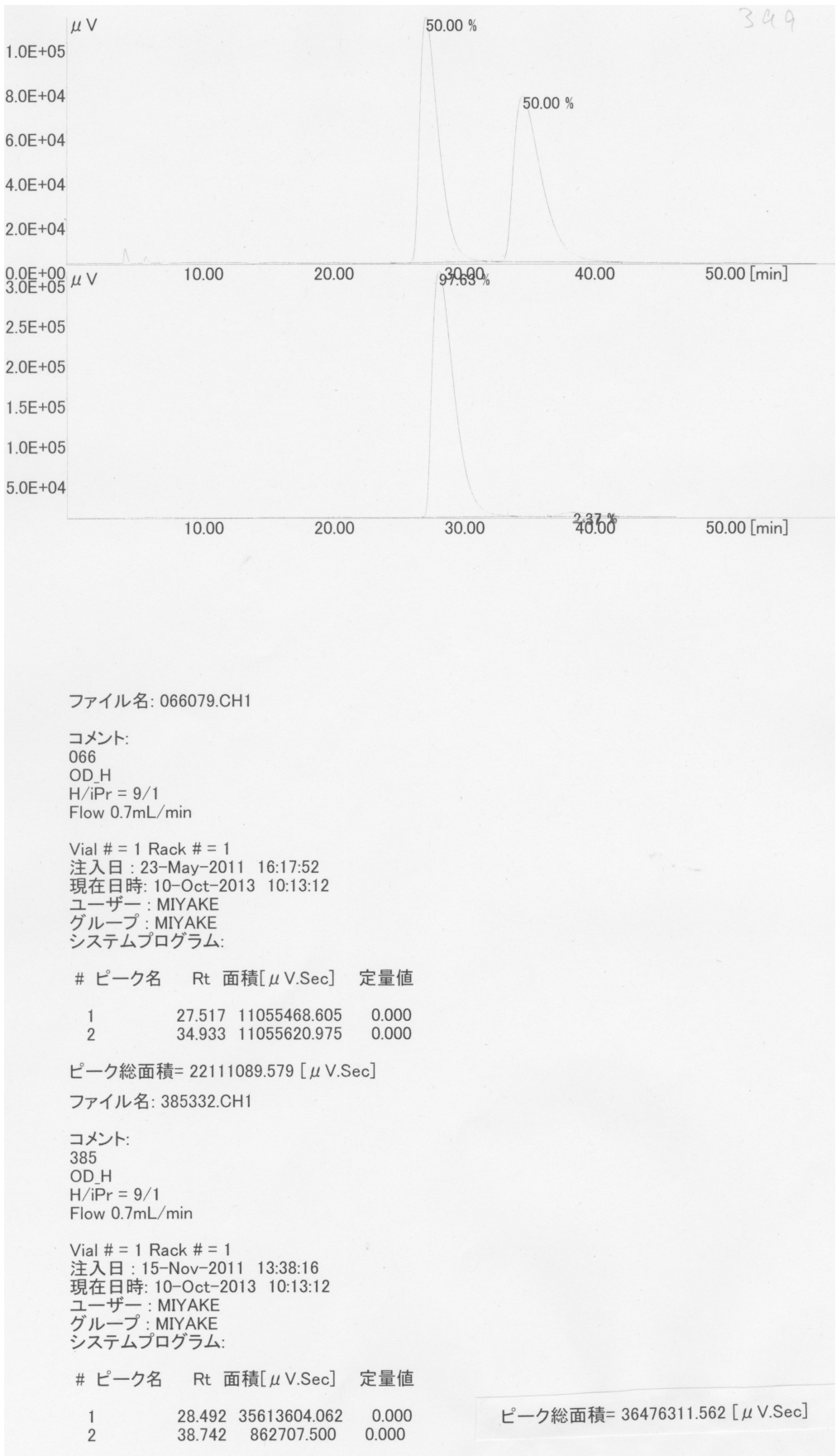
¹H NMR spectra



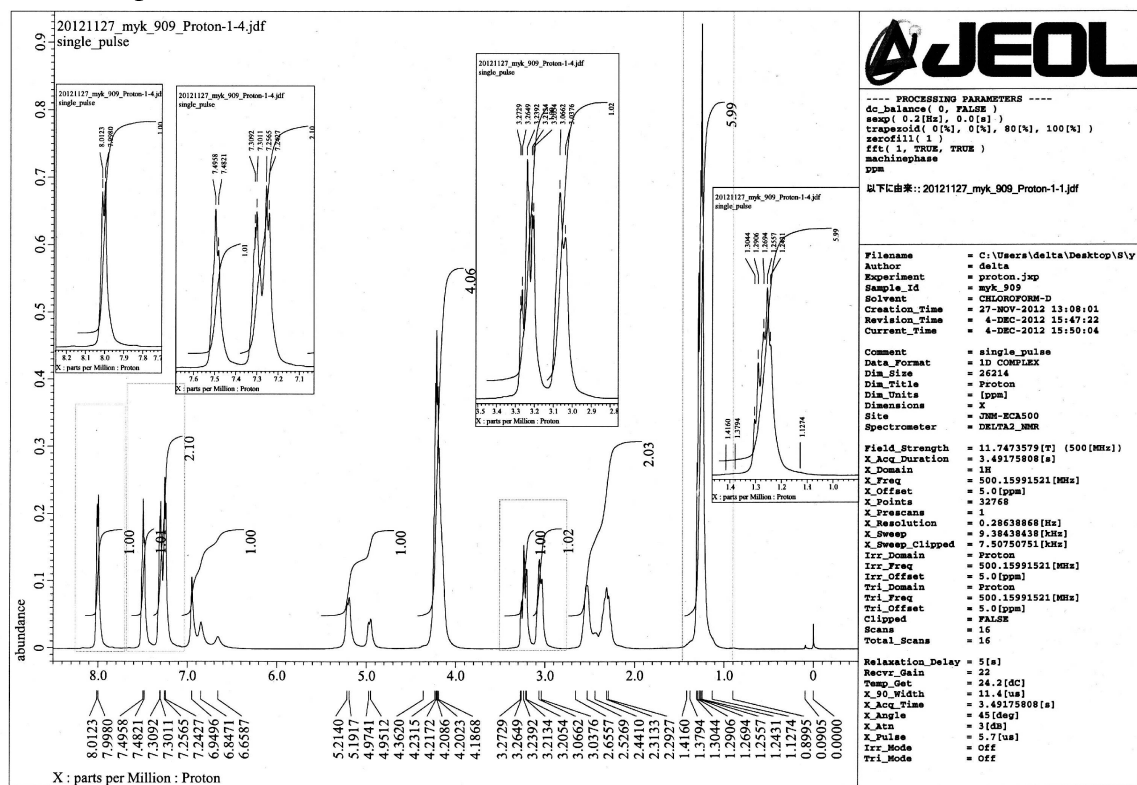
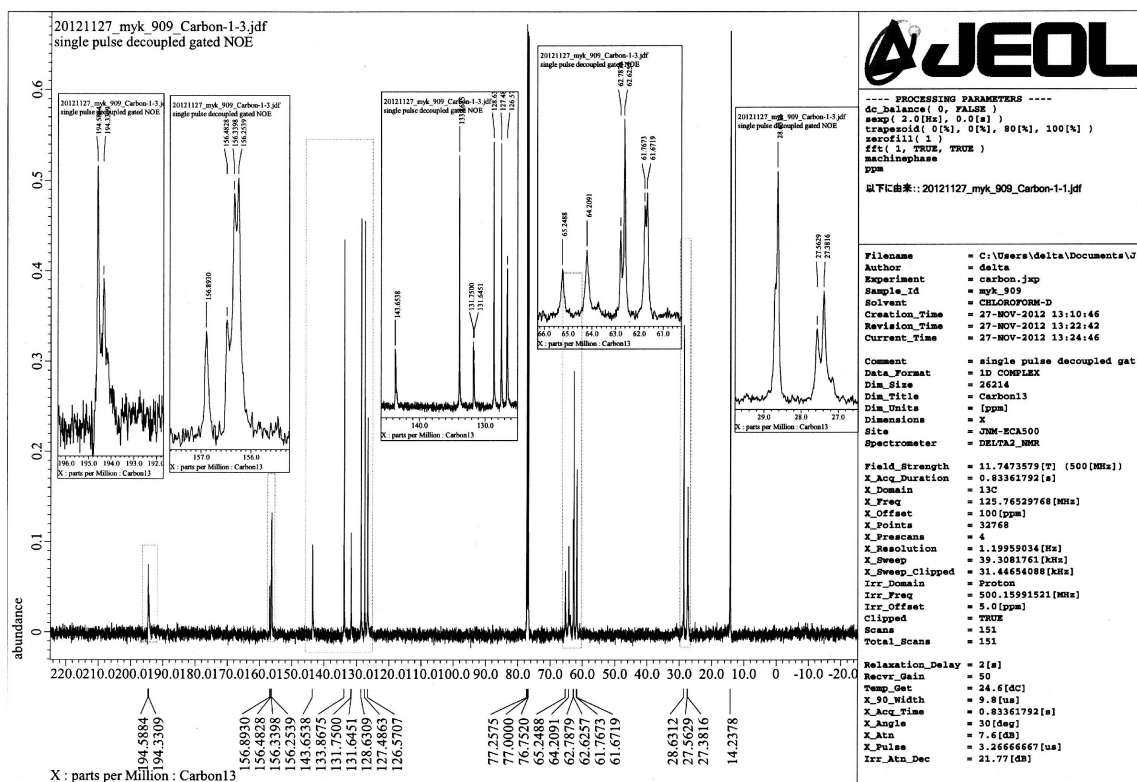
¹³C NMR spectra



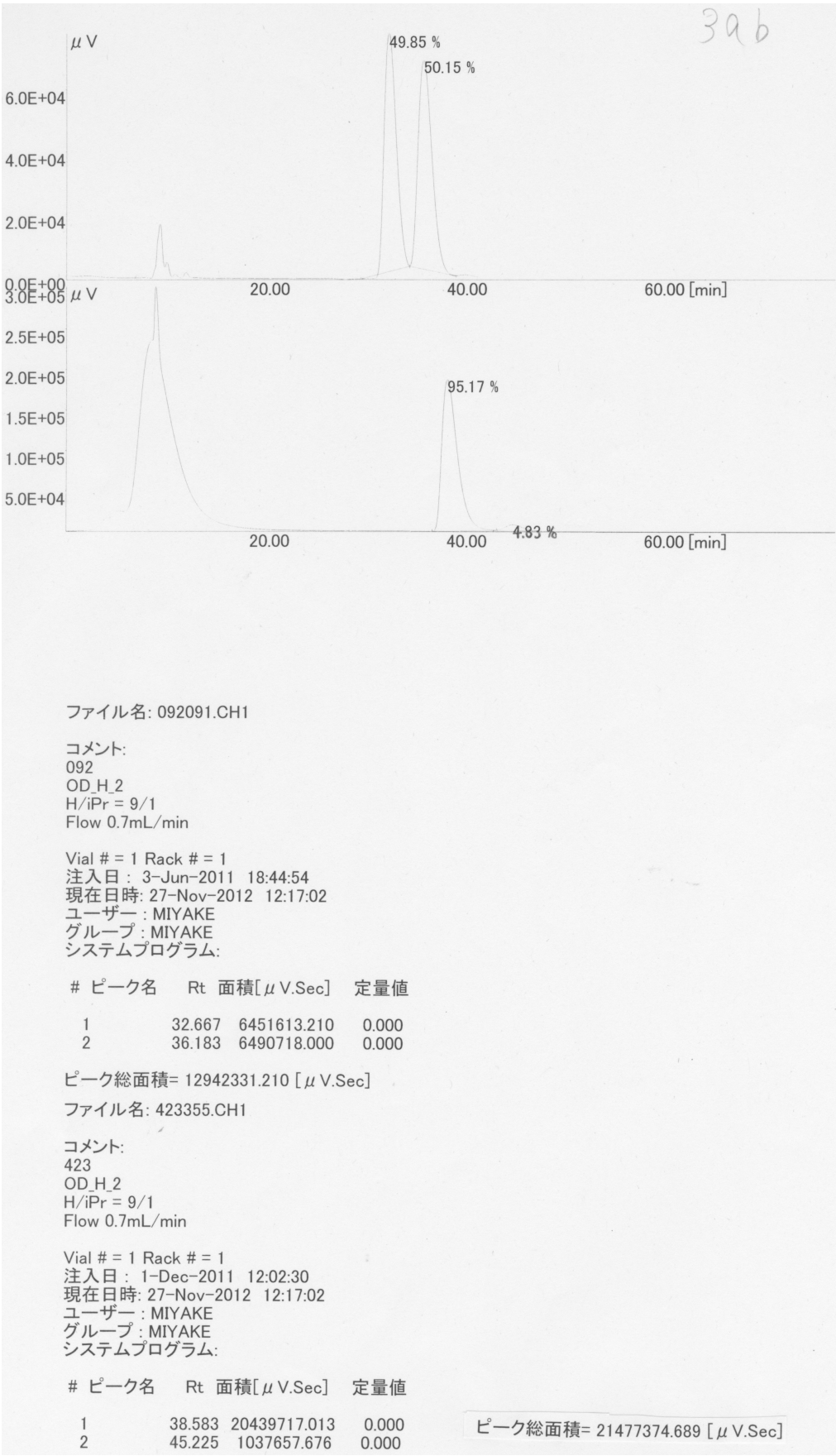
HPLC Chart

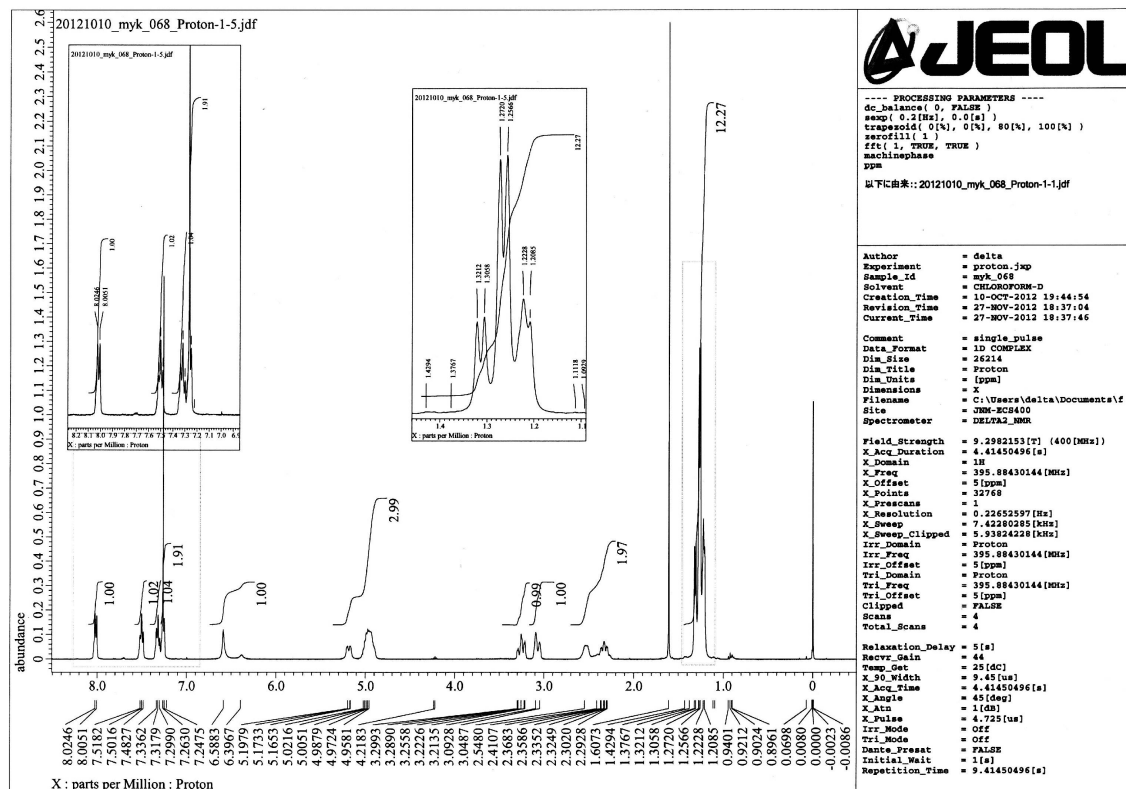
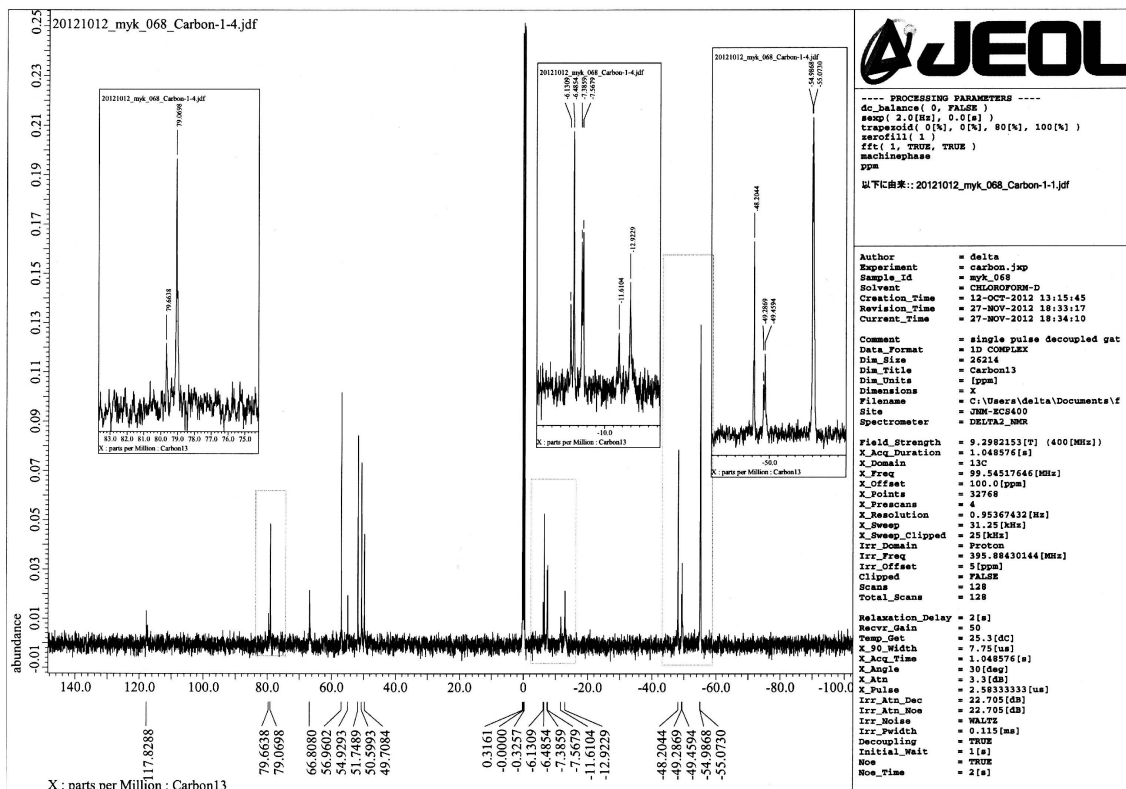


dicarboxylate (3ab, Table 2, entry 2)

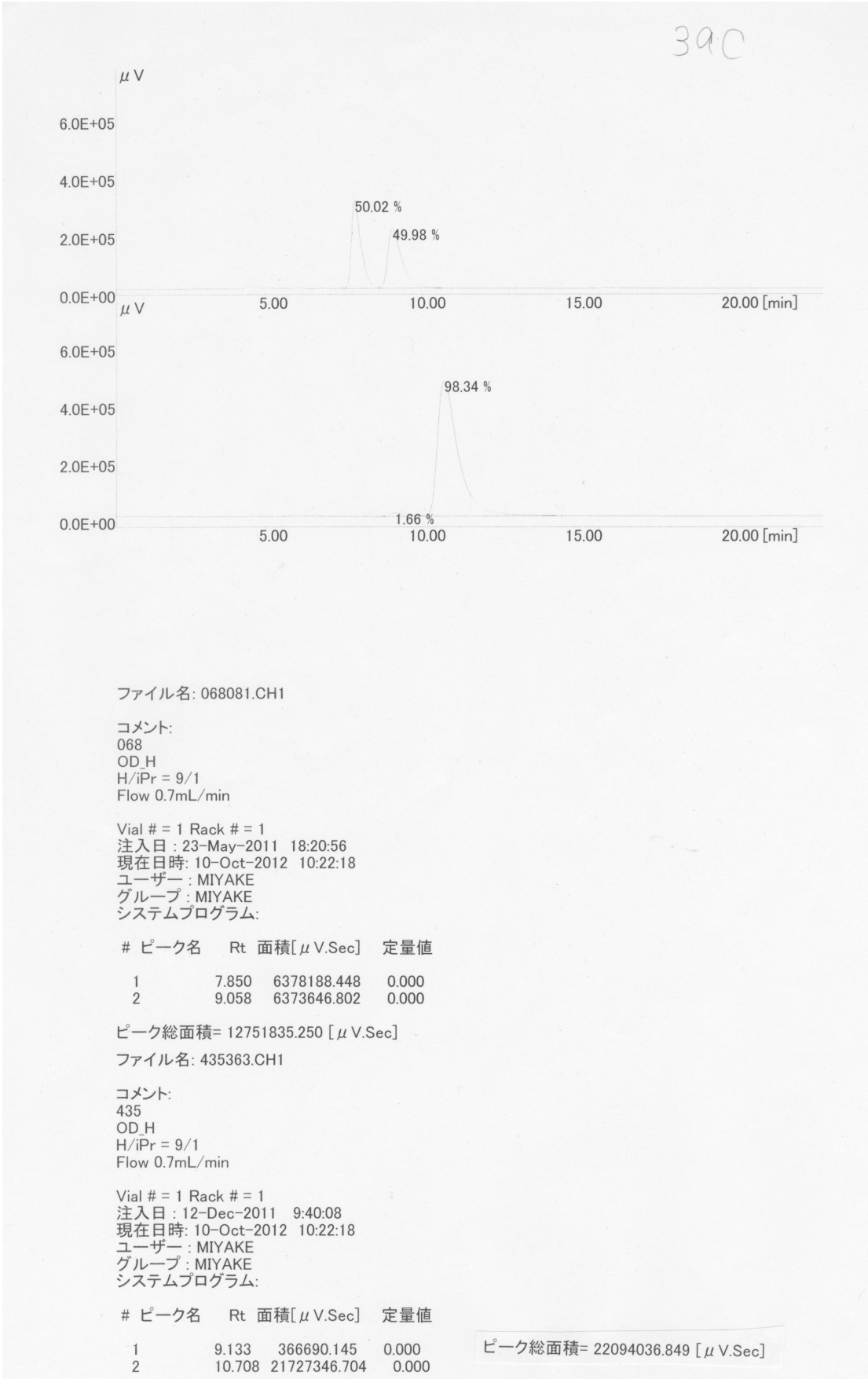
¹H NMR spectra¹³C NMR spectra

HPLC Chart



¹H NMR spectra¹³C NMR spectra

HPLC Chart



Dibenzyl 1-(1-oxo-1,2,3,4-tetrahydronaphthalen-2-yl)hydrazine-1,2-dicarboxylate (3ae, Table 2, entry 5)

20121114_myk_913_Primo-1-6.gdf
single_pulse

abundance

9.0 8.0 7.0 6.0 5.0 4.0 3.0 2.0 1.0 0

8.0020
7.9865
7.5021
7.4878
7.4729
7.3172
7.2920
7.2456
7.0790
6.9393
6.8763
6.6495
5.1997
5.1659
5.1413
5.0606
5.0365
4.9839
4.9547
3.2712
3.2437
3.2105
3.1756
3.1408
3.0993
2.5664
2.5463
2.3305
2.3076
2.06
1.2568
0.8801
0.0779
0.0000

1.00
1.24
0.88
5.32
1.04
1.04
1.05
2.06

X: parts per Million - Proton

```

##### PROCESSING PARAMETERS #####
dc.balance (0, FALSE)
smp 0.2[Hz], 0.0[s]
troughed: 0[N], 0[N], 80[N], 100[N]
zerofill(1)
fft(1, TRUE, TRUE)
machinesphase
ppm

以下由由来: 20121114_myk_913_Proton-1-1.jdf

Filename = C:\Users\delta\Desktop\A\
Author = delta
Experiment = protonen_jmp
Sample_id = myk_913
Solvent = CHLOROFORM-D
Creation_Time = 14-NOV-2013 19:19:15
Revision_Time = 27-NOV-2012 14:03:19
Current_Time = 27-NOV-2012 14:07:23

Comment = single_pulse
Data_Format = ID COMPLEX
Dir_Path =
Dir_Size = 26214
Dir_Title = Proton
Dir_Units = [ppm]
Dimensions =
Site = JNM-ECA500
Spectrometer = DELTA2X_NMR

Field_Strength = 11.7473579[T] (500[MHz])
X_Acq_Duration = 1H1759808[s]
X_Domain = 19
X_Freq = 500.15991521[MHz]
X_Pulse = 5.0[ppm]
X_Points = 32768
X_Prescans = 1
X_Resolution = 0.28638866[Hz]
X_Sweep = 9.38438438[kHz]
X_Sweep_Clippped = 7.50750751[kHz]
X_Domain = Proton
Irr_Freq = 500.15991521[MHz]
Irr_Offset = 5.0[ppm]
Irr_Domain = Proton
Tri_Freq = 500.15991521[MHz]
Tri_Offset = 5.0[ppm]
Tri_Clippped = 5.0[ppm]
Scans = 4
Total_Scans = 4

Relaxation_Delay = 5[s]
Nucvcr_Gain = 32
Temp_Set = 32.3[°C]
X_90_Width = 1.442[us]
X_90_Time = 3.49575808[s]
X_Angle = 30[°]
X_Atn = 3[dB]
X_Pulse = 5.7[us]
Irr_Mode = Z
Tri_Mode = Off

```

20121025_msk_427_Carbon-1-4.jdf

Abundance vs. X: parts per Million : Carbon13

Labeled Peaks (m/z):

- 94.5505
- 94.3014
- 156.7657
- 156.5645
- 156.1813
- 143.7461
- 143.6790
- 135.7369
- 135.5453
- 134.0891
- 131.8186
- 128.7624
- 128.4846
- 128.3643
- 127.9872
- 127.6703
- 127.5170
- 126.7889
- 77.3161
- 77.2012
- 77.0000
- 76.6743
- 68.6747
- 68.2915
- 67.6592
- 67.5538
- 65.4749
- 64.4785
- 28.7344
- 27.6805
- 27.5081

Inset 1 (Top Left): 196.0, 194.0, 196.5614, 194.5505

Inset 2 (Top Middle): 156.7657, 156.5645, 156.1813

Inset 3 (Top Right): 143.7461, 143.6790

Inset 4 (Middle): 136.0, 134.0, 136.0891, 134.0896

Inset 5 (Bottom Right): 70.0, 68.0, 70.0000, 68.2915, 67.6592, 65.4749, 64.4785

```

----- PROCESSING PARAMETERS -----
dc_balance! 0, FALSE
amp! 2.0[Hz], 0.5[Hz]
treshold! 0[Hz], 0[N], 80[N], 100[N]
zerofill! 1
fft! 1, TRUE, TRUE
machinphase
ups

以下に由来： 20121025_mysk_427_Carboten-1-1.jdf

Author = delta
Experiment = carboten_jmp
Sample_ID = mysk_427
Solvent = CHLOROFORM-D
Revision_time = ORG-2012 21:07:07
Revision_time = 27-MOV-2012 18:29:09
Current_time = 27-MOV-2012 18:30:47

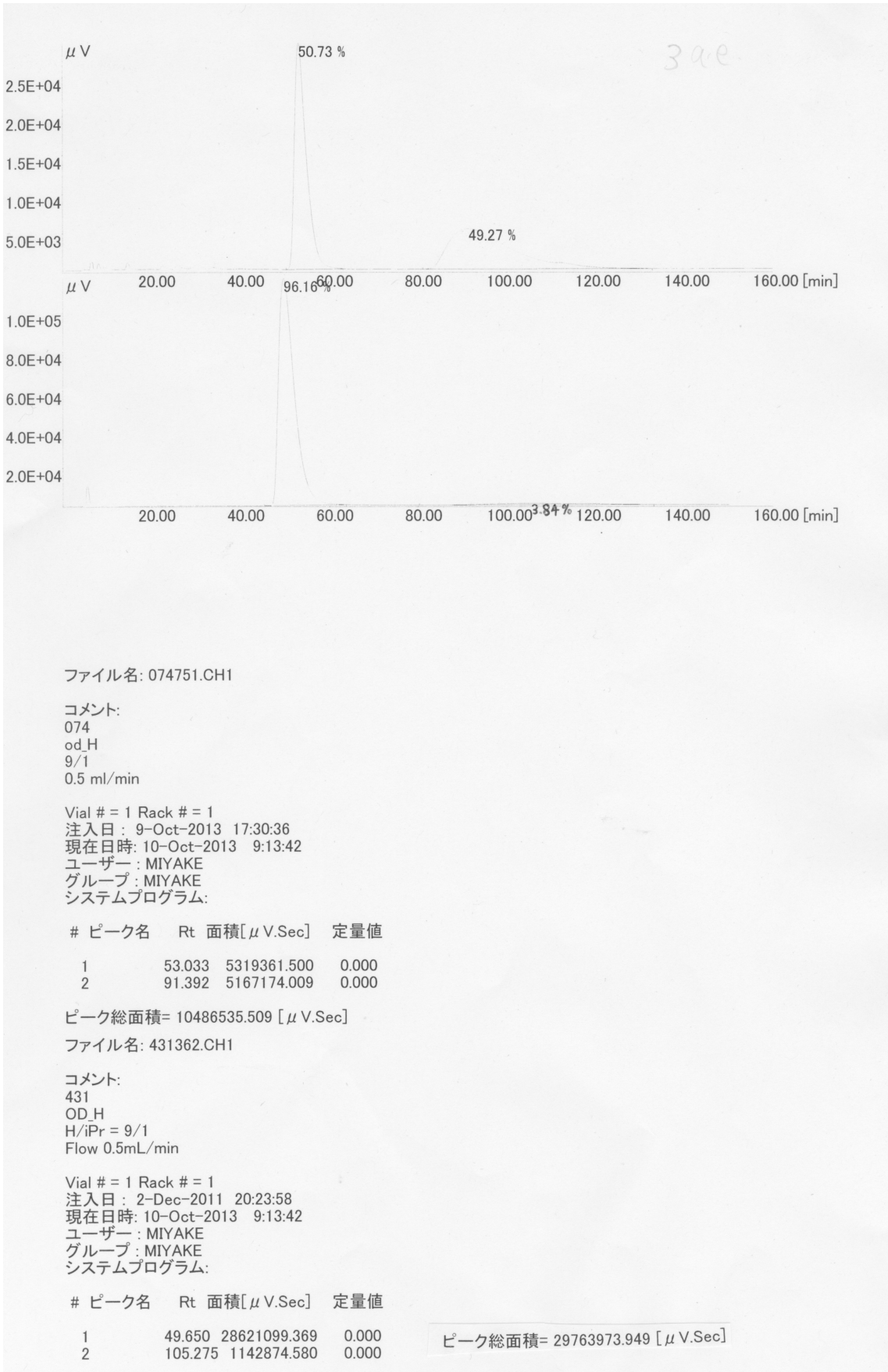
Comment = single pulse decoupled gat
Data_Format = 1D COMPLEX
Dir_Path = C:\Users\delta\Documents\1f
Dir_Unit = CHLOROACID
Dimensions = DELTA72_BMRK
Filebase = C:\Users\delta\Documents\1f
Spectrometer = DELTA72_BMRK

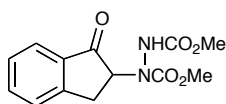
Field_Strength = 9.29825153 [7] 400[MHz]
X_Acq_Rotation_Delay = 1.048576[s]
X_Domain = 35
X_Offset = 99.54517646[MHz]
X_Fpoints = 127.00[pptm]
X_Prescans = 4
X_Resolution = 0.95367432[Hz]
X_Sweep = 31.25 [Hz/s]
X_Sweep_Clippped = 25 [Hz]
Xrr_Domain = 35
Xrr_Offset = 995.88430144[MHz]
Xrr_Clipped = 25 [ppm]
Scans = 372
Total_Scans = 372

Relaxation_Delay = 2[s]
Recur_Dein = 2[s]
Temp_Opt = 8.0[°C]
X90_Width = 7.75 [us]
X_Acq_Time = 48.64576[s]
X_Angle = 30[deg]
X_Atn = 3.3[db]
X_Pulse = 18.833333[us]
Xrr_Acta_Dec = 22.705 [dB]
Xrr_Acta_Nuc = 22.705 [dB]
Xrr_Raise = 22.705 [dB]
Xrr_Fwidth = 0.115[mHz]
Decoupling = TRUE
Initial_Wait = 1[s]
Noise = TRUE
Now_Time = 2[s]

```

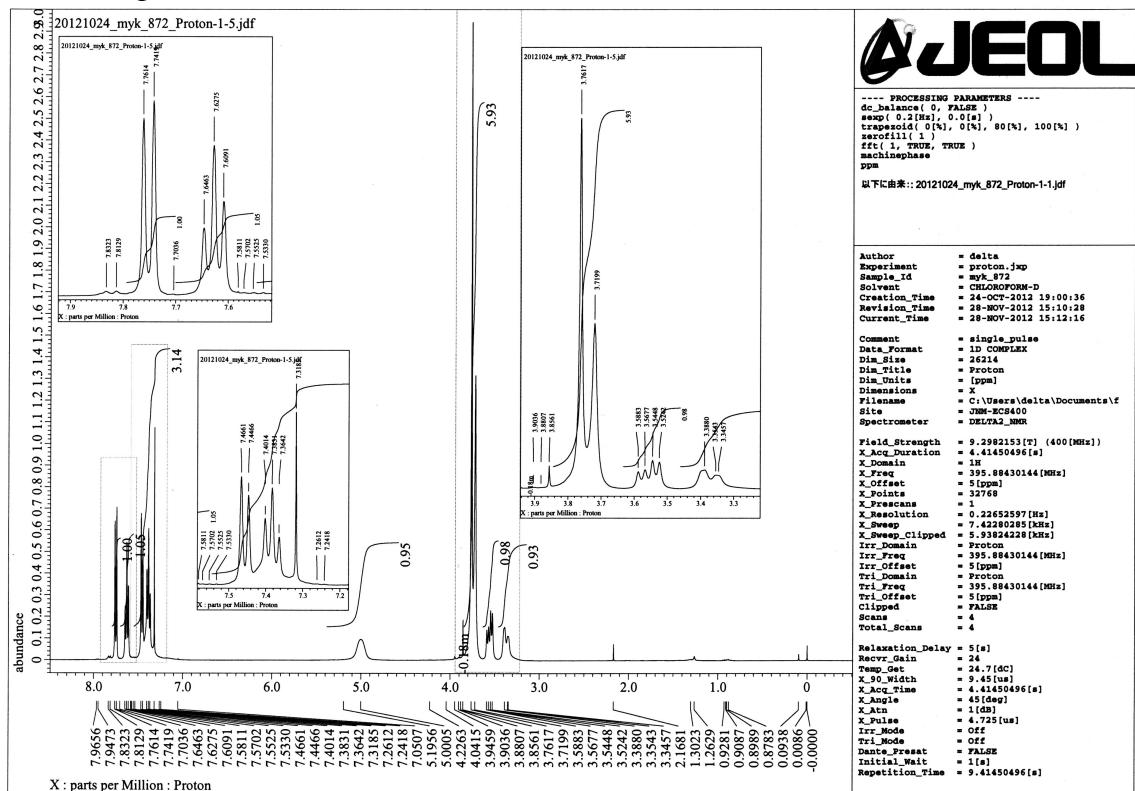

HPLC Chart



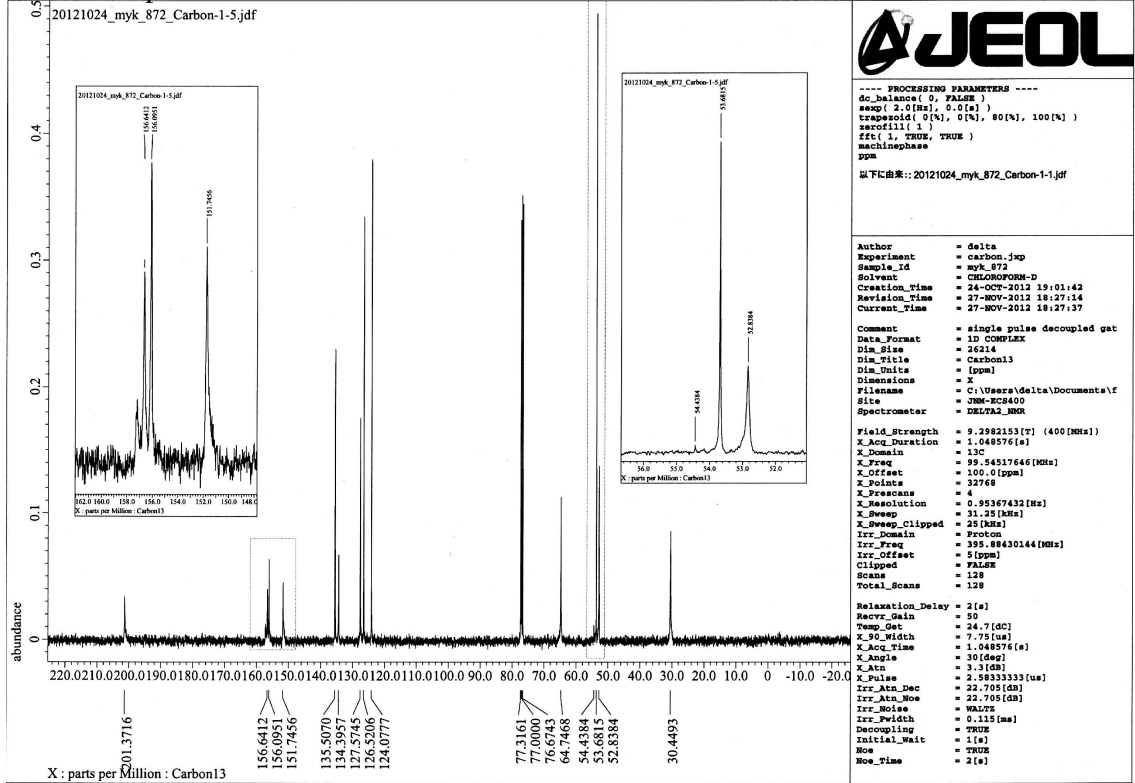


Dimethyl 1-(1-oxo-2,3-dihydro-1H-inden-2-yl)hydrazine-1,2-dicarboxylate
(3ba, Table 3, Entry 1)

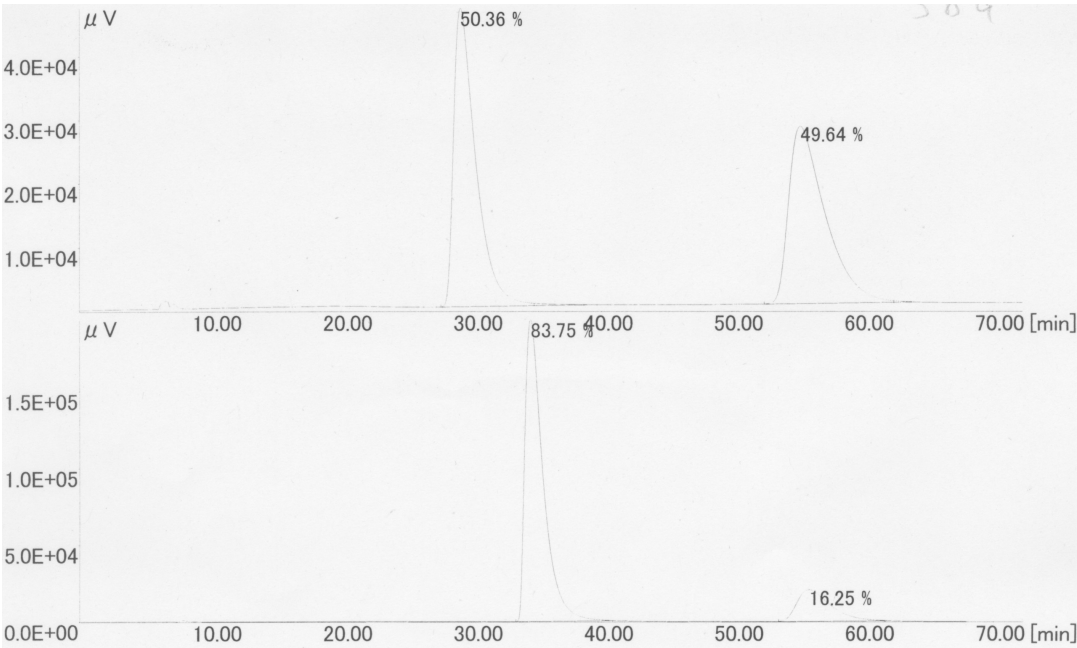
¹H NMR spectra



¹³C NMR spectra



HPLC Chart



ファイル名: 398750.CH1

コメント:

398

od_H

9/1

1.0 ml/min

Vial # = 1 Rack # = 1

注入日: 9-Oct-2013 16:08:28

現在日時: 10-Oct-2013 10:09:34

ユーザー: MIYAKE

グループ: MIYAKE

システムプログラム:

#	ピーク名	Rt	面積[μ V.Sec]	定量値
1		29.192	5458424.021	0.000
2		55.283	5379428.912	0.000

ピーク総面積= 10837852.933 [μ V.Sec]

ファイル名: 444380.CH1

コメント:

445

OD_H

H/iPr = 9/1

Flow 1.0mL/min

Vial # = 1 Rack # = 1

注入日: 13-Dec-2011 16:35:38

現在日時: 10-Oct-2013 10:09:34

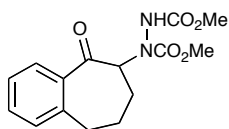
ユーザー: MIYAKE

グループ: MIYAKE

システムプログラム:

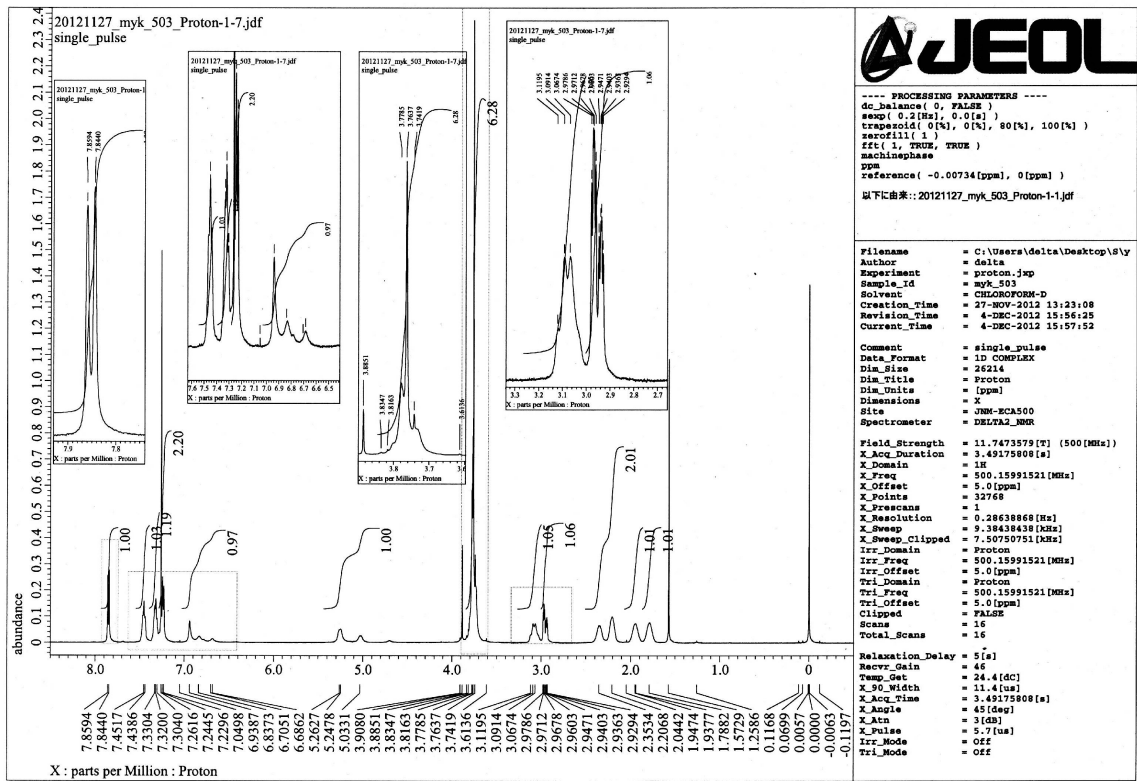
#	ピーク名	Rt	面積[μ V.Sec]	定量値
1		34.567	18837725.500	0.000
2		55.908	3654452.000	0.000

ピーク総面積= 22492177.500 [μ V.Sec]

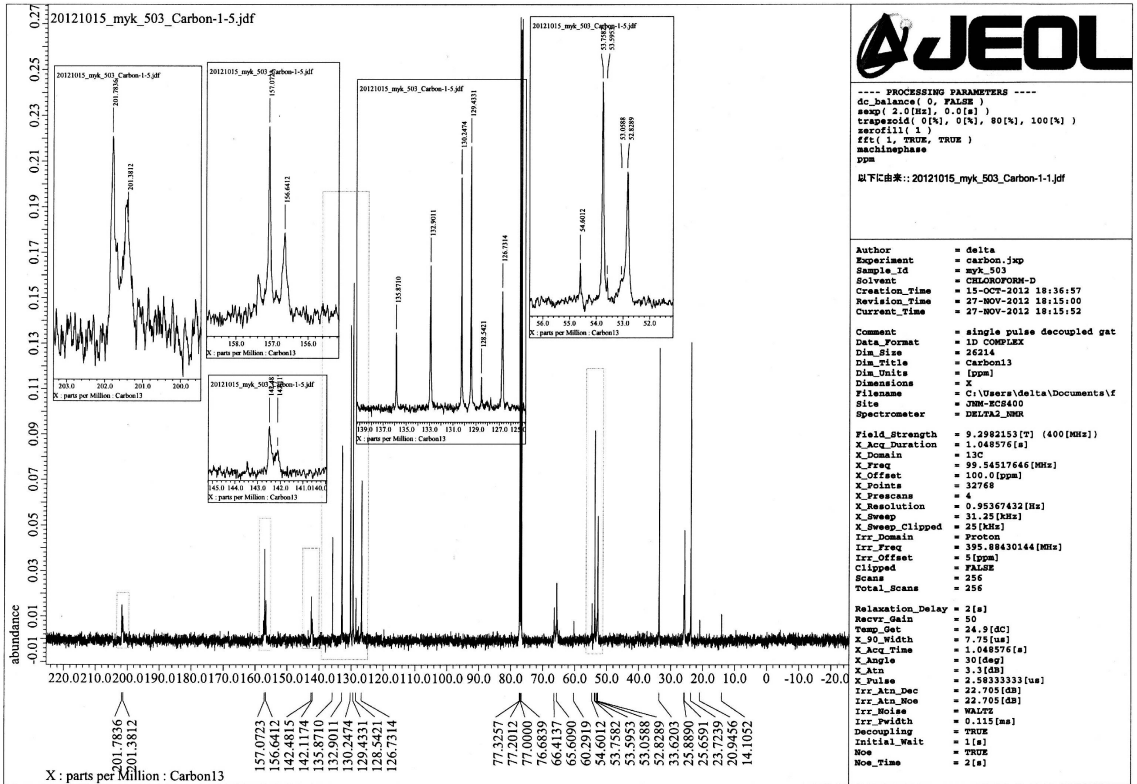


Dimethyl 1-(5-oxo-6,7,8,9-tetrahydro-5H-benzo[7]annulen-6-yl)hydrazine-1,2-dicarboxylate (3ca, Table 3, entry 4)

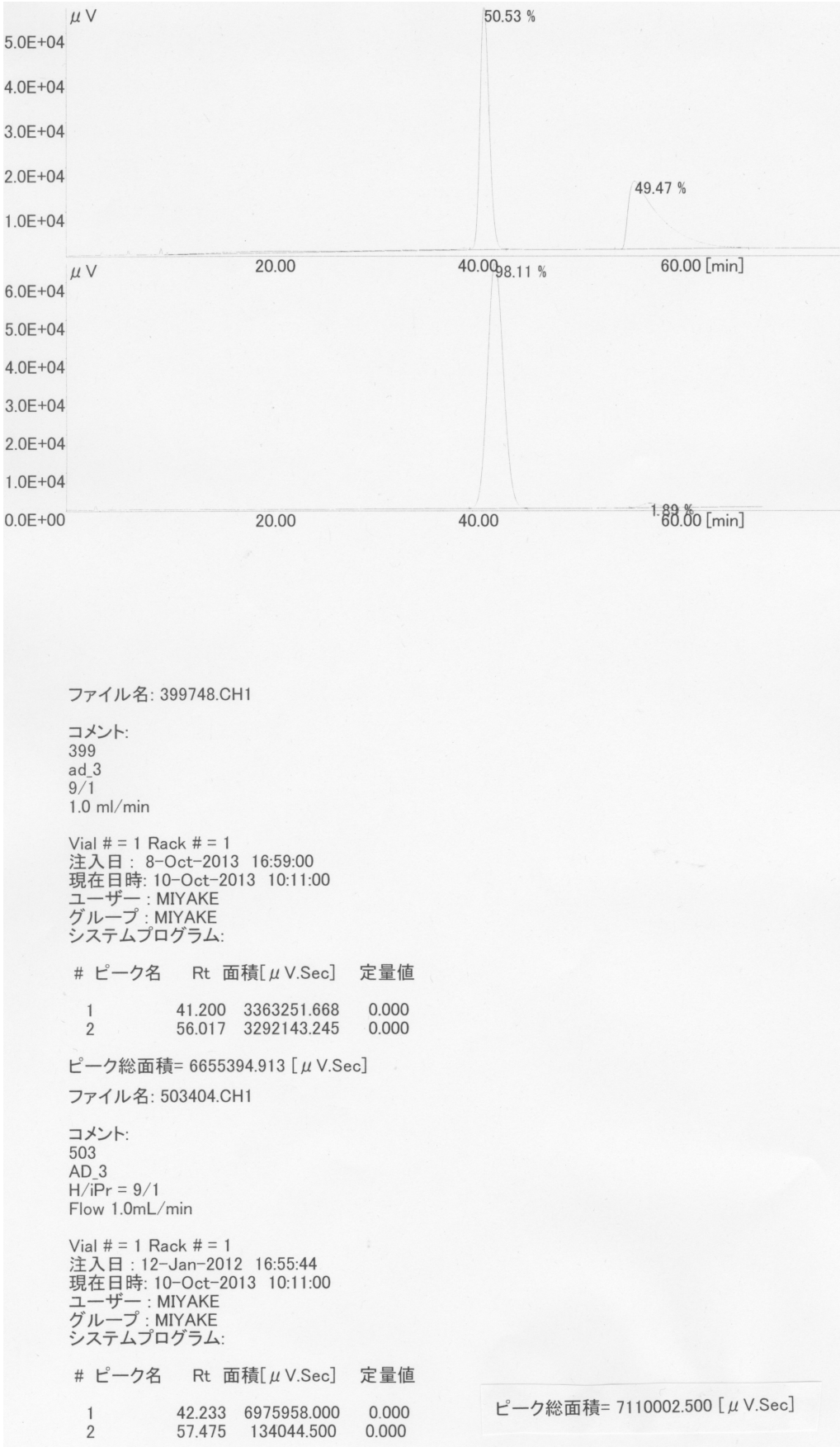
¹H NMR spectra

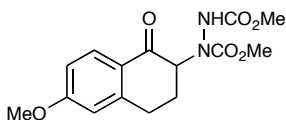


¹³C NMR spectra



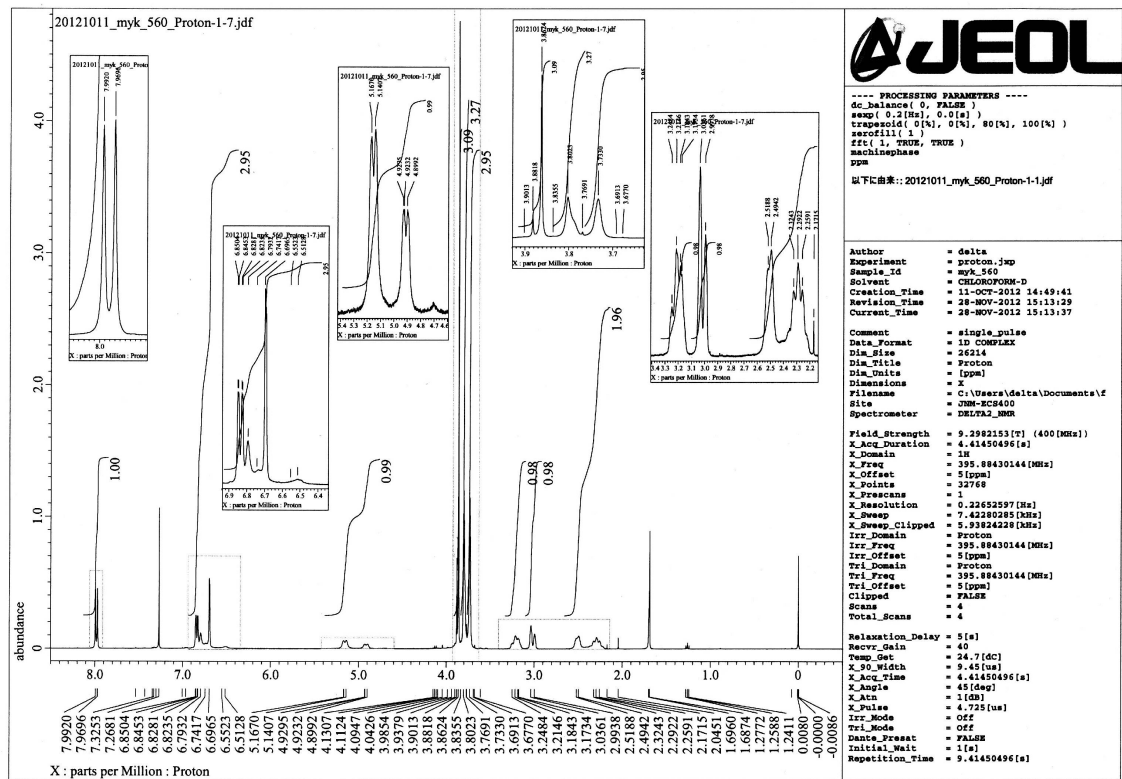
HPLC Chart



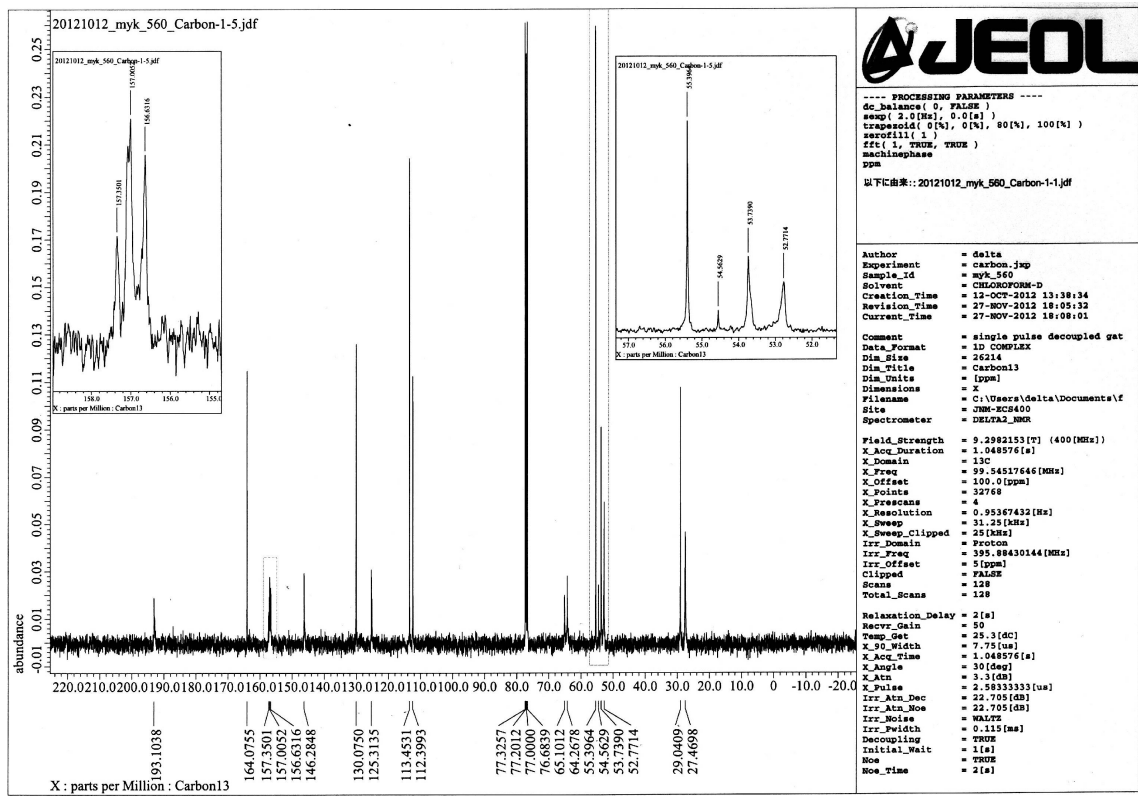


Dimethyl 1-(6-methoxy-1-oxo-1,2,3,4-tetrahydronaphthalen-2-yl)hydrazine-1,2-dicarboxylate (3da, Table 3, entry 5)

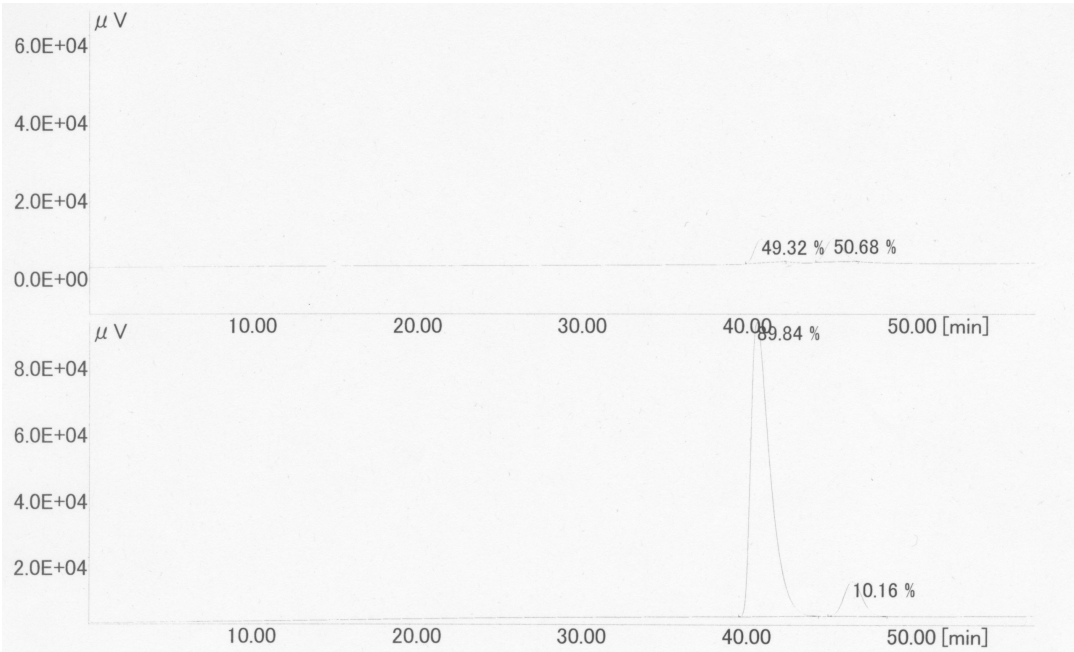
¹H NMR spectra



¹³C NMR spectra



HPLC Chart



ファイル名: 396596.CH1

コメント:

396

ad_3_ad_h

H/iPr = 3/1

Flow 1.0mL/min

Vial # = 1 Rack # = 1

注入日: 21-Nov-2012 11:13:34

現在日時: 27-Nov-2012 10:33:00

ユーザー: MIYAKE

グループ: MIYAKE

システムプログラム:

#	ピーク名	Rt	面積[μ V.Sec]	定量値
1		41.283	406338.250	0.000
2		45.650	417473.500	0.000

ピーク総面積= 823811.750 [μ V.Sec]

ファイル名: 916595.CH1

コメント:

916

ad_3_ad_h

H/iPr = 3/1

Flow 1.0mL/min

Vial # = 1 Rack # = 1

注入日: 21-Nov-2012 10:12:42

現在日時: 27-Nov-2012 10:33:00

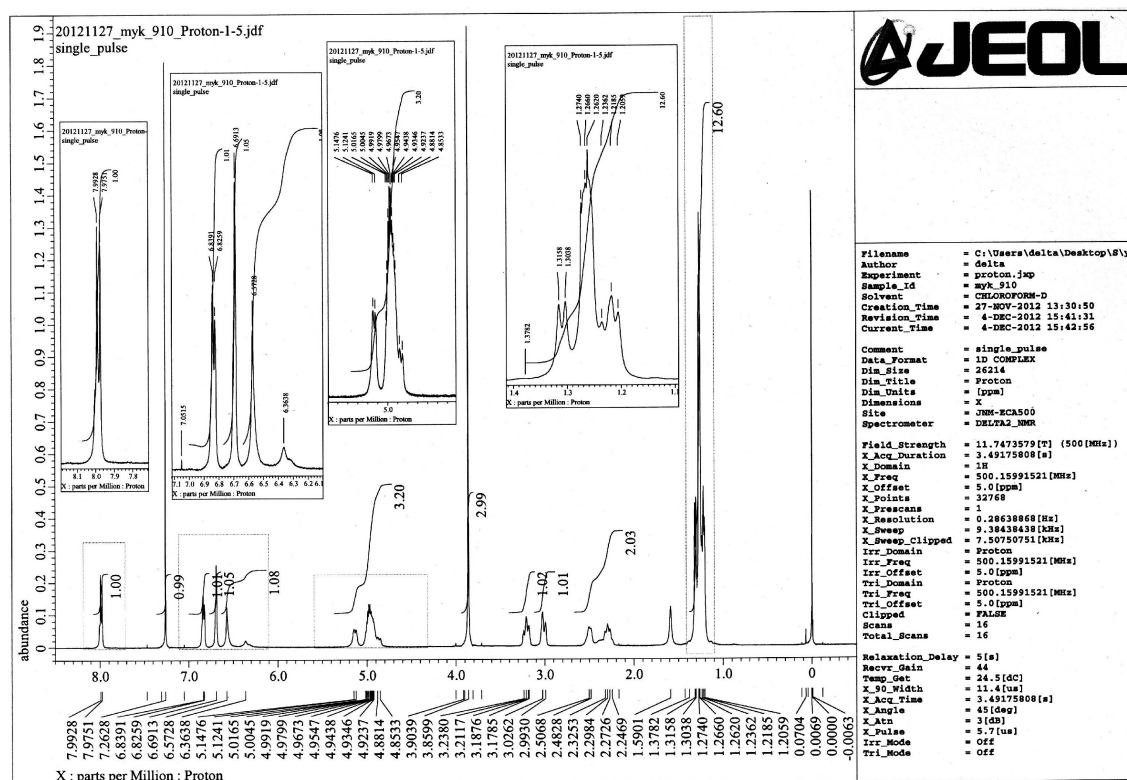
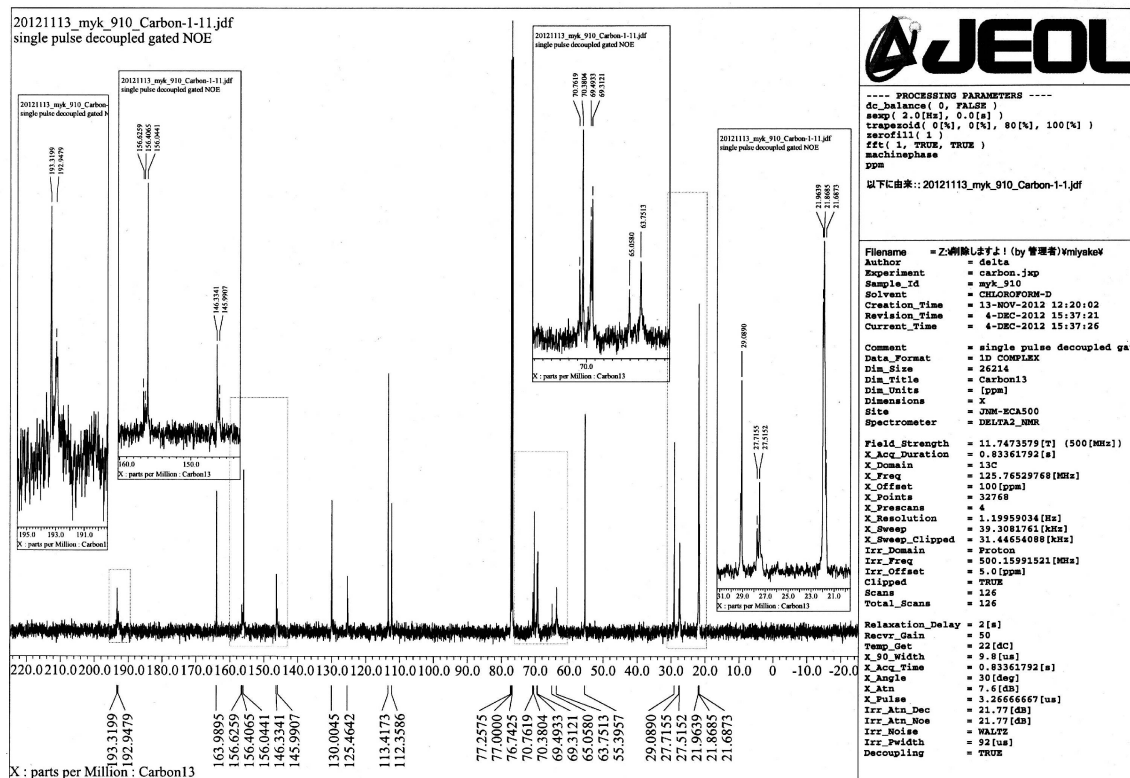
ユーザー: MIYAKE

グループ: MIYAKE

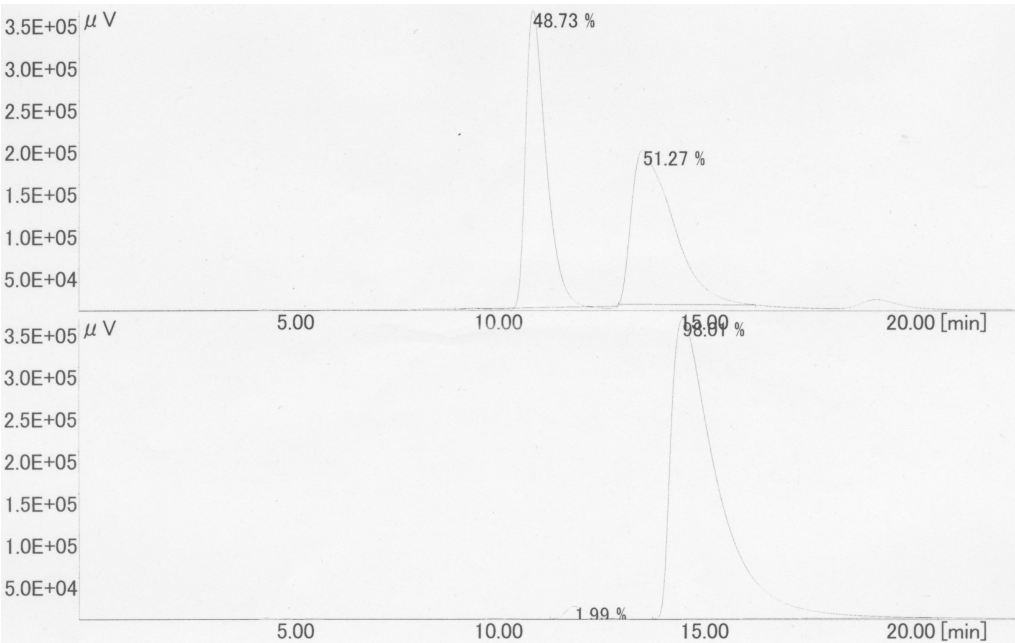
システムプログラム:

#	ピーク名	Rt	面積[μ V.Sec]	定量値
1		41.025	7043375.000	0.000
2		46.842	796658.250	0.000

ピーク総面積= 7840033.250 [μ V.Sec]

¹H NMR spectra20121113_myk_910_Carbon-1-11.jdf
single pulse decoupled gated NOE

HPLC Chart



ファイル名: 236387.CH1

コメント:

236

OD_H

H/iPr = 9/1

Flow 0.7mL/min

Vial # = 1 Rack # = 1

注入日: 5-Jan-2012 12:40:36

現在日時: 10-Oct-2013 10:37:04

ユーザー: MIYAKE

グループ: MIYAKE

システムプログラム:

#	ピーク名	Rt	面積[$\mu V \cdot Sec$]	定量値
1		11.017	12177013.311	0.000
2		13.683	12813421.500	0.000

ピーク総面積= 24990434.811 [$\mu V \cdot Sec$]

ファイル名: 502411.CH1

コメント:

502

OD_H

H/iPr = 9/1

Flow 0.7mL/min

Vial # = 1 Rack # = 1

注入日: 13-Jan-2012 3:05:36

現在日時: 10-Oct-2013 10:37:04

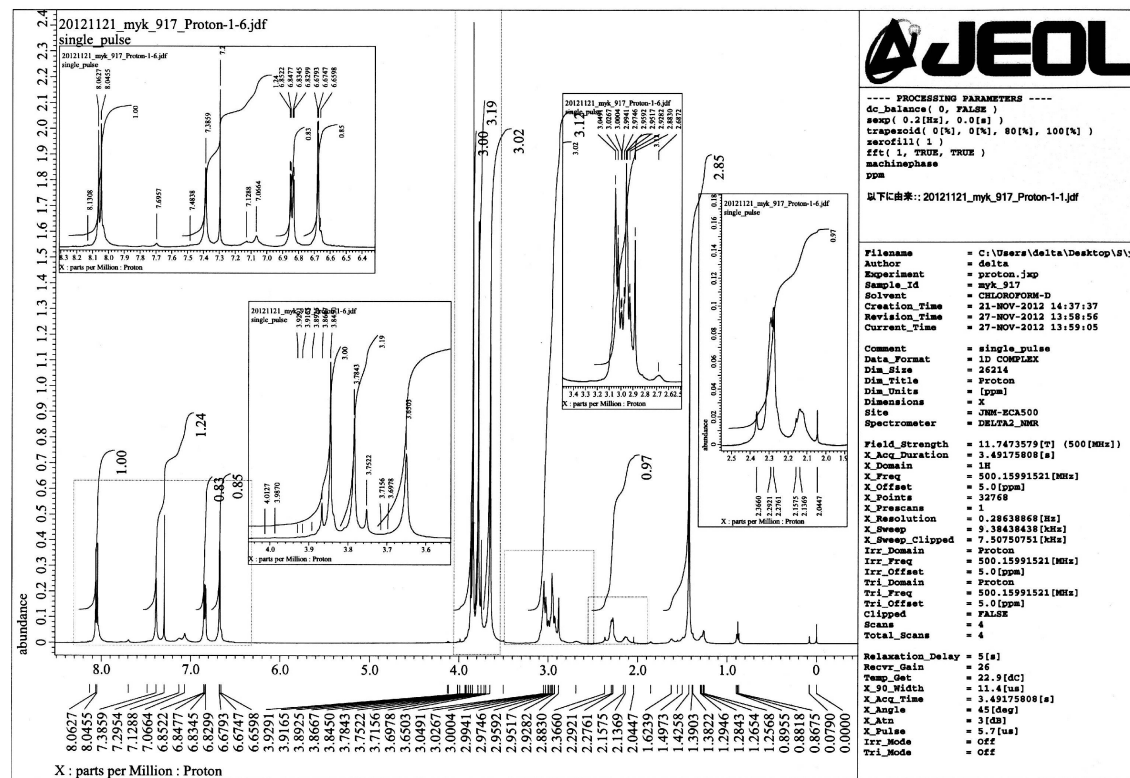
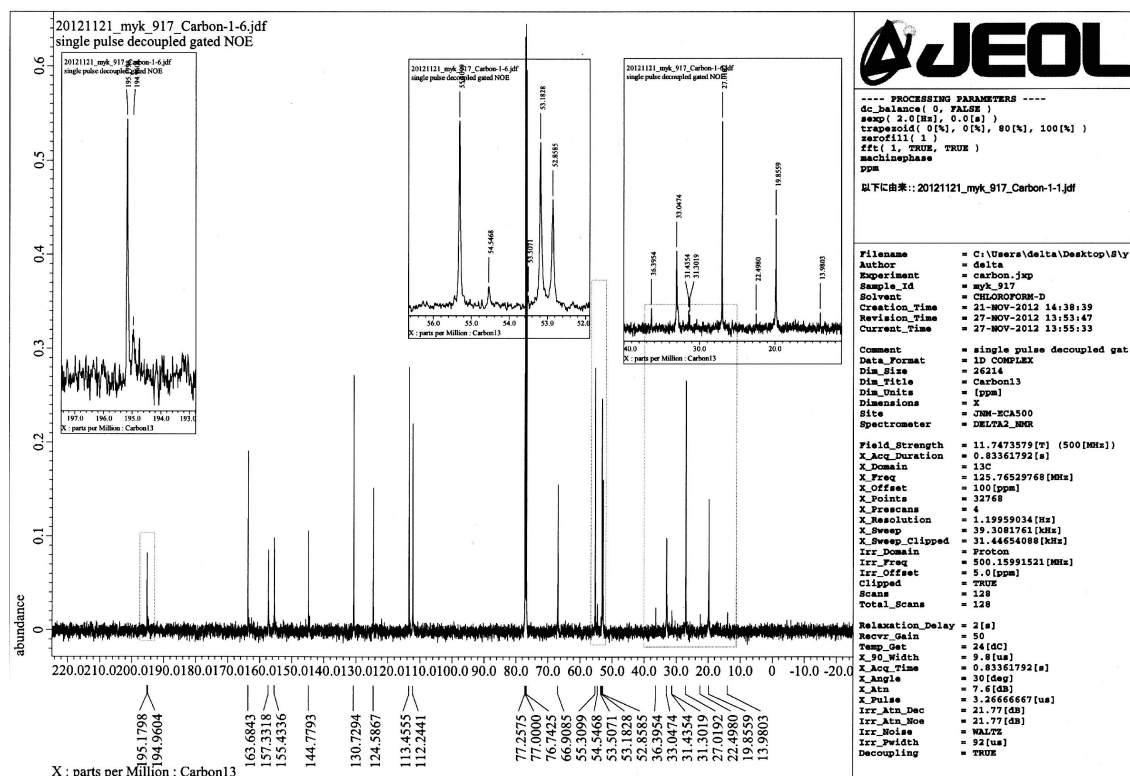
ユーザー: MIYAKE

グループ: MIYAKE

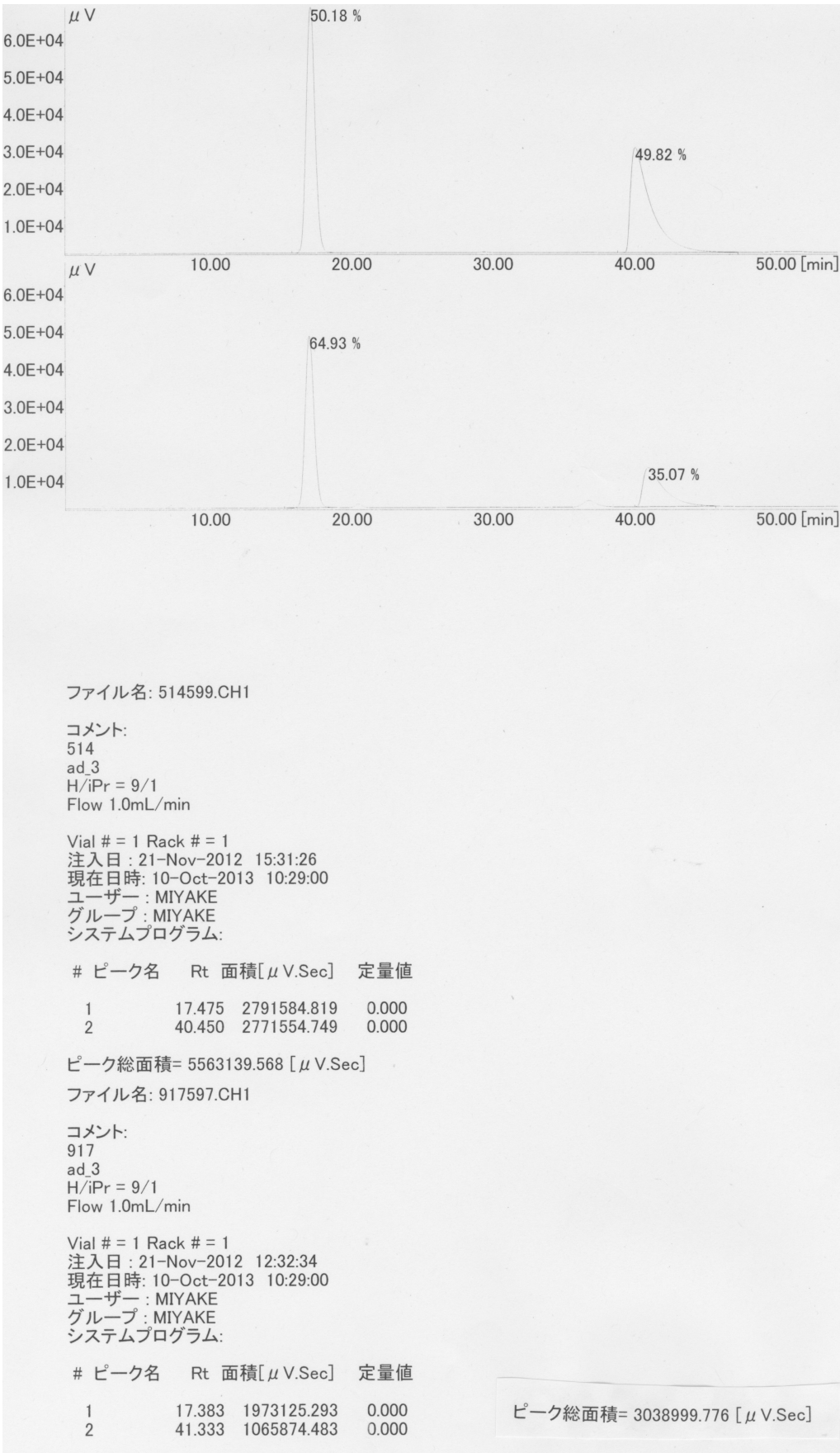
システムプログラム:

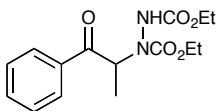
#	ピーク名	Rt	面積[$\mu V \cdot Sec$]	定量値
1		12.025	563343.000	0.000
2		14.650	27728015.872	0.000

ピーク総面積= 28291358.872 [$\mu V \cdot Sec$]

¹H NMR spectra¹³C NMR spectra

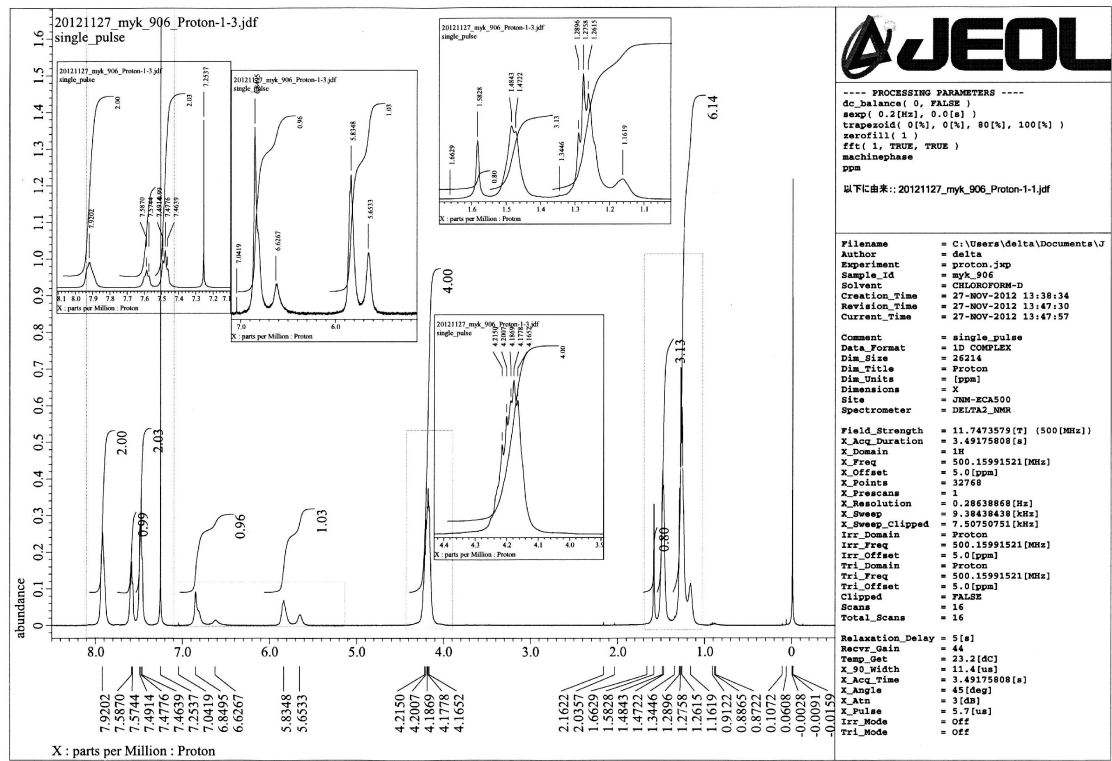
HPLC Chart



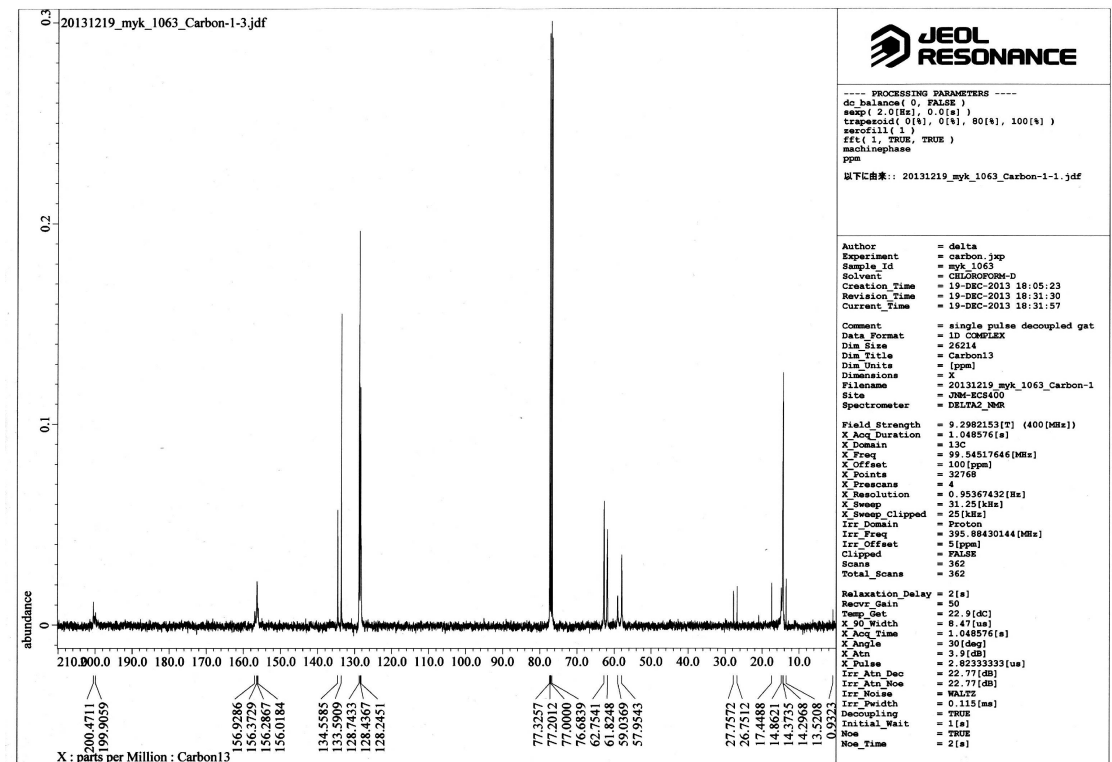


(S)-Diethyl 1-(1-oxo-1-phenylpropan-2-yl)hydrazine-1,2-dicarboxylate (3fb,
Table 3, entry 8)

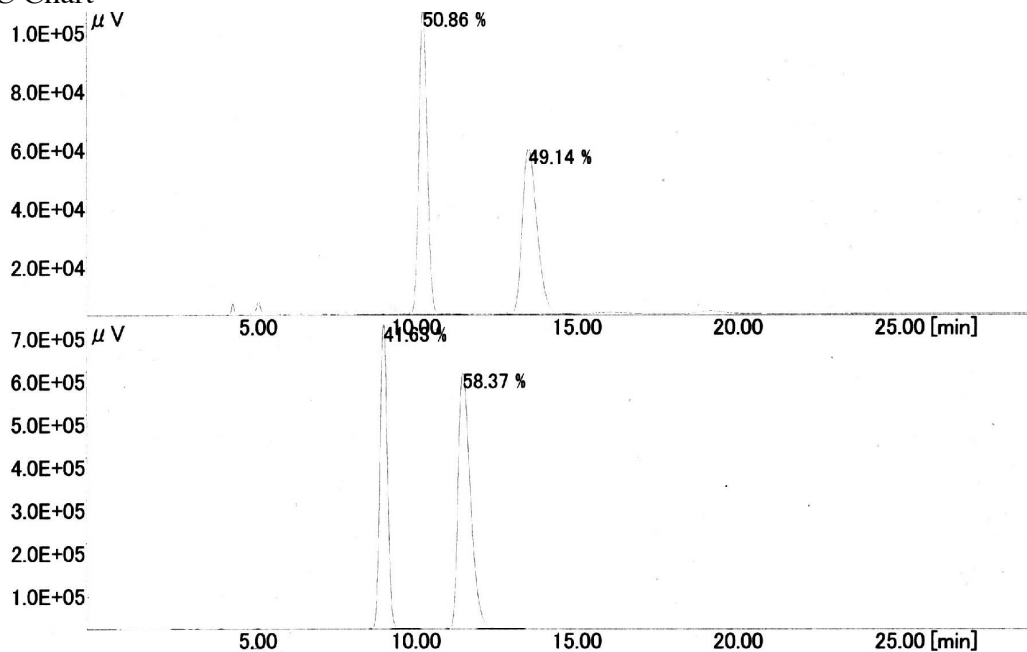
¹H NMR spectra



¹³C NMR spectra



HPLC Chart



ファイル名: 030_1777.CH1

コメント:

030

ad_3

3/1

1.0 ml/min

Vial # = 1 Rack # = 1

注入日 : 17-Dec-2013 19:03:26

現在日時: 27-Dec-2013 14:09:08

ユーザー : MIYAKE

グループ : MIYAKE

システムプログラム:

#	ピーク名	Rt	面積[μ V.Sec]	定量値
1		10.442	1832765.168	0.000
2		13.717	1770756.000	0.000

ピーク総面積= 3603521.168 [μ V.Sec]

ファイル名: 918594.CH1

コメント:

918

ad_3

H/iPr = 3/1

Flow 1.0mL/min

Vial # = 1 Rack # = 1

注入日 : 16-Nov-2012 16:14:04

現在日時: 27-Dec-2013 14:09:08

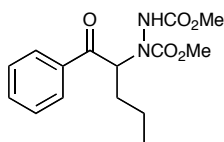
ユーザー : MIYAKE

グループ : MIYAKE

システムプログラム:

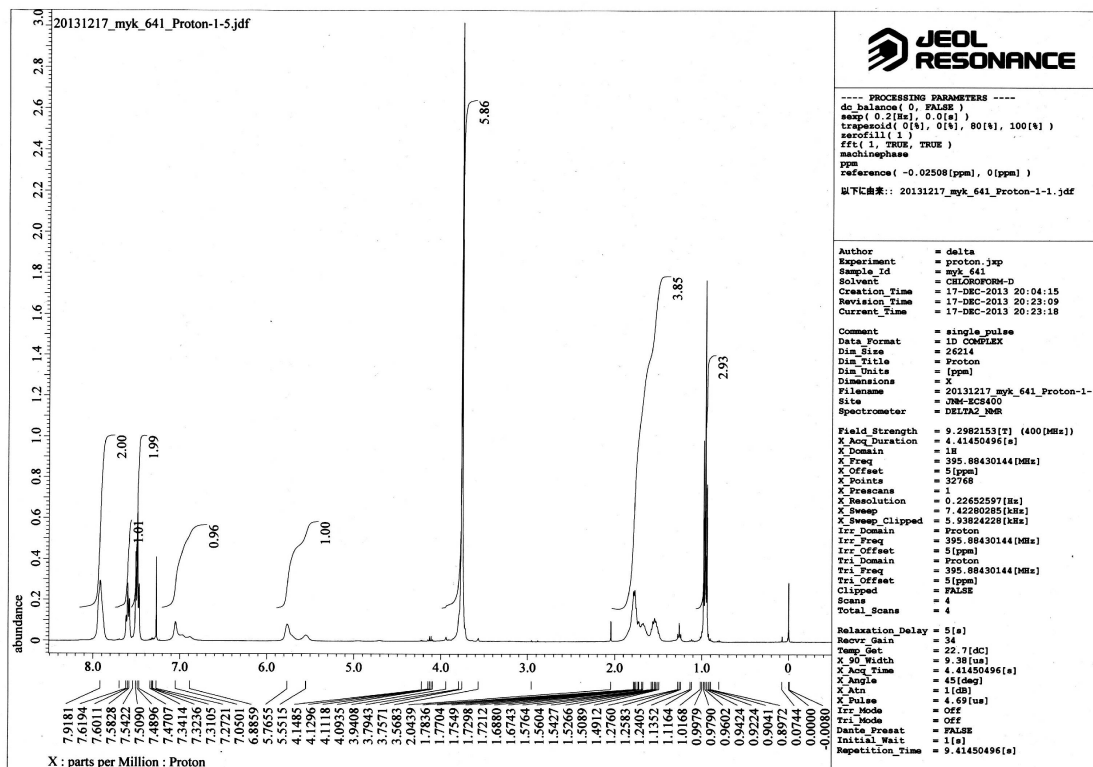
#	ピーク名	Rt	面積[μ V.Sec]	定量値
1		9.217	11070282.750	0.000
2		11.675	15522125.500	0.000

ピーク総面積= 26592408.250 [μ V.Sec]

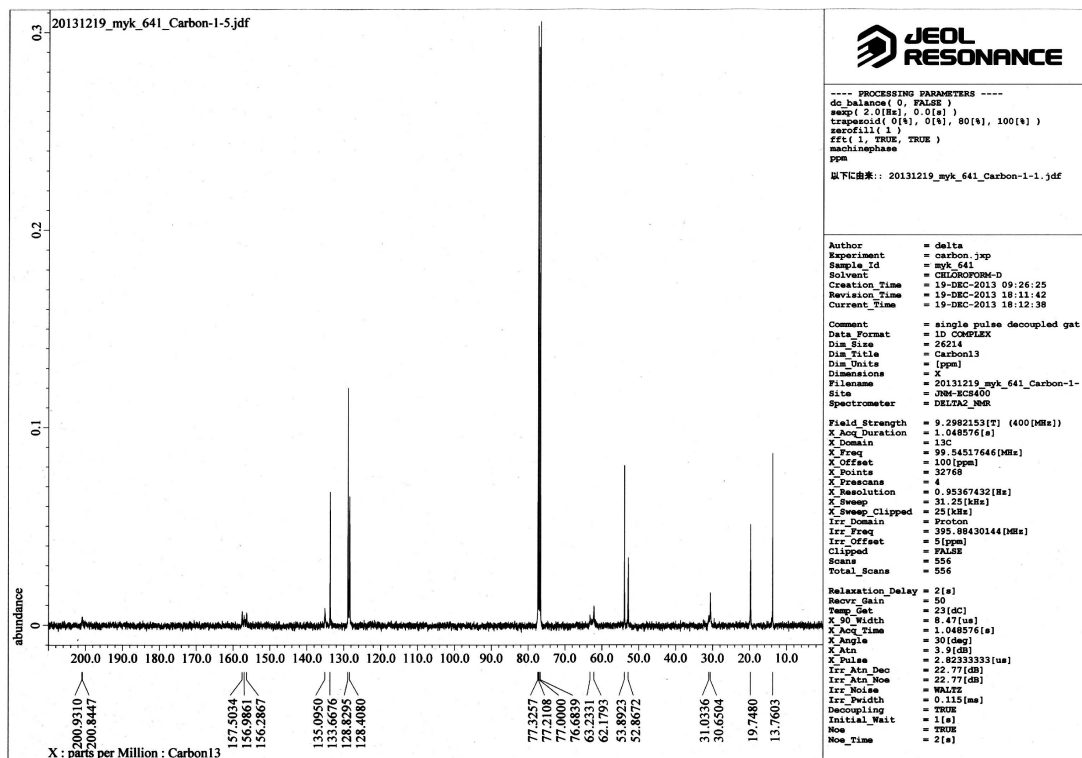


Dimethyl 1-(1-oxo-1-phenylpentan-2-yl)hydrazine-1,2-dicarboxylate (3ga,
Table 3, entry 9)

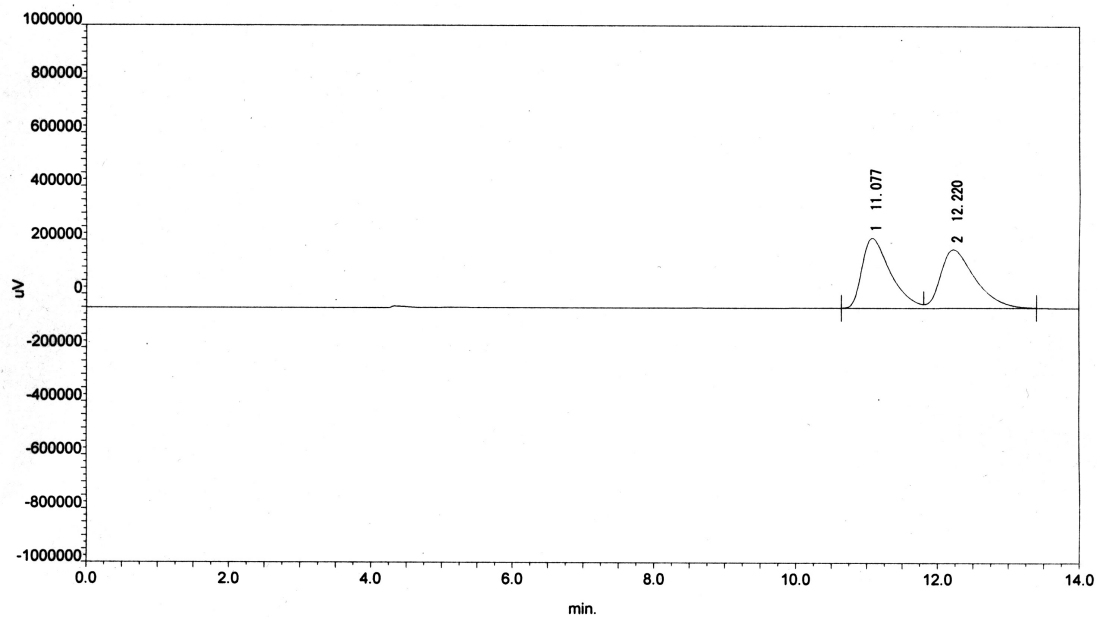
¹H NMR spectra



¹³C NMR spectra

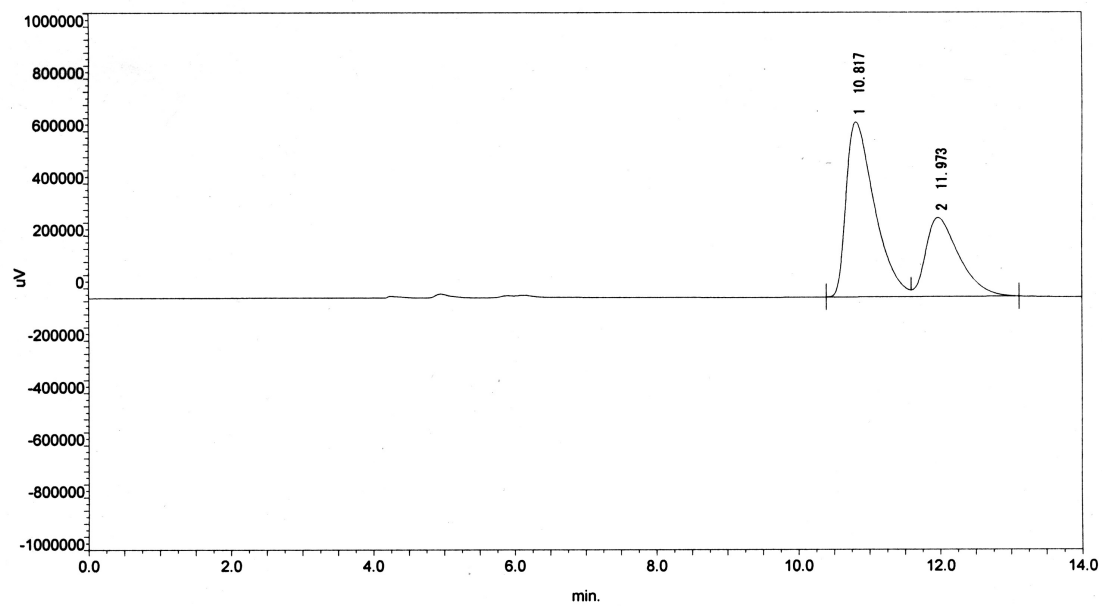


HPLC Chart



解析結果

No.	Rt(min)	ピーク名	面積	面積(%)	高さ	NTP	対称性	分離度
1	11.08		7269428.818	49.5760	259943	3548.6	-----	1.388
2	12.22		7393762.182	50.4240	217568	2924.7	-----	-----
			14663191.000	100.0000	477511			



解析結果

No.	Rt(min)	ピーク名	面積	面積(%)	高さ	NTP	対称性	分離度
1	10.82		18612502.949	65.4289	667640	3405.2	1.788	1.438
2	11.97		9834424.051	34.5711	301411	3046.0	-----	-----
			28446927.000	100.0000	969051			