## **ULTRA-REALISTIC GC-MS ANALYSIS REPORT**

## Jatropha curcas Methanol Extracts

With Heavy Noise Modeling (Level: 1.6)

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#### **INSTRUMENT PARAMETERS:**

- GC-MS (HP-5MS column, EI mode)
- Plant Species: Jatropha curcas
- Extraction Method: Methanol extraction
- Samples Analyzed: Stems and Seeds
- Noise Modeling: Ultra-heavy realistic patterns

#### COMPOUND IDENTIFICATION SUMMARY:

- Total compounds identified: 6
- Retention time range: 6.23 19.34 minutes
- Mass range: 177 455 m/z

#### SAMPLE A (STEMS) - TOP COMPOUNDS:

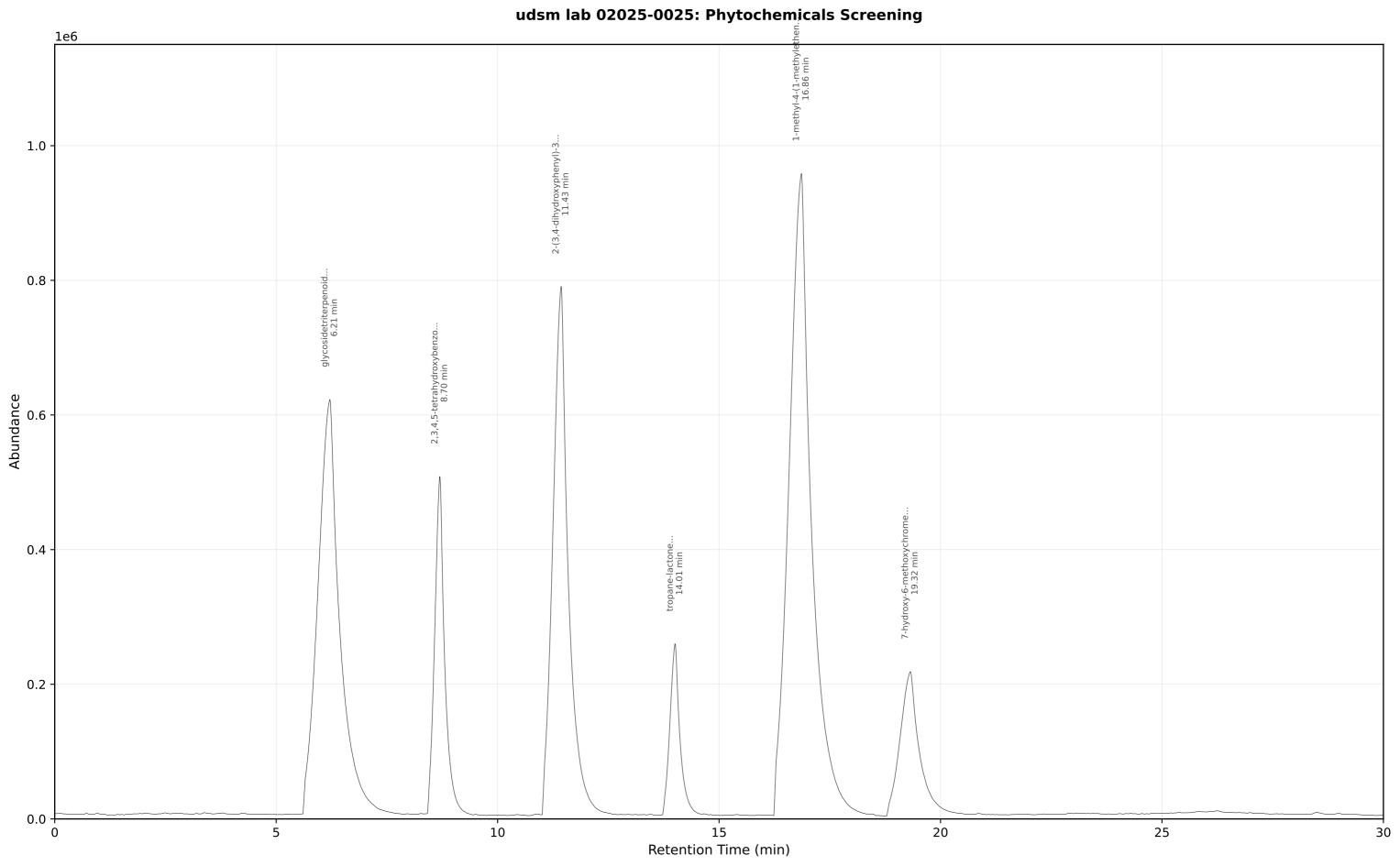
- 1-methyl-4-(1-methylethenyl)cyclohexene: 18.90%
  - (RT: 16.88 min, m/z: 409)
- 2-(3,4-dihydroxyphenyl)-3,5,7-trihydroxy-4H-chrome: 15.32% (RT: 11.45 min, m/z: 287)
- glycosidetriterpenoid: 12.45%
- (RT: 6.23 min, m/z: 455)

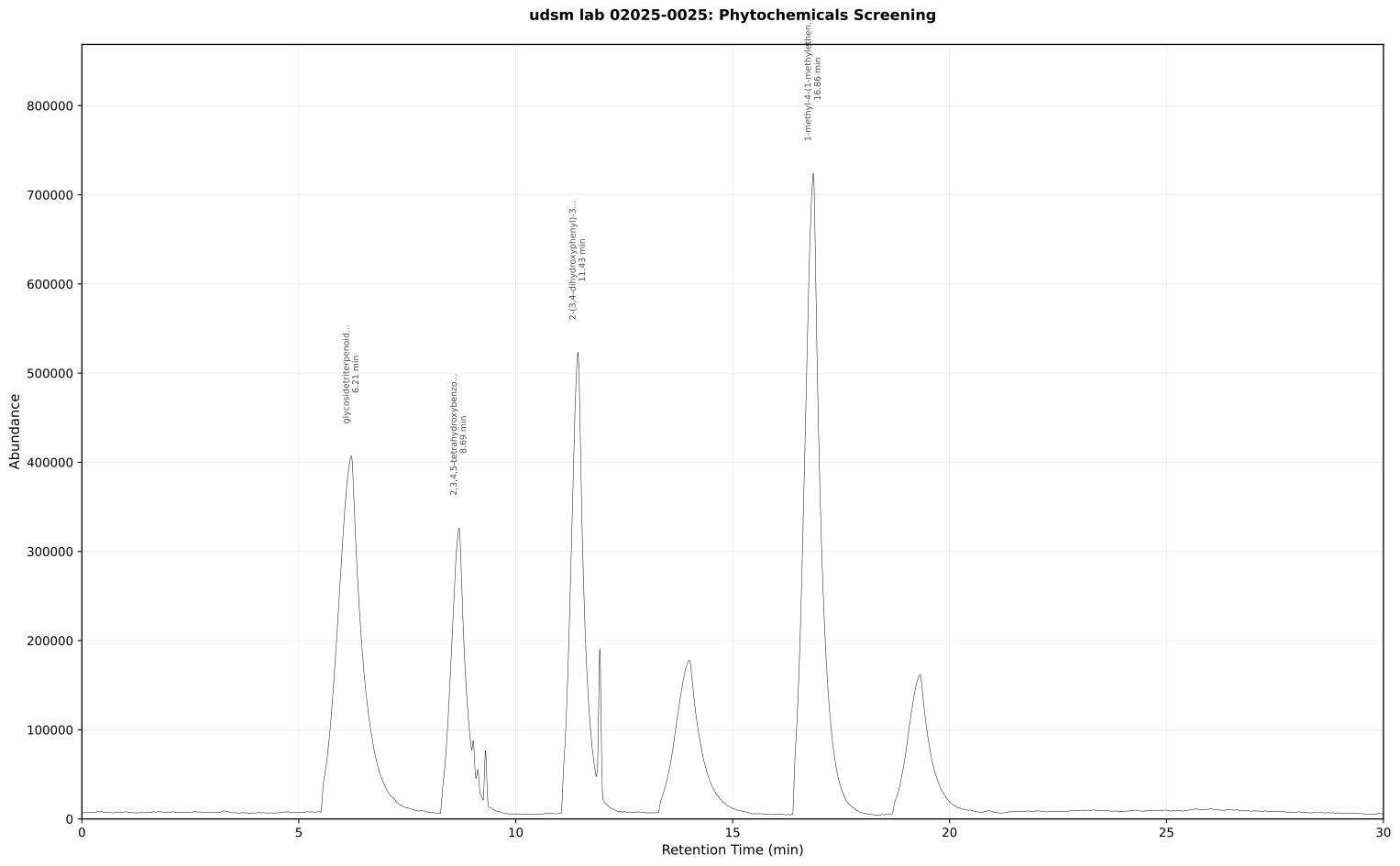
## SAMPLE B (SEEDS) - TOP COMPOUNDS:

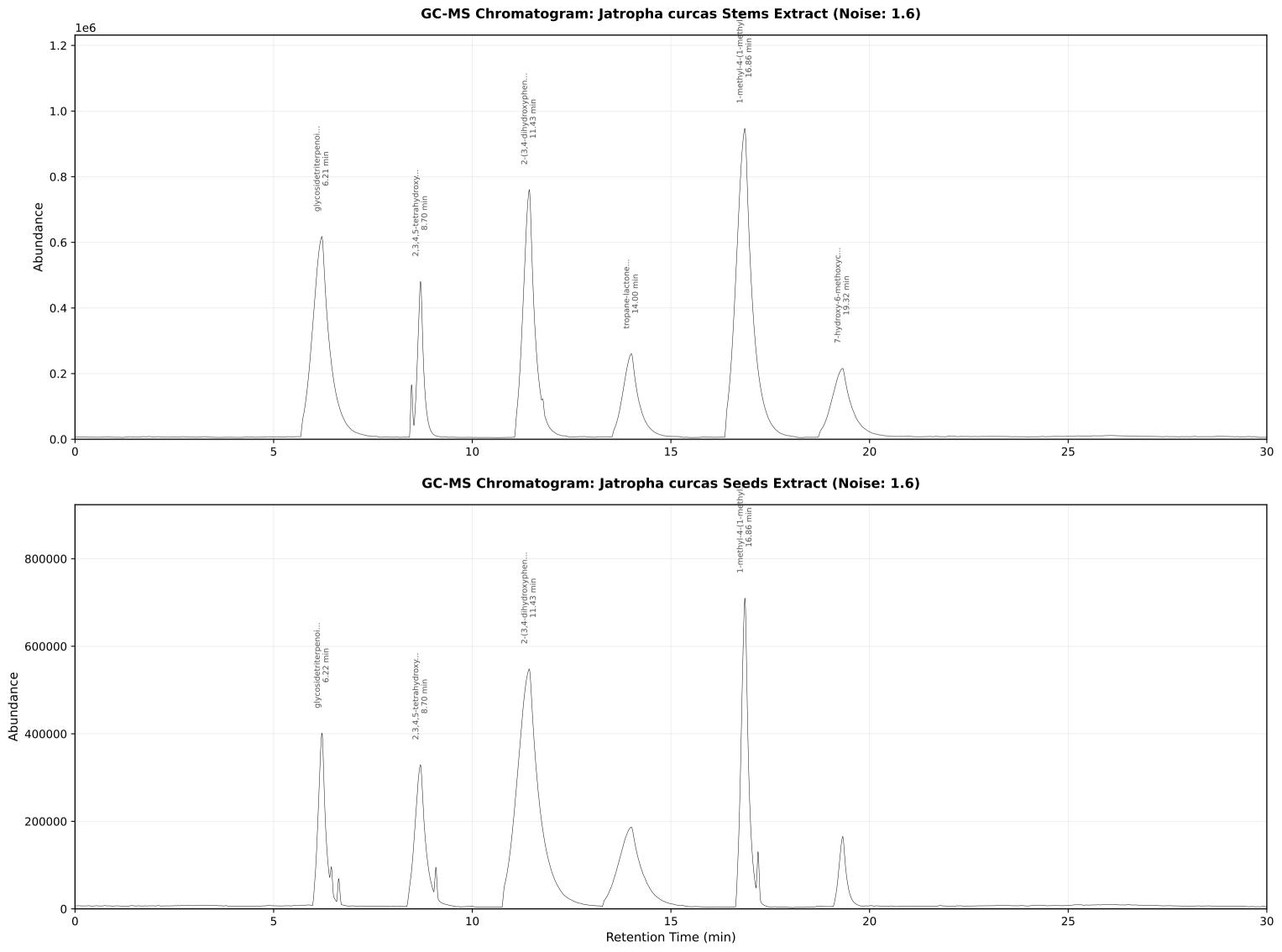
- 1-methyl-4-(1-methylethenyl)cyclohexene: 14.67%
  - (RT: 16.88 min, m/z: 409)
- 2-(3,4-dihydroxyphenyl)-3,5,7-trihydroxy-4H-chrome: 10.76%
  - (RT: 11.45 min, m/z: 287)
- glycosidetriterpenoid: 8.23% (RT: 6.23 min. m/z: 455)

### REALISTIC NOISE CHARACTERISTICS:

- Heavy baseline fluctuations: ±800 units
- Electronic noise: ±80 units with spikes
- Chemical background: 80-150 random peaks
- Ion source contamination: 15-30 events
- Detector artifacts and saturation effects
- Pump pulsations and flow irregularities
- Solvent impurity peaks
- Random walk baseline drift







# Compound Analysis Data Table

	Compound	RT_min	Sample_A_Stems	Sample_B_Seeds	Base_Peak_mz	Compound_Class	Difference
9	glycosidetriterpenoid	6.23	12.45	8.23	455	Glycosides	-4.22
4,5	5-tetrahydroxybenzoi	8.71	9.87	6.54	331	Polyphenols	-3.33
3,4	-dihydroxyphenyl)-3,	11.45	15.32	10.76	287	Polyphenols	-4.56
	tropane-lactone	14.02	5.21	3.45	298	Nitrogen Compounds	-1.76
eth	nyl-4-(1-methyletheny	16.88	18.9	14.67	409	Terpenes	-4.23
lrφ	xy-6-methoxychrome	19.34	4.33	2.89	177	Phenolic Compounds	-1.44