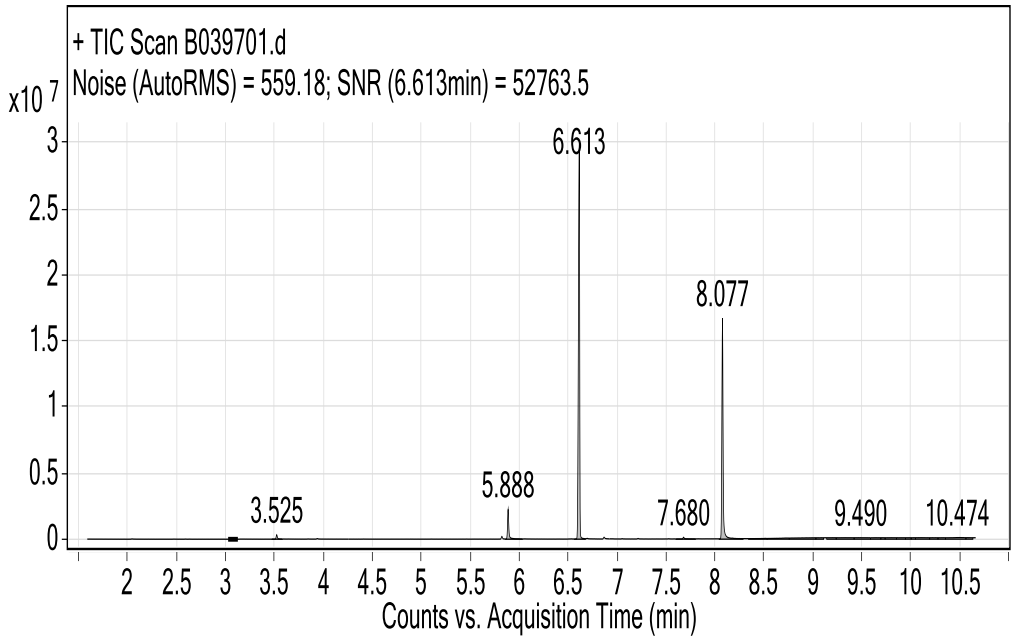
Politie, eenheid Amsterdam
DRR Laboratorium Forensische Opsporing
Instrument: **GC-MS 2 / AI156**

Sample Name **B039701**
Sample ID

Data File D:\data\overige2021\rk\20200201rk_cocacase28x\B039701.D
Inst Method algemeen.M
Proc Method algemeen.M
Acquisition Date 2/1/2021 6:43:51 PM
Operator RK
Comment RK project Stein
Sample Name B039701 **ID of Sample**
Inweeg [g] 0.0220 **Sample Type**
Vial Number 61 **Inj Vol [µl]** 1.00



Chromatogram Peak List

RT [min.]	Peak Area	Area [%]	AreaΣ [%]	S/N	FWHM
3.53	335,213	1.22	0.57	608.60	0.015
5.89	2,217,261	8.09	3.80	4045.10	0.015
6.61	27,392,990	100.00	46.97	52763.50	0.015
7.68	368,043	1.34	0.63	249.30	0.035
8.08	15,299,078	55.85	26.23	29954.00	0.015
9.17	607,177	2.22	1.04	195.90	0.092
9.49	1,040,226	3.80	1.78	196.50	0.112
10.08	914,241	3.34	1.57	195.30	0.101
10.47	1,358,983	4.96	2.33	211.70	0.139
10.60	453,703	1.66	0.78	196.20	0.051

Compound Table

#	RT [min.]	Mass	Name	CAS#	Match
1	3.524	181.00	Anhydroecgonine methyl ester	43021-26-7	***97.06***
2	5.888	204.10	Levamisole	14769-73-4	***99.31***
3	6.615	303.10	Cocaine	50-36-2	***99.43***
4	8.077	394.50	Octacosane	630-02-4	***98.74***

RT [min.]	Name	Formula	Match	Mass	Library
3.52	Anhydroecgonine methyl ester	C10H15NO2	***97.06***	181.00	SWGDRUG_112916.L
3.52	Anhydroecgonine methyl ester	C10H15NO2	***97.06***	181.00	SWGDRUG.L
3.52	2-Carbomethoxy-8-methyl-8-azabicyclo[3.2.1]oct-2-ene	C10H15NO2	***97.06***	181.10	NIST11.L
3.52	Ecgonidine, methyl ester	C10H15NO2	***90.39***	181.10	NIST11.L
3.52	2-Carbomethoxy-8-methyl-8-azabicyclo[3.2.1]oct-2-ene	C10H15NO2	85.20	181.10	NIST11.L

RT [min.]	Name	Formula	Match	Mass	Library
5.89	Levamisole	C11H12N2S	***99.31***	204.10	NIST11.L
5.89	Levamisole	C11H12N2S	***99.00***	204.00	SWGDRUG.L
5.89	Levamisole	C11H12N2S	***99.00***	204.00	SWGDRUG_112916.L
5.89	(-)-Tetramisole	C11H12N2S	***98.85***	204.10	NIST11.L
5.89	Levamisol	C11H12N2S	***98.56***	204.00	POLLAB.L

RT [min.]	Name	Formula	Match	Mass	Library
6.62	Cocaine	C17H21NO4	***99.43***	303.10	NIST11.L
6.62	Cocaine [PL940]	C17H21NO4	***98.94***	303.20	REFSPEC.L
6.62	Cocaine, Ph.Eur		***98.63***		POLLAB.L
6.62	Cocaine	C17H21NO4	***98.37***	303.10	NIST11.L
6.62	Cocaine	C17H21NO4	***98.09***	303.10	NIST11.L

RT [min.]	Name	Formula	Match	Mass	Library
8.08	Octacosane	C28H58	***98.74***	394.50	NIST11.L
8.08	n-Octacosane [PL1090]	C28H58	***97.30***	394.50	REFSPEC.L
8.08	Hentriacontane	C31H64	***95.04***	436.50	NIST11.L
8.08	Tetracosaan	C24H50	***94.76***	338.00	POLLAB.L
8.08	Octacosane	C28H58	***94.62***	394.50	NIST11.L

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