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EUROPEAN PROJECT RESPONSE 2 to challenges in forensic drugs analyses
Co-funded by the Internal Security Found of the European Union

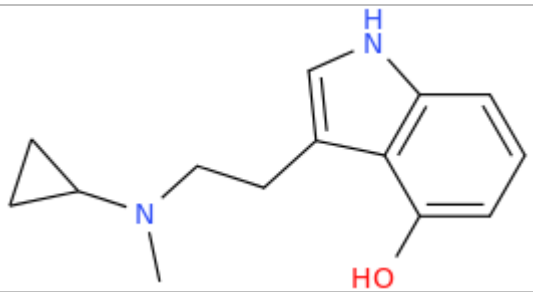
ANALYTICAL REPORT

4-HO-McPT (C14H18N2O)

3-{2-[cyclopropyl(methyl)amino]ethyl}-1H-indol-4-ol

Remark – other NPS detected: none

Sample ID:	1961-18
Sample description:	powder
Sample type:	test purchase /ISF projekt (NFL-SI)
Date of sample receipt (DD/MM/YYYY):	21/08/2018
Date of entry (DD/MM/YYYY) into NFL database:	10/09/2018
Report updates (if any) will be published here:	http://www.policija.si/apps/nfl_response_web/seznam.php

Substance identified - structure[1] (base form)	
Systematic name	3-{2-[cyclopropyl(methyl)amino]ethyl}-1H-indol-4-ol
Other names	
Formula (per base form)	C14H18N2O
M_w (g/mol)	230,31
Salt form/anions detected	fumarate
StdInChIKey (per base form)	GFVJBFIXZYLVPO-UHFFFAOYSA-N
Other NPS detected	none
Additional info (purity..)	> 93 % based on H1 NMR

[1] Created by OPSIN free tool: http://opsin.ch.cam.ac.uk/ **DOI:** 10.1021/ci100384d

Report updates

date	comments (explanation)

Instrumental methods (if applied) in NFL

1. GC-MS (Agilent): GC-method is RT locked to tetracosane (9.258 min). Injection volume 1 ml and split mode (1:50). Injector temperature: 280 0C. Chromatographic separation: on column HP1-MS (100% dimethylpolysiloxane), length 30 m, internal diameter 0.25 mm, film thickens 0.25 µm. Carrier gas He: flow-rate 1.2 ml/min. GC oven program: 170 ^{0}C for 1 min, followed by heating up to 190 ^{0}C at rate 8 ^{0}C/min, then heating up to 293 0C at a rate of 18 ^{0}C/min, hold for 7.1 min, then heating at 50 ^{0}C/min up to 325 ^{0}C and finally 6.1 min isothermal. MSD source EI = 70 eV. GC-MS transfer line T= 235^{0}C, source and quadropole temperatures 280^{0}C and 180^{0}C, respectively. Scan range m/z scan range: from 50 (30 until 6 min.) to 550 (300 until 6 min) amu.

2. HPLC-TOF (Agilent): 6230B TOF with Agilent 1260 Infinity HPLC with binary pump, column: Zorbax Eclipse XDB-C18, 50 x 4.6 mm, 1.8 micron. Mobile phases (A) 0.1% formic acid and 1mM ammonium formate in water; (B) 0.1% formic acid in methanol (B). Gradient: starting at 5% B, changing to 40% B over 4 min, then to 70% over 2 min and in 5 min to 100%, hold 1 min and back to 5%, equilibration for 1.7 min. The flow rate: 1.0 ml/min; Injection volume 1 µl. MS parameters: 2GHz, Extended Dynamic range mode to a maximum of 1700 amu, acquisition rate 1.30 spectra/sec. Sample ionisation: by Agilent Jet Stream technology (Dual AJS ESI). Ion source: positive ion scan mode with mass scanning from 82 to 1000 amu. Other TOF parameters: drying gas (N2) and sheath temperature 325 °C; drying gas flow rate 6 l/min; sheath gas flow rate 8 l/min; nebulizer 25 psig; Vcap. 4000 V; nozzle 2000 V; skimmer 65 V; fragmentor 175 V and Octopole RF 750 V.

3.FTIR-ATR (Perkin Elmer): scan range 4000-400 cm-1; resolution 4cm-1

4. GC- (MS)-IR condensed phase (GC-MS (Agilent) & IR (Spectra analyses-Danny)
GC-method: Injection volume 1 ml and split mode (1:5). Injector temperature 280 ^{0}C. Chromatographic separation as above **(1).**
Split MS : IR = 1: 9.
MSD source EI = 70 eV. GC-MS transfer line T= 235^{0}C, source and quadropole temperatures 280^{0}C and 180^{0}C, respectively. Scan range m/z scan range: from 50 (30 until 6 min.) to 550 (300) amu.
IR (condesed (solid) phase): IR scan range 4000 to 650, resolution 4 cm^{-1} .

5. **IC** (anions) (Thermo Scientific, Dionex ICS 2100), Column: IonPac AS19, 2 x 250mm; Eluent: 10mM KOH from 0 to 10 min, 10-58 mM from 10 to 40min; Flow rate: 0.25 ml/min; Temperature: 30˚C; Suppressor: AERS 500 2mm, suppressor current 13mA; Inj. Volume: 25 µl

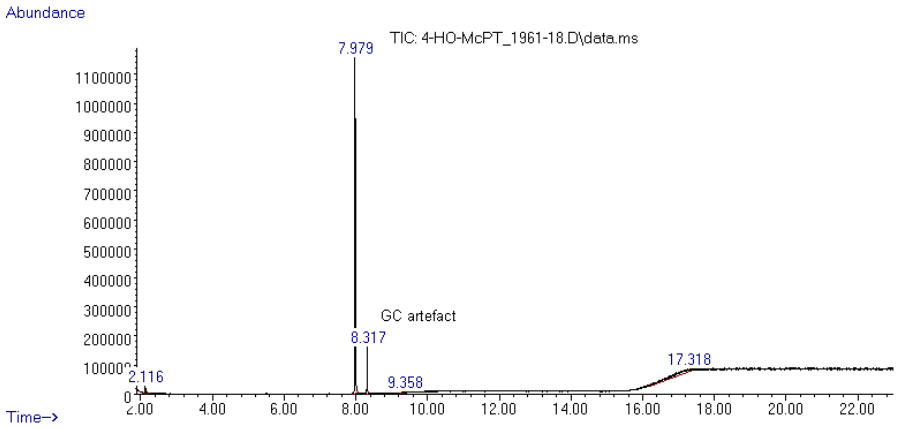
Supporting information

Solubility in	result/remark
CH_2Cl_2	partially
MeOH	partially
H_2O	/

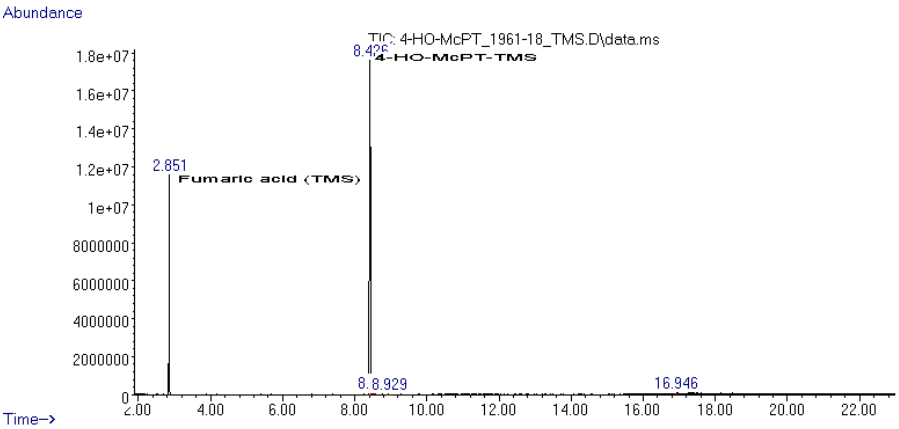
Analytical technique:	applied	remarks
GC-MS (EI ionization)	+	NFL GC-RT (min): 7,98 BP(1): 84; BP(2): 159,BP(3) :230,
HPLC-TOF	+	Exact mass (theoretical): 230,1419; measured value Δppm:-0,3; formula:C14H18N2O
FTIR-ATR	+	direct measurement (sample as received)
FTIR (solid phase) always as base form	+	
IC (anions)	+	
NMR (in FKKT)	+	
validation		
other		

ANALYTICAL RESULTS

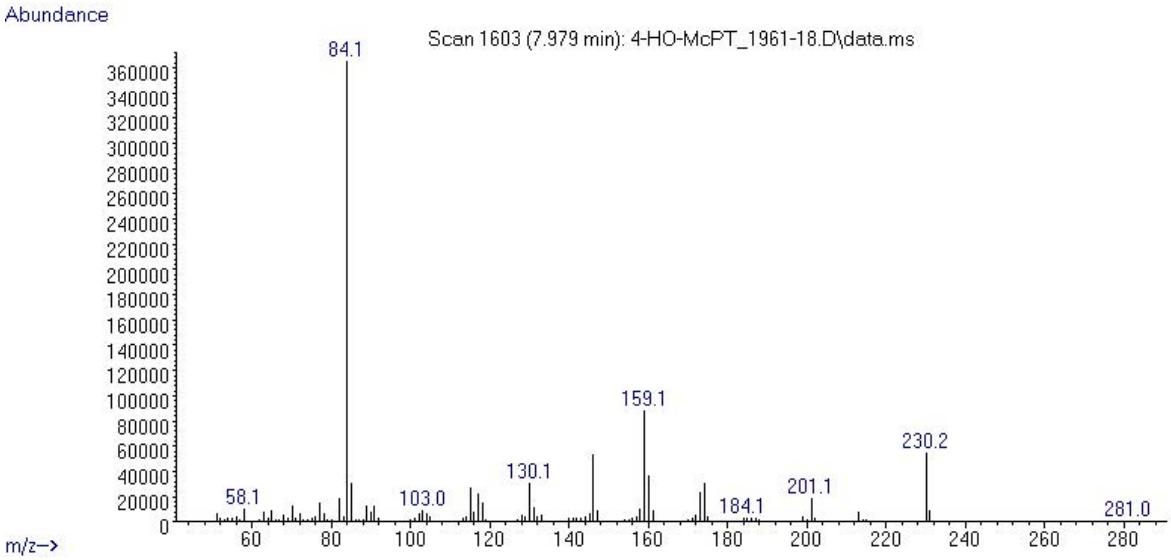
GC-MS chromatogram (non derivatized sample)



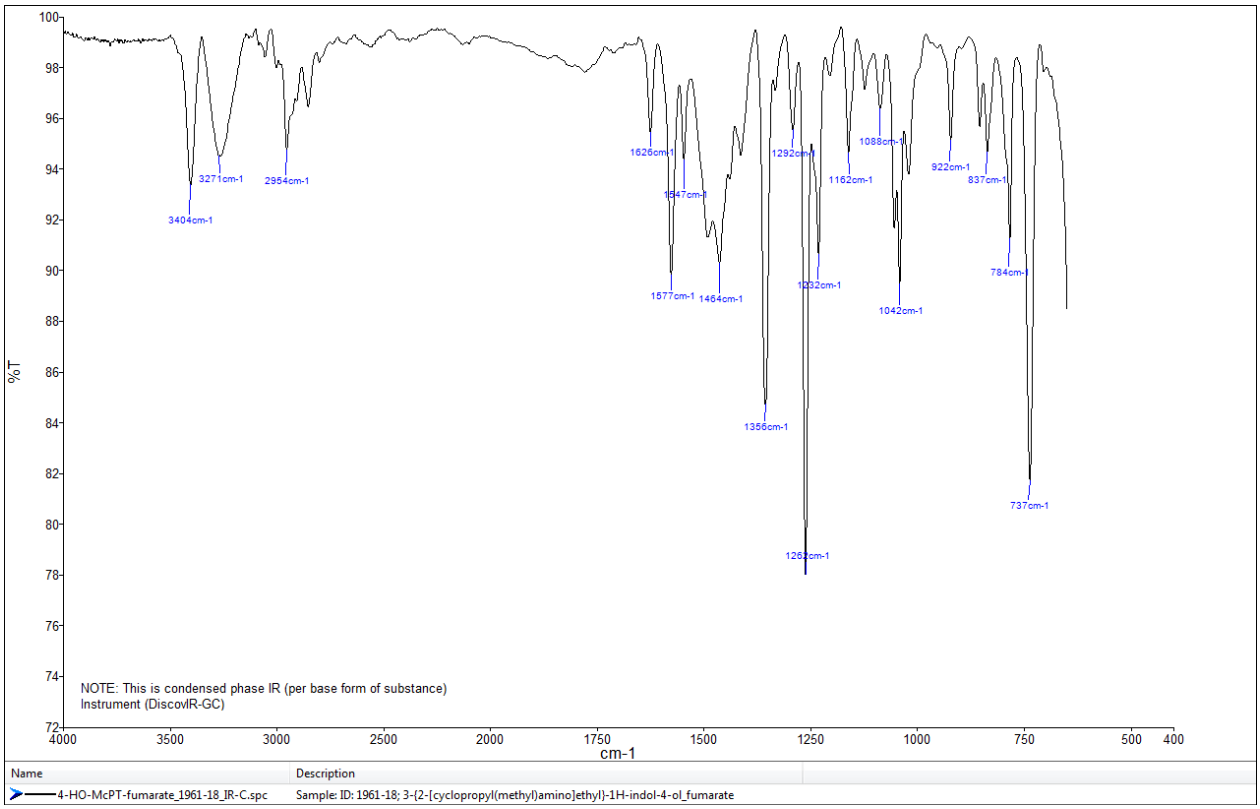
GC-MS chromatogram – TMS derivative



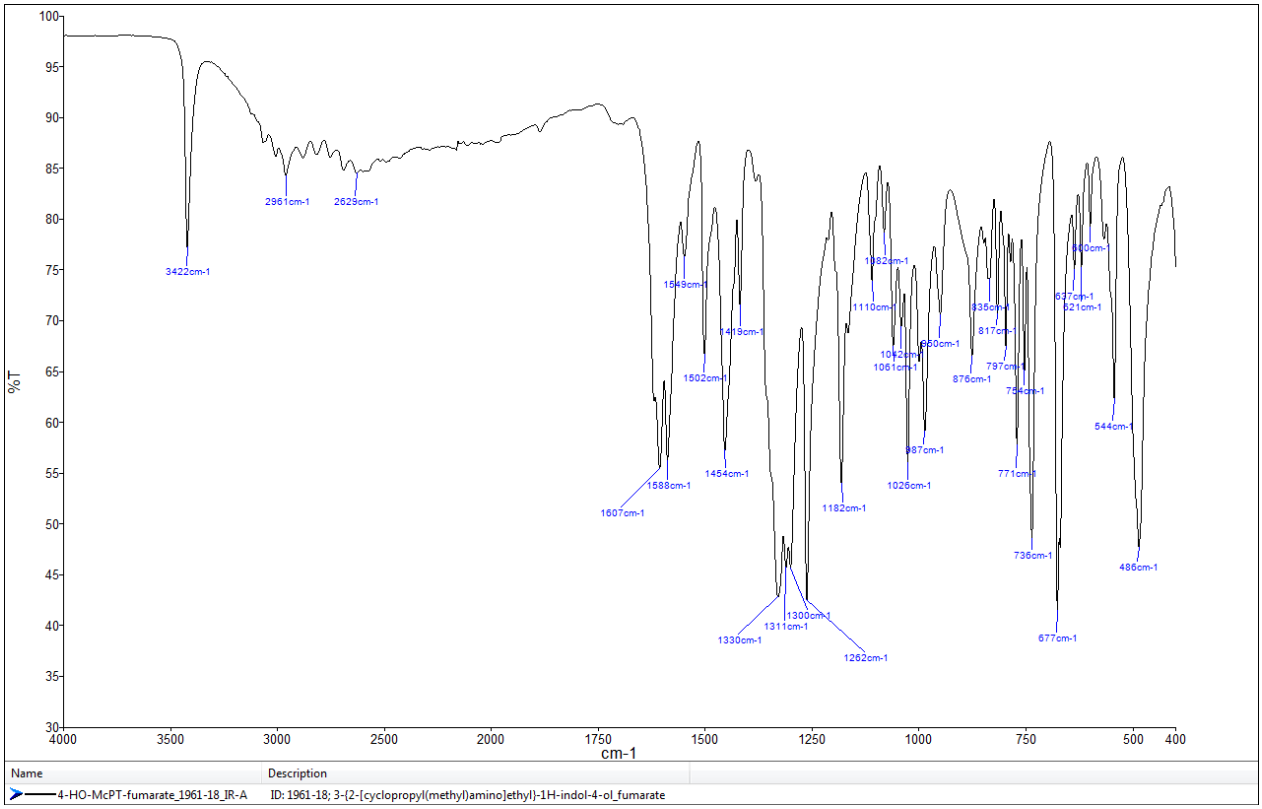
MS (EI)



FTIR-ATR - direct measurement (sample as received fumarate salt)



IR (solid phase – after chromatographic separation)



TOF REPORT

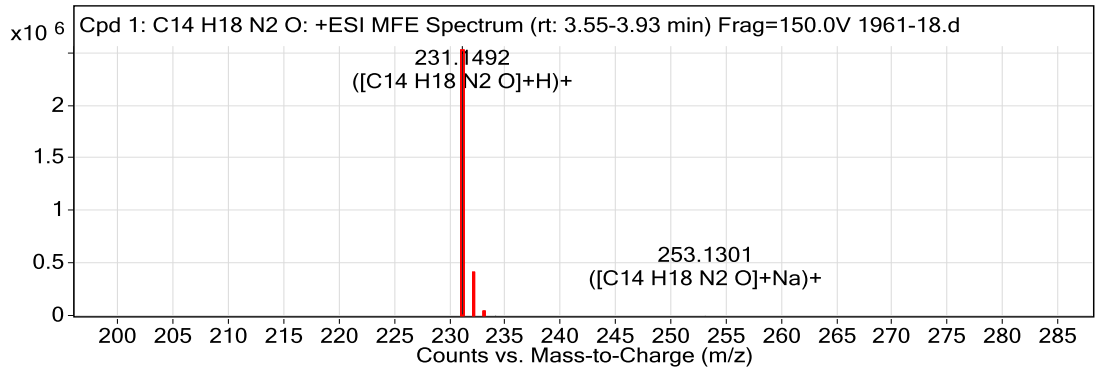
Data File	1961-18.d	**Sample Name**	ID 1961-18
Sample Type	Sample	**Position**	P1-C7
Instrument Name	6230B TOF LC-MS	**User Name**	TG
Acq Method	general-04_12_2017-XDB-C18-ESI+.m	**Acquired Time**	8/22/2018 4:23:52 PM
IRM Calibration Status	Success	**DA Method**	a-Drugs_NFL.m
Comment	MeOH		

Compound Table

Label	MFG Formula	Obs. RT	Obs. Mass
Cpd 1: C14 H18 N2 O	C14 H18 N2 O	3.6	230.142

Obs. m/z	Obs. RT	Obs. Mass
231.1492	3.6	230.142

MFE MS Zoomed Spectrum



MS Spectrum Peak List

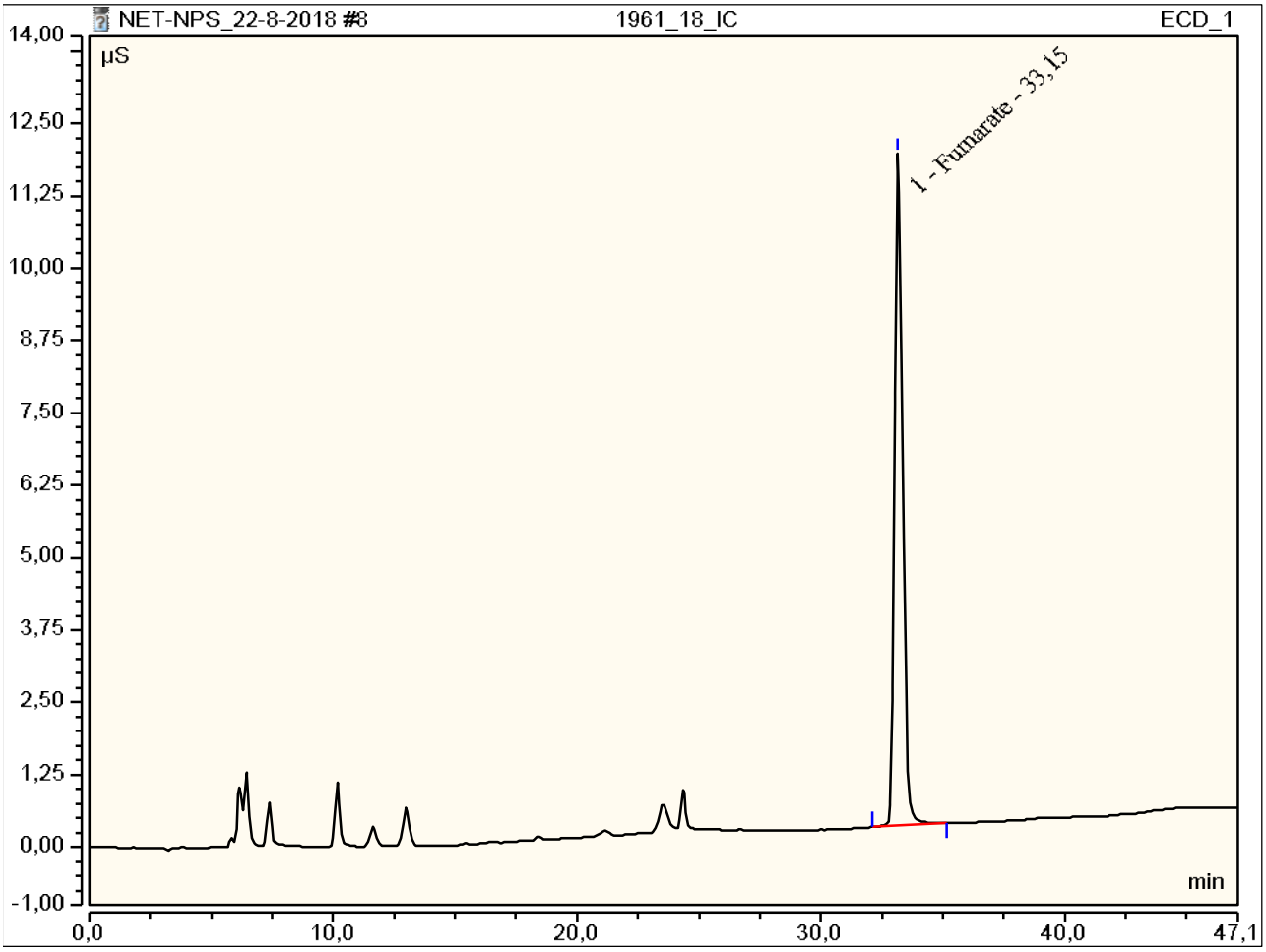
Obs. m/z	Charge	Abund	Formula	Ion/Isotope
231.1492	1	2563623.5	C14 H18 N2 O	(M+H)+
232.1526	1	370515.27	C14 H18 N2 O	(M+H)+
233.155	1	32647	C14 H18 N2 O	(M+H)+
234.1577	1	2186.53	C14 H18 N2 O	(M+H)+
253.1301	1	1010.84	C14 H18 N2 O	(M+Na)+

--- End Of Report ---

Peak Integration Report

Sample Name:	**1961_18_IC**	*Inj. Vol.:*	**25,00**
Injection Type:	**Unknown**	*Dilution Factor:*	**1,0000**
Program:	**ANIONI - julij 2018**	*Operator:*	**kemija**
Inj. Date / Time:	**23-avg-2018 / 09:00**	*Run Time:*	**47,10**

No.	Time min	Peak Name	Peak Type	Area µS*min	Height µS	Amount mg/L
1,00	33,15	**Fumarate**	BMB	4,49	11,62	n.a.
		TOTAL:		4,49	11,62	0,00



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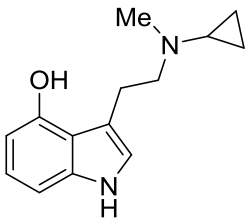
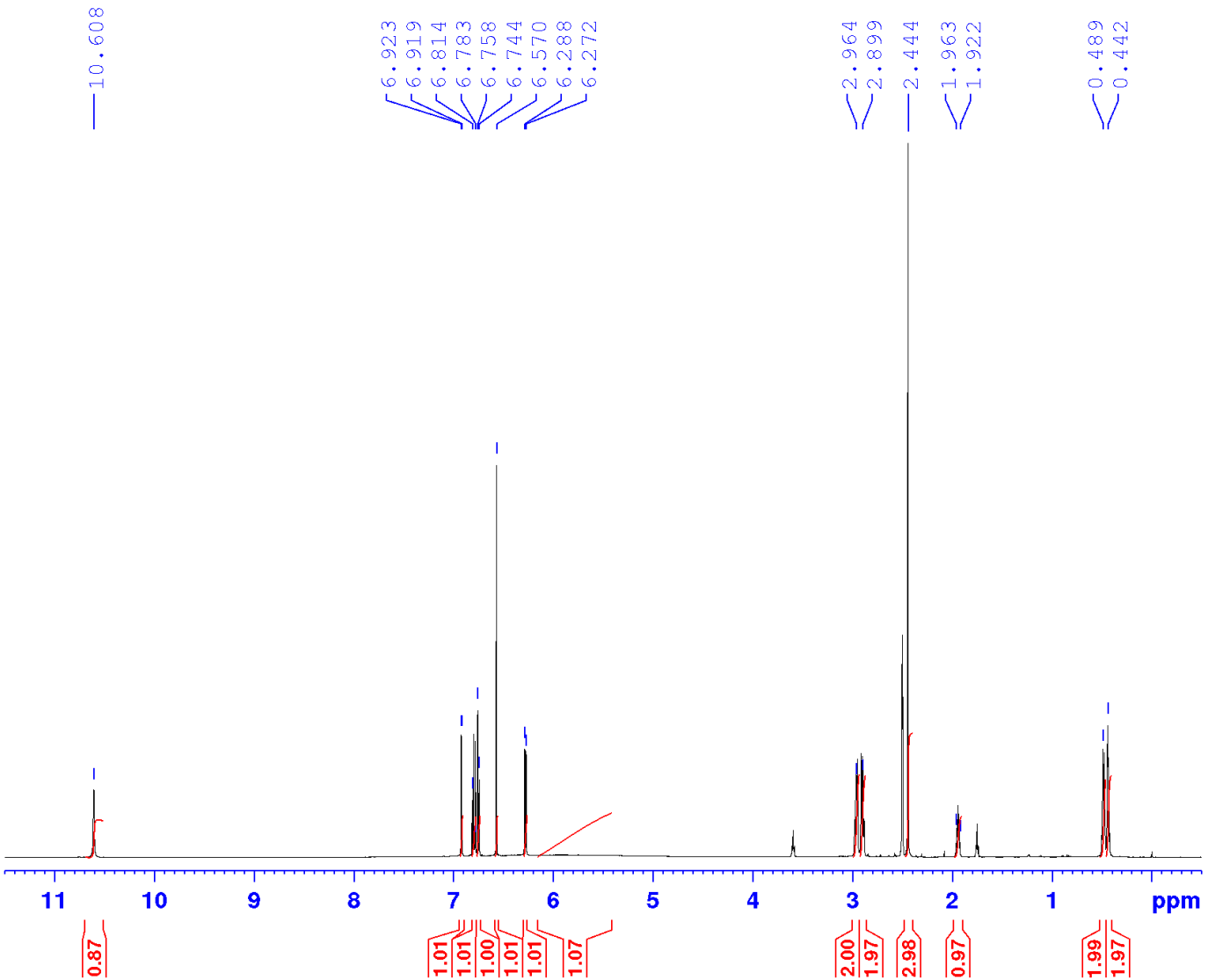
R E P O R T

Contract No.	C1714-17-460078 (Republic of Slovenia, Ministry of the Interior, POLICE)
Sample ID:	**1961-18**
Received date:	August 29, 2018
Our notebook code:	NFL-1961-18
NMR sample preparation:	22.2 mg dissolved in 0.7 mL DMSO-d_6
NMR experiments:	^{1}H, ^{13}C, ^{1}H–^{1}H *gs*-COSY, ^{1}H–^{13}C *gs*-HSQC, ^{1}H–^{13}C *gs*-HMBC, ^{1}H–^{15}N *gs*-HMBC
Proposed structure with formula, exact mass, molecular weight:	 Chemical Formula: $C_{14}H_{18}N_2O$ Exact Mass: 230,1419 Molecular Weight: 230,3110
Chemical name:	3-(2-(cyclopropyl(methyl)amino)ethyl)-1*H*-indol-4-ol
Comments:	- Structure elucidation based on 1D and 2D NMR spectra and HRMS. - The sample consists of herein identified compound and fumarate in molar ratio of 1 : 0.5, based on ^{1}H NMR spectrum. - >93% purity of a sample based on ^{1}H NMR spectrum
Supporting information:	Copies of ^{1}H and ^{13}C NMR spectra, ^{1}H and ^{13}C FIDs.
Principal investigator:	Prof. Dr. Janez Košmrlj
Date of report:	September 7, 2018

NFL-1961-18
1H

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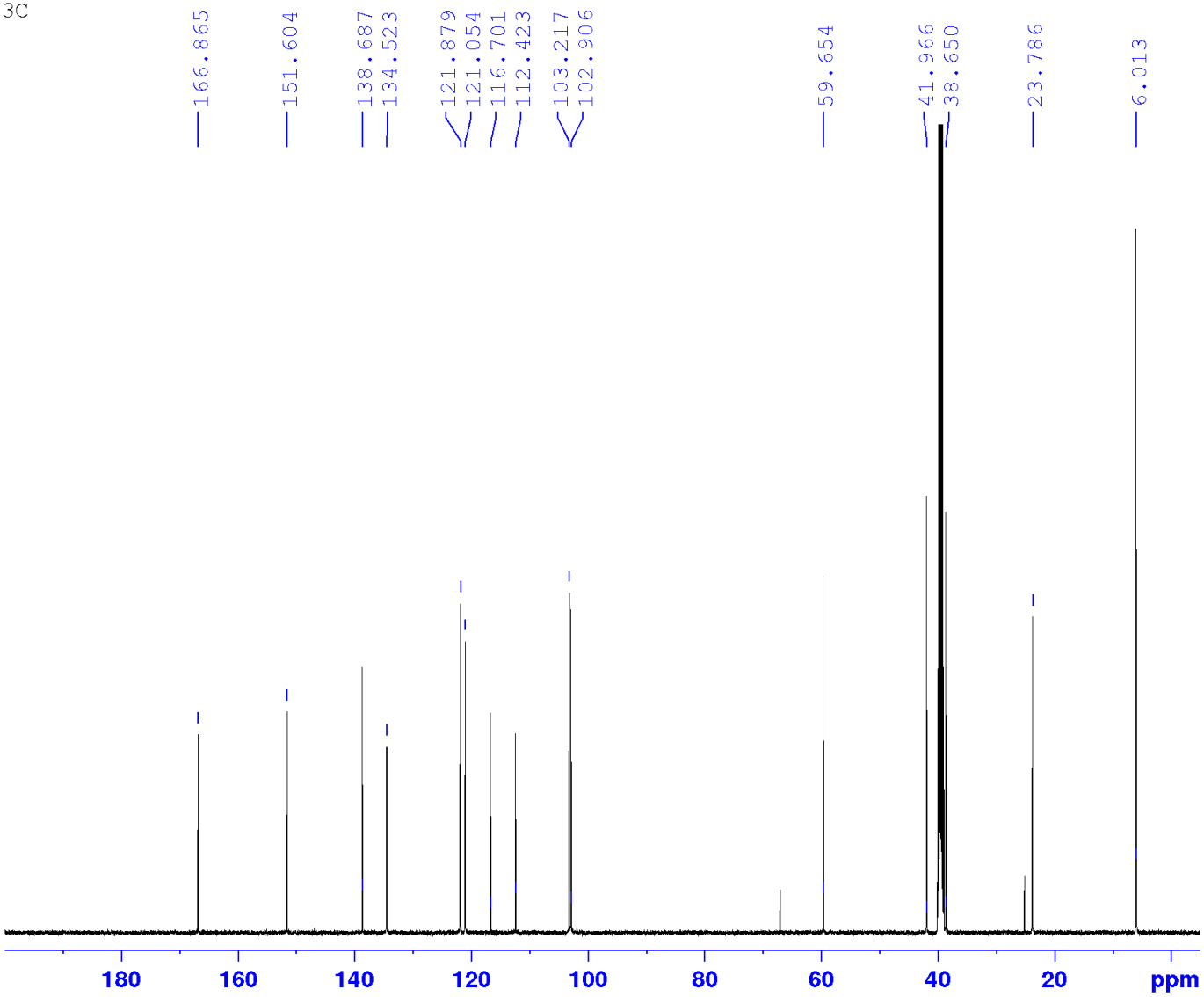
Current Data Parameters
NAME          NFL-1961-18
EXPNO                  10
PROCNO                  1

F2 - Acquisition Parameters
Date_            20180830
Time                7.26
INSTRUM            spect
PROBHD   5 mm PABBO BB-
PULPROG             zg30
TD                 65536
SOLVENT             DMSO
NS                    16
DS                     2
SWH           10000.000 Hz
FIDRES         0.152588 Hz
AQ            3.2767999 sec
RG                    64
DW               50.000 usec
DE                6.50 usec
TE                296.0 K
D1           1.00000000 sec
TD0                    1

======== CHANNEL f1 ========
SFO1         500.1330885 MHz
NUC1                   1H
P1                 8.70 usec
PLW1        26.00000000 W

F2 - Processing parameters
SI                 65536
SF           500.1300041 MHz
WDW                   EM
SSB       0
LB                0.30 Hz
GB        0
PC                1.00
```

NFL-1961-18
13C




```
Current Data Parameters
NAME         NFL-1961-18
EXPNO                  3
PROCNO                 1

F2 - Acquisition Parameters
Date_           20180829
Time              23.07
INSTRUM           spect
PROBHD   5 mm PABBO BB-
PULPROG         zgpg30
TD               65536
SOLVENT            DMSO
NS                 3072
DS                    4
SWH           29761.904 Hz
FIDRES         0.454131 Hz
AQ            1.1010048 sec
RG                 2050
DW               16.800 usec
DE                6.50 usec
TE               296.0 K
D1           1.00000000 sec
D11          0.03000000 sec
TD0                   1

======== CHANNEL f1 ========
SFO1        125.7703637 MHz
NUC1                13C
P1                 8.70 usec
PLW1       122.00000000 W

======== CHANNEL f2 ========
SFO2        500.1320005 MHz
NUC2                1H
CPDPRG[2         waltz16
PCPD2             80.00 usec
PLW2         26.00000000 W
PLW12         0.30046001 W
PLW13         0.15113001 W

F2 - Processing parameters
SI                 32768
SF          125.7578424 MHz
WDW                  EM
SSB      0
LB                1.00 Hz
GB       0
PC                1.40
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