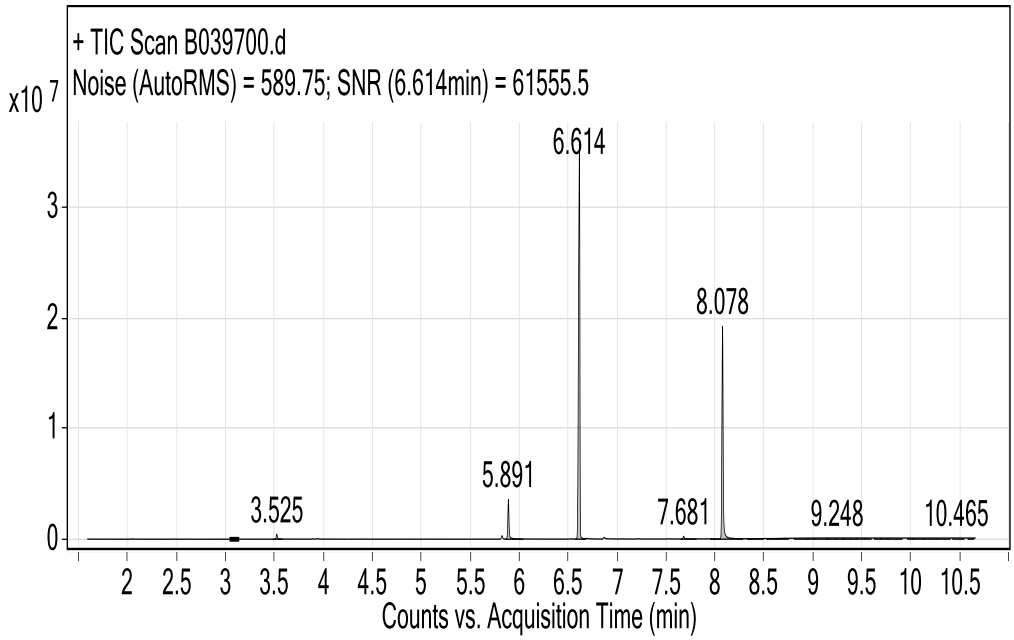
Politie, eenheid Amsterdam
DRR Laboratorium Forensische Opsporing
Instrument: **GC-MS 2 / AI156**

Sample Name **B039700**
Sample ID

Data File D:\data\overige2021\rk\20200201rk_cocacase28x\B039700.D
Inst Method algemeen.M
Proc Method algemeen.M
Acquisition Date 2/1/2021 6:27:12 PM
Operator RK
Comment RK project Stein

Sample Name B039700 **ID of Sample**
Inweeg [g] 0.0220 **Sample Type**
Vial Number 60 **Inj Vol [µl]** 1.00



Chromatogram Peak List

RT [min.]	Peak Area	Area [%]	AreaΣ [%]	S/N	FWHM
3.53	430,238	1.24	0.62	742.60	0.016
5.89	3,374,426	9.71	4.90	6156.40	0.015
6.61	34,740,194	100.00	50.44	61555.50	0.016
7.68	469,536	1.35	0.68	426.10	0.026
8.08	18,300,831	52.68	26.57	32716.40	0.015
9.20	555,419	1.60	0.81	187.20	0.060
9.25	469,149	1.35	0.68	190.30	0.051
9.41	409,050	1.18	0.59	185.30	0.048
9.75	458,312	1.32	0.67	185.50	0.055
10.47	723,988	2.08	1.05	188.80	0.079

Compound Table

#	RT [min.]	Mass	Name	CAS#	Match
1	3.525	181.10	2-Carbomethoxy-8-methyl-8-azabicyclo[3.2.1]oct-2-ene	43021-26-7	***97.49***
2	5.891	204.10	Levamisole	14769-73-4	***99.29***
3	6.616	303.10	Cocaine	50-36-2	***99.50***
4	8.078	394.50	Octacosane	630-02-4	***98.49***

RT [min.]	Name	Formula	Match	Mass	Library
3.53	2-Carbomethoxy-8-methyl-8-azabicyclo[3.2.1]oct-2-ene	C10H15NO2	***97.49***	181.10	NIST11.L
3.53	Anhydroecgonine methyl ester	C10H15NO2	***97.49***	181.00	SWGDRUG.L
3.53	Anhydroecgonine methyl ester	C10H15NO2	***97.49***	181.00	SWGDRUG_112916.L
3.53	Ecgonidine, methyl ester	C10H15NO2	***90.35***	181.10	NIST11.L
3.53	2-Carbomethoxy-8-methyl-8-azabicyclo[3.2.1]oct-2-ene	C10H15NO2	84.93	181.10	NIST11.L

RT [min.]	Name	Formula	Match	Mass	Library
5.89	Levamisole	C11H12N2S	***99.29***	204.10	NIST11.L
5.89	Levamisole	C11H12N2S	***98.85***	204.10	NIST11.L
5.89	(-)-Tetramisole	C11H12N2S	***98.84***	204.10	NIST11.L
5.89	Tetramisole	C11H12N2S	***98.84***	204.00	SWGDRUG.L
5.89	Tetramisole	C11H12N2S	***98.84***	204.00	SWGDRUG_112916.L

RT [min.]	Name	Formula	Match	Mass	Library
6.62	Cocaine	C17H21NO4	***99.50***	303.10	NIST11.L
6.62	Cocaine [PL940]	C17H21NO4	***99.04***	303.20	REFSPEC.L
6.62	Cocaine, Ph.Eur		***98.62***		POLLAB.L
6.62	Cocaine	C17H21NO4	***98.43***	303.10	NIST11.L
6.62	Cocaine	C17H21NO4	***98.23***	303.10	NIST11.L

RT [min.]	Name	Formula	Match	Mass	Library
8.08	Octacosane	C28H58	***98.49***	394.50	NIST11.L
8.08	n-Octacosane [PL1090]	C28H58	***97.53***	394.50	REFSPEC.L
8.08	Octacosane	C28H58	***95.42***	394.50	NIST11.L
8.08	Hentriacontane	C31H64	***94.98***	436.50	NIST11.L
8.08	Tetracosaan	C24H50	***94.41***	338.00	POLLAB.L

--- End Of Report ---