ULTRA-REALISTIC GC-MS ANALYSIS REPORT

Jatropha curcas Methanol Extracts

With Heavy Noise Modeling (Level: 1.4)

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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: 8
- Retention time range: 6.45 25.87 minutes
- Mass range: 179 457 m/z

SAMPLE A (STEMS) - TOP COMPOUNDS:

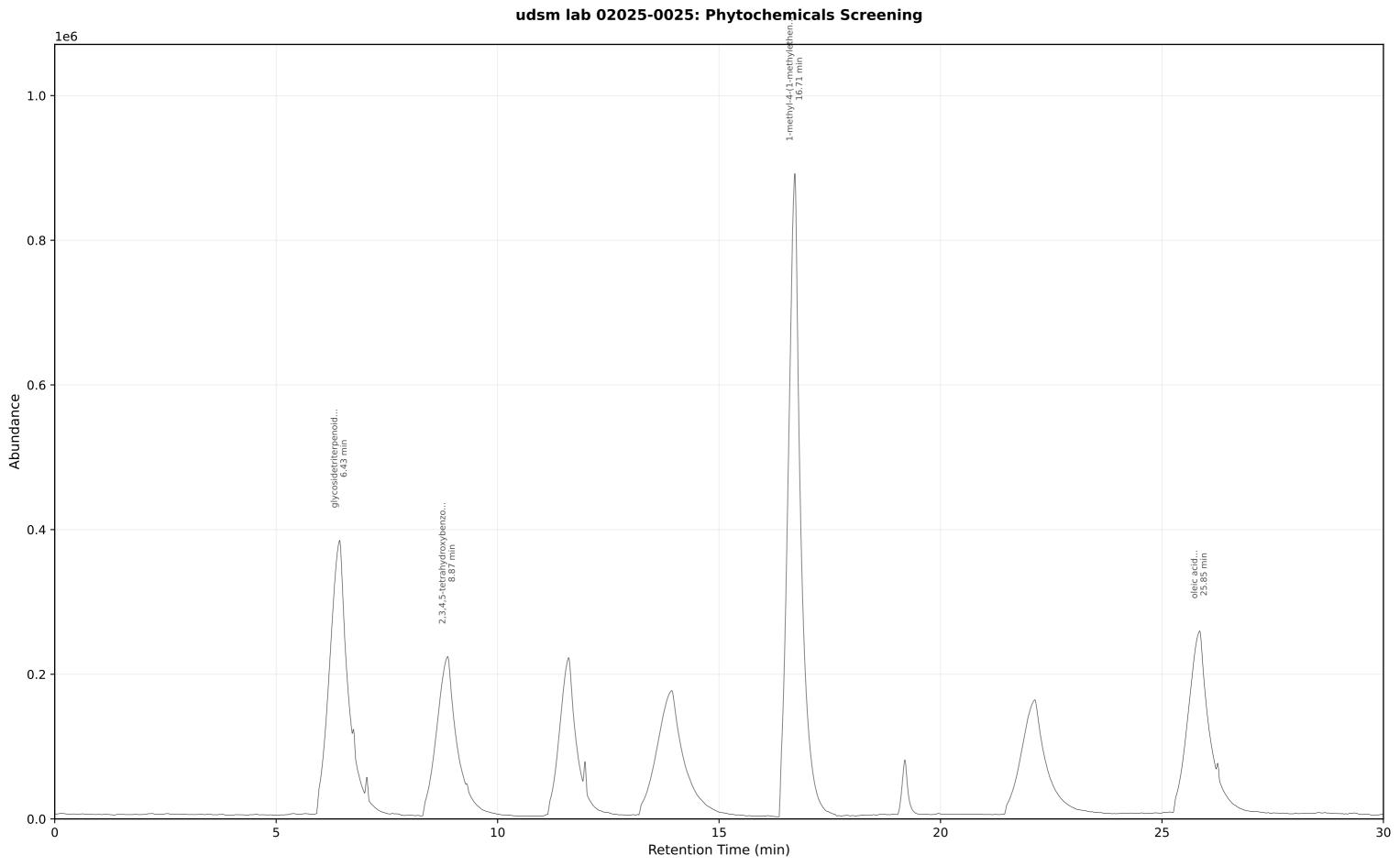
- 1-methyl-4-(1-methylethenyl)cyclohexene: 17.45%
 - (RT: 16.73 min, m/z: 411)
- glycosidetriterpenoid: 7.84%
- (RT: 6.45 min, m/z: 457)
- oleic acid: 5.67%
- (RT: 25.87 min, m/z: 284)

SAMPLE B (SEEDS) - TOP COMPOUNDS:

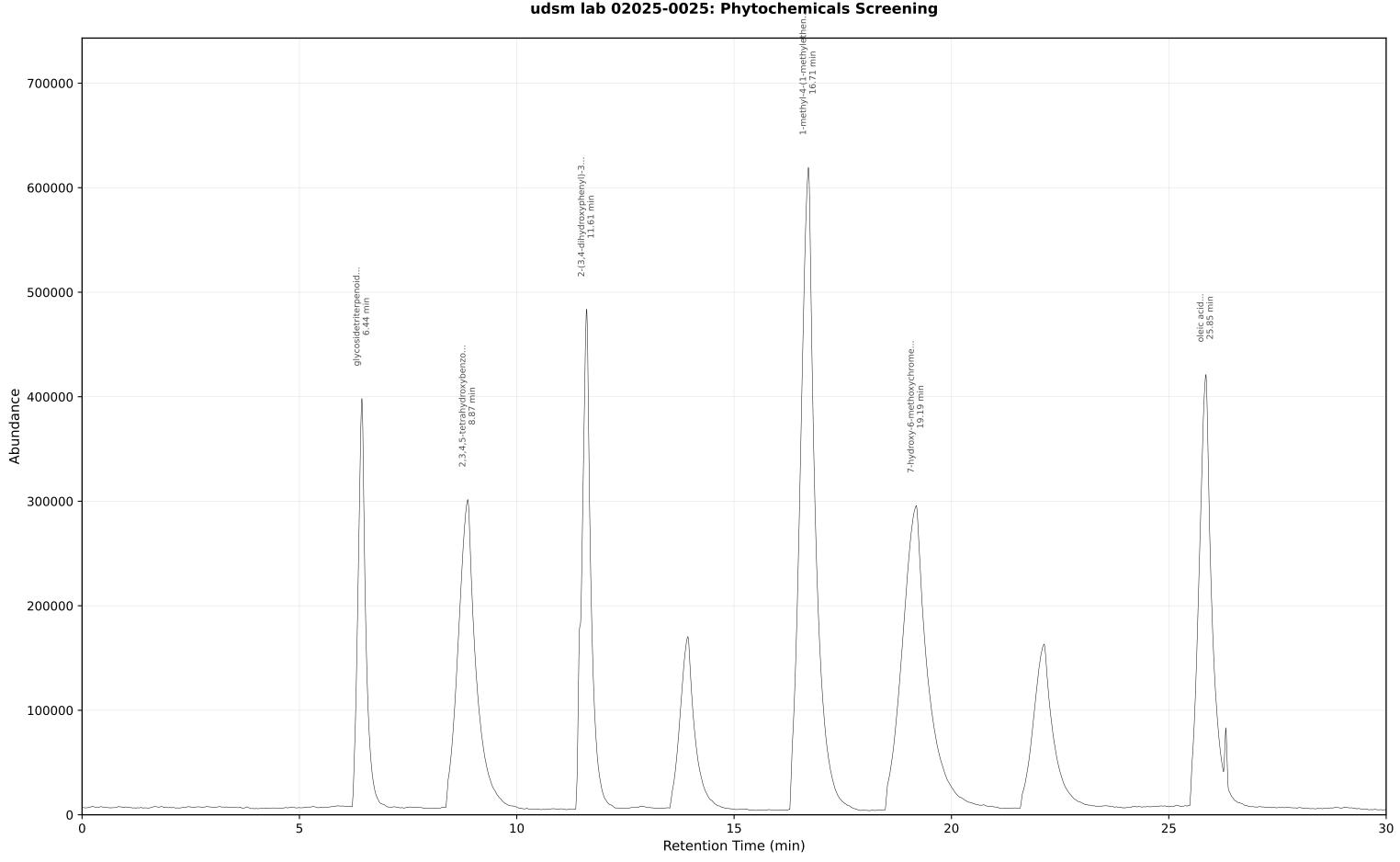
- 1-methyl-4-(1-methylethenyl)cyclohexene: 12.98%
 - (RT: 16.73 min, m/z: 411)
- 2-(3,4-dihydroxyphenyl)-3,5,7-trihydroxy-4H-chrome: 9.41% (RT: 11.62 min, m/z: 289)
- oleic acid: 8.34%
 - (RT: 25.87 min, m/z: 284)

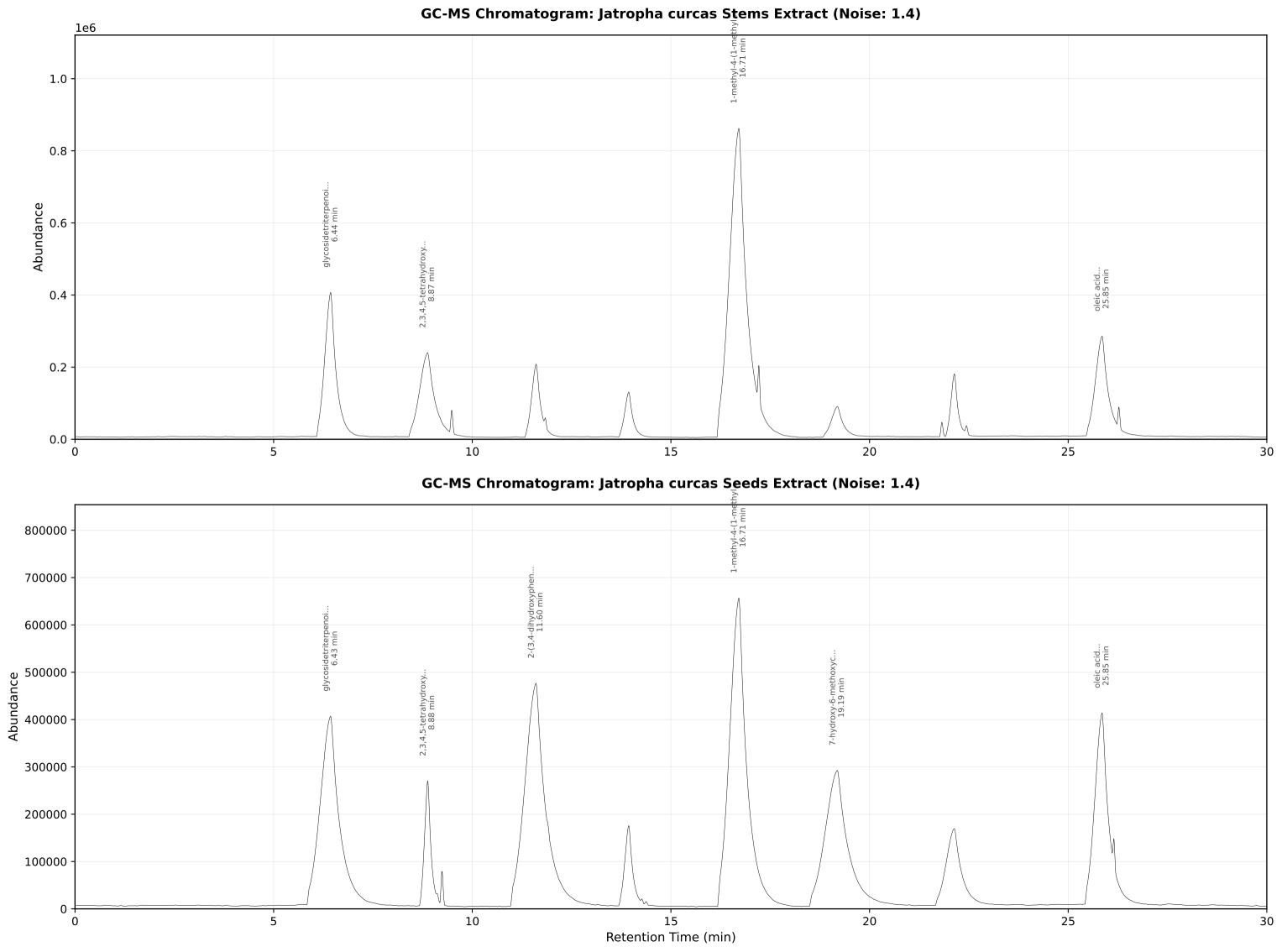
REALISTIC NOISE CHARACTERISTICS:

- 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 effectsPump pulsations and flow irregularities
- Solvent impurity peaks
- Random walk baseline drift



udsm lab 02025-0025: Phytochemicals Screening





Compound Analysis Data Table

Compound	RT_min	Sample_A_Stems	Sample_B_Seeds	Base_Peak_mz	Compound_Class	Difference
glycosidetriterpenoid	6.45	7.84	7.69	457	Glycosides	-0.15
4,5-tetrahydroxybenzoi	8.89	4.56	5.47	333	Polyphenols	0.91
,4-dihydroxyphenyl)-3,	11.62	3.92	9.41	289	Polyphenols	5.49
tropane-lactone	13.95	2.87	3.11	300	Nitrogen Compounds	0.24
ethyl-4-(1-methyletheny	16.73	17.45	12.98	411	Terpenes	-4.47
roxy-6-methoxychrome	19.21	1.92	5.63	179	Phenolic Compounds	3.71
palmitic acid	22.15	3.21	2.89	195	Acids	-0.32
oleic acid	25.87	5.67	8.34	284	Lipids	2.67