CHEMISTRY
A EUROPEAN JOURNAL

Supporting Information

Polysulfurated pyrene-cored dendrimers: luminescent and electrochromic properties

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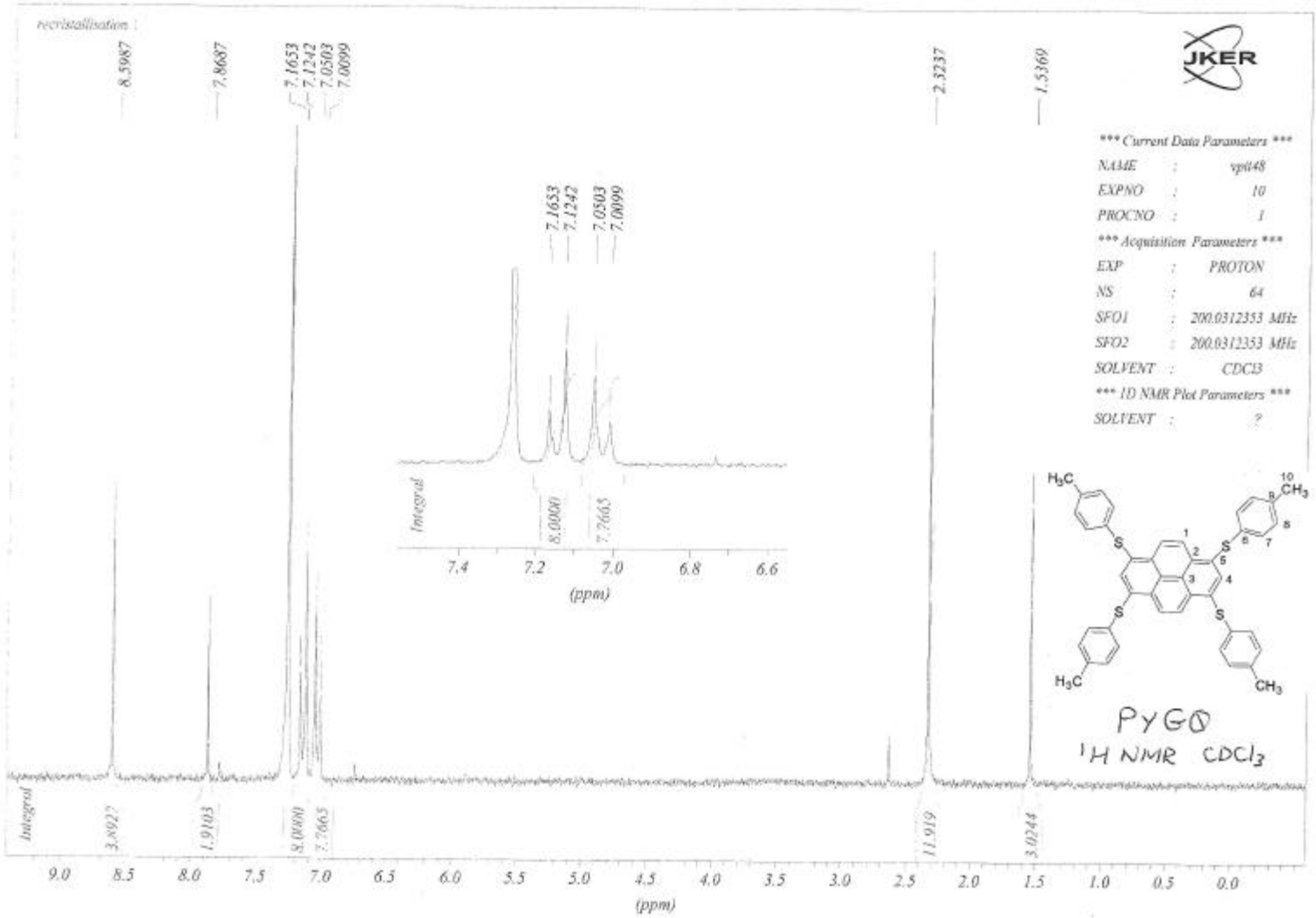
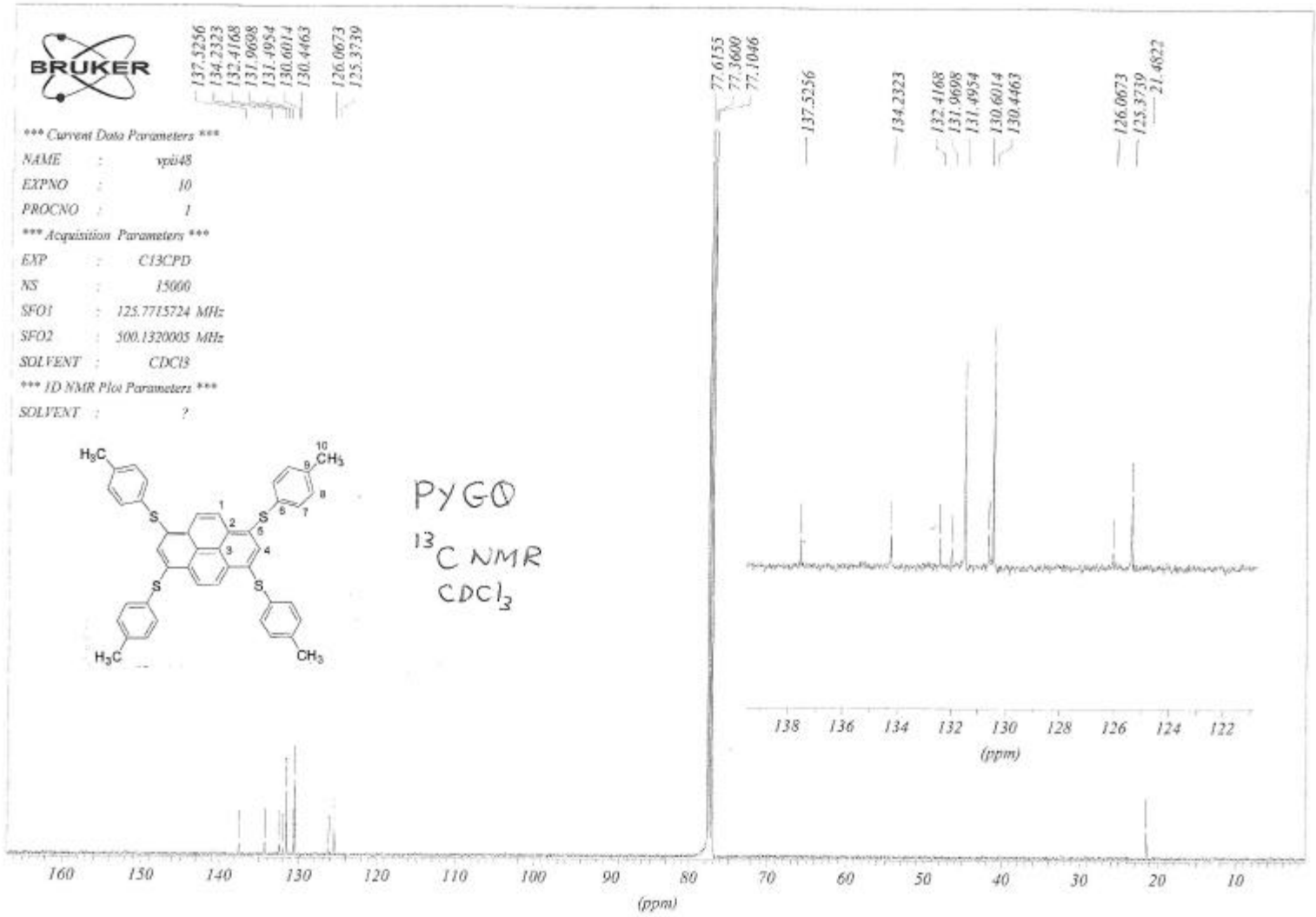
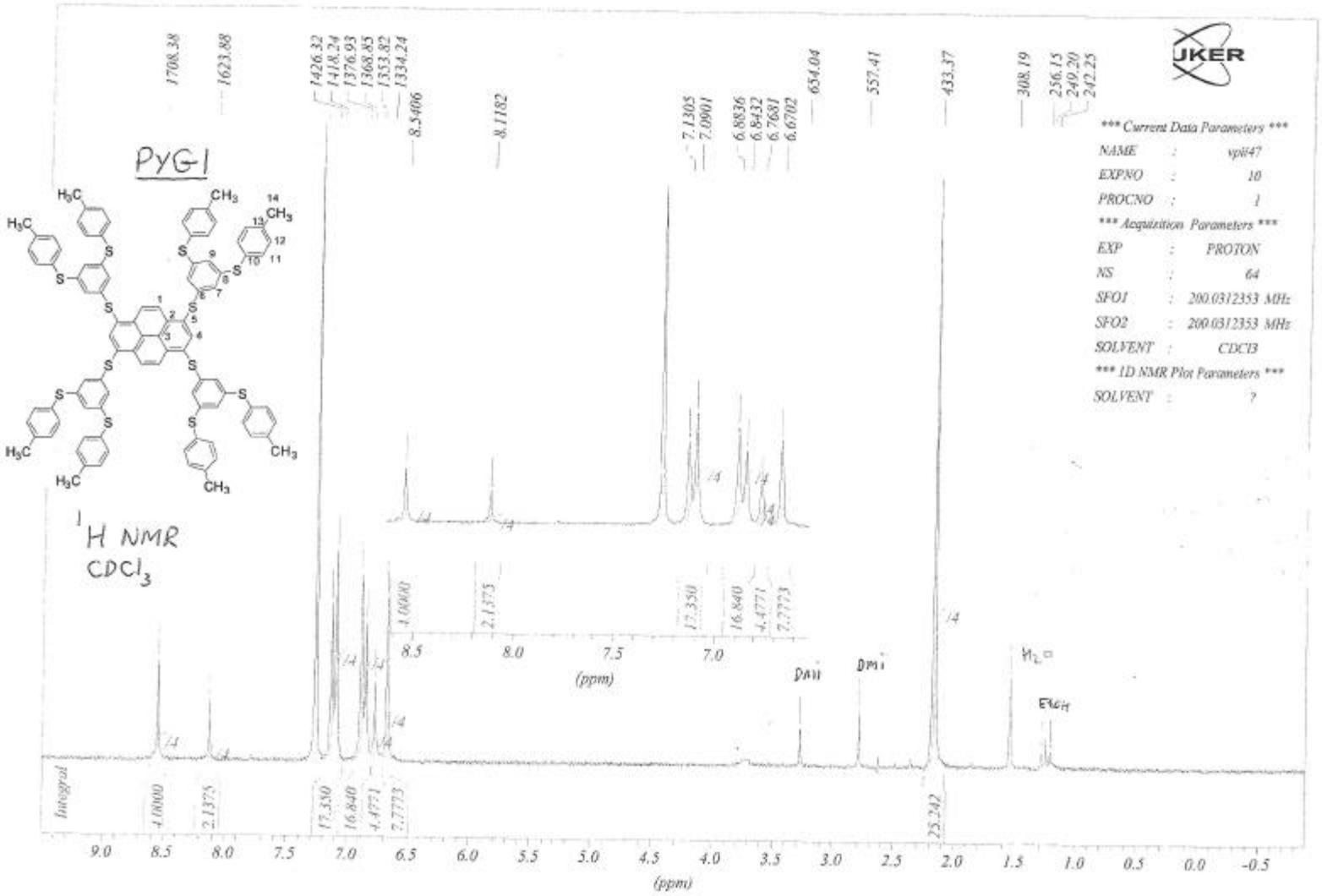
a) ^{1}H and ^{13}C NMR spectra of polysulfurated pyrene-cored dendrimers: (see below)

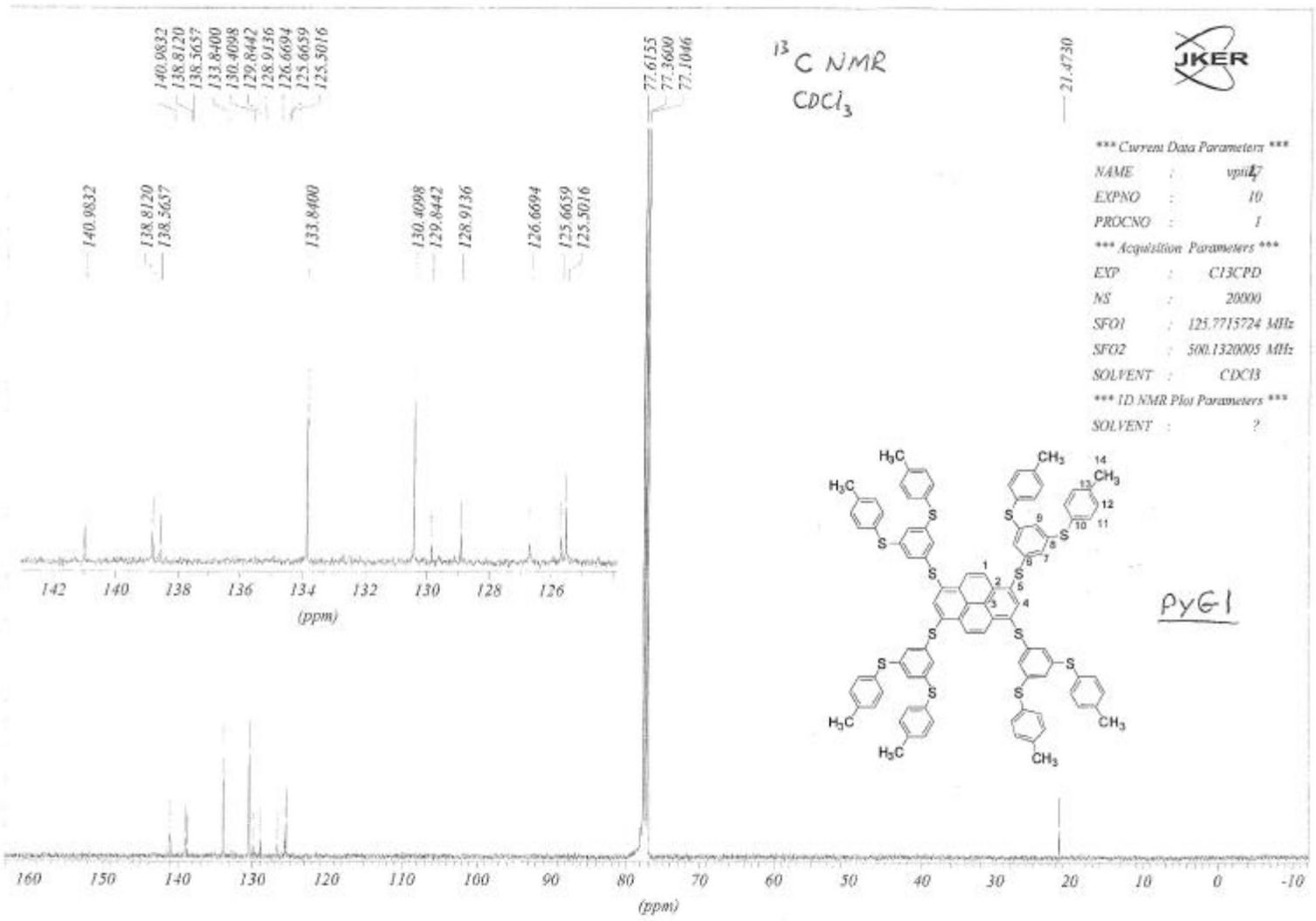
b) General procedure for ESI-MS by chemical ionisation with AgOTf

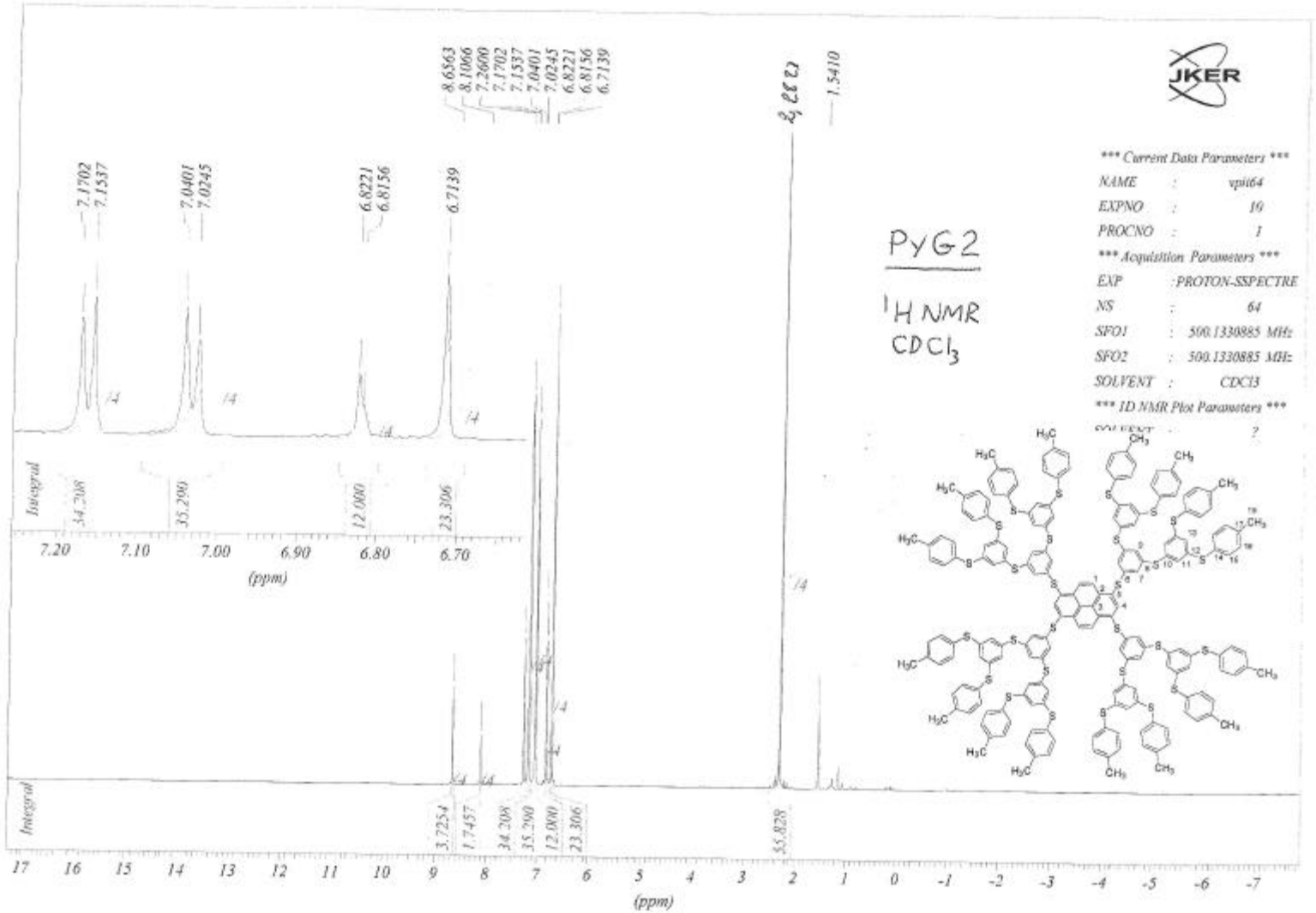
The pyrene sample (2 mg) was dissolved in dichloromethane (250 μL). This solution was diluted to 1/100 in dichloromethane. Another solution of AgOTf (1.7 mg) was prepared in acetonitrile (500 μL). The latter solution was diluted to 1/10 in acetonitrile. The "working solution" was prepared while mixing the pyrene solution (200 μL) to the AgOTf solution (200 μL). This solution was doped with double internal standards (poly(propylene glycol of known masses) and was introduced in the ionisation source (5500 V) at a flow rate of 5 μL/min.


recristallisation :
8.5987
7.8687
7.1653
7.1242
7.0503
7.0099
2.3237
1.5369
*** Current Data Parameters ***
NAME : vpt48
EXPNO : 10
PROCNO : 1
*** Acquisition Parameters ***
EXP : PROTON
NS : 64
SFO1 : 200.0312353 MHz
SFO2 : 200.0312353 MHz
SOLVENT : CDCl3
*** 1D NMR Plot Parameters ***
SOLVENT : ?
Integral
8.0000
7.7665
(ppm)
Integral
3.8927
1.9103
8.0000
7.7665
11.919
3.0244
(ppm)
PYG0
1H NMR CDCl3


BRUKER
137.5256
134.2323
132.4168
131.9698
131.4954
130.6014
130.4463
126.0673
125.3739
77.6155
77.3600
77.1046
21.4822
*** Current Data Parameters ***
NAME : vpi48
EXPNO : 10
PROCNO : 1
*** Acquisition Parameters ***
EXP : C13CPD
NS : 15000
SFO1 : 125.7715724 MHz
SFO2 : 500.1320005 MHz
SOLVENT : CDCl3
*** 1D NMR Plot Parameters ***
SOLVENT : ?
PyGO
13C NMR
CDCl3
(ppm)


PYG1
^{1}H NMR
$CDCl_3$
1708.38
1623.88
1426.32
1418.24
1376.93
1368.85
1353.82
1334.24
8.5406
8.1182
7.1305
7.0901
6.8836
6.8432
6.7681
6.6702
654.04
557.41
433.37
308.19
256.15
249.20
242.25
*** Current Data Parameters ***
NAME : vpii47
EXPNO : 10
PROCNO : 1
*** Acquisition Parameters ***
EXP : PROTON
NS : 64
SFO1 : 200.0312353 MHz
SFO2 : 200.0312353 MHz
SOLVENT : CDCl3
*** 1D NMR Plot Parameters ***
SOLVENT : ?
4.0000
2.1375
17.350
16.840
4.4771
7.7773
25.242
(ppm)
DMF
H_2O
EtOH


140.9832
138.8120
138.5657
133.8400
130.4098
129.8442
128.9136
126.6694
125.6659
125.5016
77.6155
77.3600
77.1046
21.4730
13C NMR
CDCl3
JKER
*** Current Data Parameters ***
NAME : vpiii7
EXPNO : 10
PROCNO : 1
*** Acquisition Parameters ***
EXP : C13CPD
NS : 20000
SFO1 : 125.7715724 MHz
SFO2 : 500.1320005 MHz
SOLVENT : CDCl3
*** 1D NMR Plot Parameters ***
SOLVENT : ?
PyG1
(ppm)
(ppm)


PyG2
^{1}H NMR
$CDCl_3$
JKER
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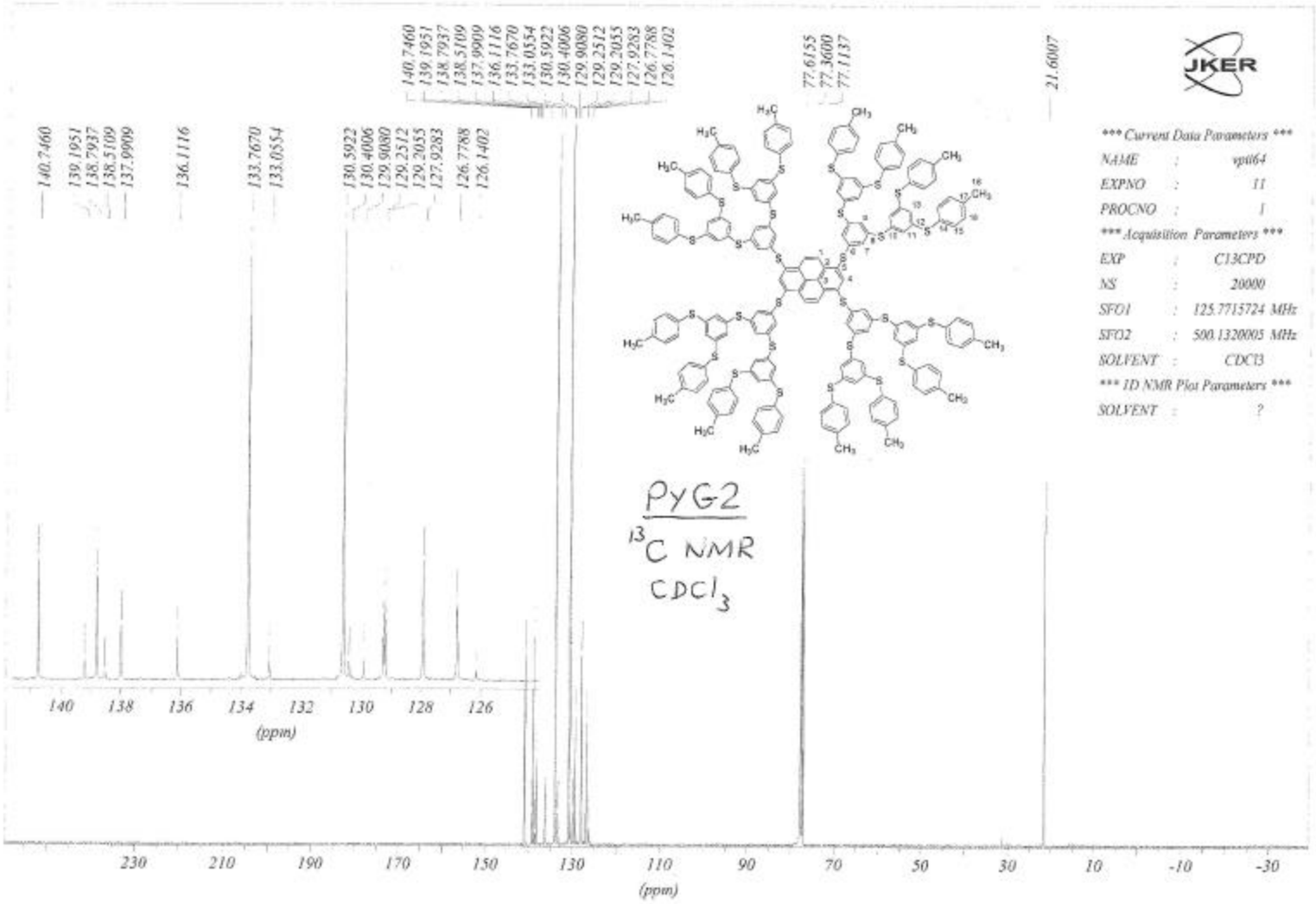
NAME : vpii64
EXPNO : 10
PROCNO : 1
*** Acquisition Parameters ***
EXP : PROTON-SSPECTRE
NS : 64
SFO1 : 500.1330885 MHz
SFO2 : 500.1330885 MHz
SOLVENT : CDCl3
*** 1D NMR Plot Parameters ***
8.6563
8.1066
7.2600
7.1702
7.1537
7.0401
7.0245
6.8221
6.8156
6.7139
2.2812
1.5410
Integral
3.7254
1.7457
34.208
35.290
12.000
23.306
55.828
(ppm)


140.7460
139.1951
138.7937
138.5109
137.9909
136.1116
133.7670
133.0554
130.5922
130.4006
129.9080
129.2512
129.2055
127.9283
126.7788
126.1402
77.6155
77.3600
77.1137
21.6007
JKER
*** Current Data Parameters ***
NAME : vpti64
EXPNO : 11
PROCNO : 1
*** Acquisition Parameters ***
EXP : C13CPD
NS : 20000
SFO1 : 125.7715724 MHz
SFO2 : 500.1320005 MHz
SOLVENT : CDCl3
*** 1D NMR Plot Parameters ***
SOLVENT : ?
PyG2
13C NMR
CDCl3
140 138 136 134 132 130 128 126
(ppm)
230 210 190 170 150 130 110 90 70 50 30 10 -10 -30
(ppm)