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GC-MS Analysis of Bioactive Compound in Bark Extract of *Butea monosperma* (Lam) Taub. in Jharkhand

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ABSTRACT

Plants produce a diverse range of bioactive molecules, making them rich sources of different types of medicine. An essential part in the investigation of plant is the identification of the biologically active compounds present in plant leading to further biological and pharmacological studies. The bark of *Butea monosperma* was collected from the Kulhi Jungle at Panchyat Kuchu of Block Ormanjhi in Ranchi District of Jharkhand. 1.5 ml of extract in a GC vial was subjected to the GC-MS equipment. The biomolecules investigation of organic extract was performed by GC-MS equipment (GCMS-QP2010 Ultra). Bis(2-ethylhexyl) phthalate and Lup-20(29)-en-3-one were abundant in the organic extract which also express apoptosis inhibitor and anti-HIV properties and anti-bacterial activities respectively.

Keywords: Bark, Butea monosperma, GCMS, Bis(2-ethylhexyl) phthalate, Lup-20(29)-en-3-one

INTRODUCTION

Plants produce a diverse range of bioactive molecules, making them rich sources of different types of medicine (Nair *et al.*, 2017). An essential part in the investigation of plant is the identification of the biologically active compounds present in plant leading to further biological and pharmacological studies (Farid *et al.*, 2015; Guo *et al.*, 2013). *Butea monosperma* also called Palash, is commonly known as the Flame of the forest, belongs to the family Fabaceae (Patil *et al.*, 2014). Bark is having anti-microbial properties, anti-fungal activity and used in tumors, bleeding piles, ulcers (Tambekar and Khante, 2010; Singh, 2011; Patil *et al.*, 2006). Secondary metabolites are the important source with a variety of structural arrangements and properties (Pavitra and Vadivukarasi, 2012). In the last few years, gas chromatography mass spectrometry (GC-MS) has become firmly established as a key technological platform for secondary metabolite

profiling in both plant and non-plant species (Gani, 2003; Fernie *et al*, 2004; Kell *et al.*, 2005). Volatile compounds are identified by the GC-MS analysis (Hassanpouraghdam, 2009). GC-MS is the best technique to identify the bioactive constituents of long chain hydrocarbons, alcohols, acids, esters, alkaloids, steroids, amino and nitro compounds etc. (Vinodh *et al.*, 2013).

Present study comprises the bioactive compound identification of Bark extract of Butea monosperma.

MATERIAL AND METHOD

Collection of Plant Material

The bark of *Butea monosperma* was collected from the Kulhi Jungle at Panchyat Kuchu of Block Ormanjhi in Ranchi District of Jharkhand.

Preparation of extract

The barks were dried in a dark place at

room temperature and processed for fine powder mixture-grinder followed by sieving and stored in an air tight container. Then, the fine powder of barks was subjected to methanol as organic solvent. The extract was collected using the soxhlet apparatus.

GC-MS analysis

1.5 ml of extract in a GC vial was subjected to the GC-MS equipment. The biomolecules investigation of organic extract was performed by GC-MS equipment (GCMS-QP2010 Ultra) followed by Gas Chromatography through Column Oven Temp. :100.0 °C, Injection Temp. : 260.00 °C, Injection Mode :Split, Flow Control Mode :Linear Velocity, Pressure :90.5 kPa, Total Flow :16.3 mL/min, Column Flow :1.21 mL/ min, Linear Velocity :40.9 cm/sec, Purge Flow :3.0 mL/ min, Split Ratio :10.0, Gas : Helium(He) and Mass Spectroscopy through Start Time :4.00min, End Time : 39.98 min, ACQ Mode :Scan, Event Time :0.20sec,Scan Speed :3333, Start m/z : 40.00 and End m/z :650.00. A mass spectrum observed.

	Sample Information
Analyzed by	: \$Admn.\$
Analyzed	: 1/20/2021 5:01:50 PM
Sample Type	: \$Organic\$
Sample Name	:5
Method File	: D:\GCMS\GCMS METHOD\Organic\Extract.qgm

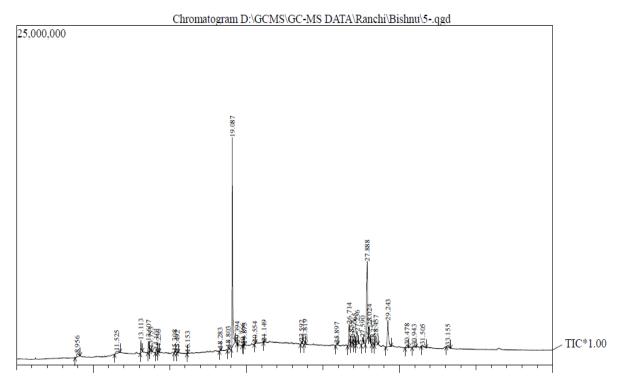


Figure 1: Chromatogram of GC-MS analysis of Bark Extract of Butea monosperma

RESULTS AND DISCUSSION

Interpretation of mass spectrum was done using database of NIST and WILEY Library. The mass spectrum of unknown compound was compared with the spectrum of known compound stored in the NIST'S and WILEY's Library. Compounds were identified by with authentic standards and by with recorded from computerized libraries.

Table 1: Peaks with Retention time against Area in GC-MS analysis of Bark Extract of Butea monosperma.

Peak#	R.Time	Area	Area%	Name			
1	8.956	939073	1.67	2,7-Nonadienoic acid, 3,8-dimethyl-, methyl ester, (Z)-			
2	11.525	869572	1.55	MOME INOSITOL			
3	13.113	1341580	2.39	1,2-Benzenedicarboxylic acid, bis(2-methylpropyl) este			
4	13.607	1239423	2.21	1,2-Benzenedicarboxylic acid, bis(2-methylpropyl) es			
5	13.753	331055	0.59	1,2-Benzenedicarboxylic acid, bis(2-methylpropyl) es			
6	14.101	387728	0.69	Dibutyl phthalate			
7	14.256	128822	0.23	1,2-BENZENEDICARBOXYLIC ACID, BIS(2-METHO2			
8	15.308	119033	0.21	TETRADECANE, 1-CHLORO-			
9	15.492	114419	0.20	2,5-Hexanediamine, 2,5-dimethyl-			
10	16.153	43221	0.08	Cyclononasiloxane, octadecamethyl-			
11	18.283	177843	0.32	Eicosanal-			
12	18.803	333973	0.60	1-Heptacosanol			
13	19.087	17868876	31.86	BIS(2-ETHYLHEXYL) PHTHALATE			
14	19.394	50996	0.09	1H-PURIN-6-AMINE, [(2-FLUOROPHENYL)METHYL			
15	19.766	79125	0.14	1-Methyl-7-azabicyclo[4.1.0]hepta-2,4-diene-7-carboxylic			
16	19.875	87933	0.16	Hexacosanal			
17	20.554	67201	0.12	Tetracosanoic acid, methyl ester			
18	21.149	57845	0.10	2,6,10,14,18,22-TETRACOSAHEXAENE, 2,6,10,15,19,2			
19	23.592	231660	0.41	Stigmasta-5,22-dien-3-ol, acetate, (3.beta.)-			
20	23.819	310385	0.55	STIGMAST-5-EN-3-OL, OLEAT			
21	25.897	265690	0.47	Stigmasterol			
22	26.714	3000264	5.35	Lup-20(29)-en-3-one			
23	26.859	298391	0.53	.gammaSitosterol			
24	27.065	1312884	2.34	.betaAmyrone			
25	27.236	2061807	3.68	Lupeol			
26	27.590	1040289	1.86	4-Campestene-3-one			
27	27.888	12590340	22.45	Lup-20(29)-en-3-one			
28	28.024	1506994	2.69	4,22-Stigmastadiene-3-one			
29	28.257	362926	0.65	9,19-Cyclolanost-24-en-3-ol, (3.beta.)-			
30	28.457	1832999	3.27	METHYL COMMATE B			
31	29.243	5148253	9.18	.gammaSitostenone			
32	30.478	593401	1.06	hexadecanamide, N-methyl-N-(1-oxohexadecyl)-			
33	30.943	402936	0.72	17-[5-(1-HYDROXY-1-METHYL-ETHYL)-2-METHYL-			
34	31.505	383427	0.68	Stigmasta-7,22-dien-3-ol, acetate, (3.beta.,5.alpha.,22E)-			
35	33.155	499603	0.89	Stigmastane-3,6-dione, (5.alpha.)-			
		56079967	100.00				

Mention in the Figure 2 confirms the compounds present in the organic extract. There were 14 peaks observed out of 35 peaks those were greater than 1% in area also mentioned in the Table 1 that comprises Name of the Compound, Molecular Formula, Molecular weight, Peak area (%) and Molecular structure. Peak 19.087 expressed the highest area percent that is 31.86 among them and identified as bis (2-ethylhexyl) phthalate. Bis (2-ethylhexyl) phthalate is a phthalate ester that is the bis(2-ethylhexyl) ester. It has a role as an apoptosis inhibitor, an androstane receptor agonist (<u>PUBCHEM</u>). Lup-20(29)-en-3-one was the second most abundant compound in the organic extract of Bark of Butea monosperma which was found to express the antiHIV properties and anti bacterial activities as well (Callies *et al.*, 2015, Madureira *et al.*, 2003; François *et al.*, 2020).

Table 2: Name of the Compound, Molecular Formula, Molecular weight, Peak area(%) and Molecular structure expressed in Bark's extract of *Butea monosperma*.

SI. No.	Name of the Compound	Molecular Formula	Molecular Weight	Peak Area %	Molecular Structure
1	2,7-Nonadienoic acid, 3,8- dimethyl-, methyl ester, (Z)-	C12H20O2	196	2.49	
2	MOME INOSITOL	C7H14O6	194	2.31	но он он
3	1,2-Benzenedicarboxylic acid, bis(2-methylpropyl) ester	C ₁₆ H ₂₂ O ₄	278	3.56	
4	1,2-Benzenedicarboxylic acid, bis(2-methylpropyl) ester	$C_{16}H_{22}O_4$	278	3.29	
5	Dibutyl phthalate	$C_{16}H_{22}O_4$	278	1.03	
6	TETRADECANE, 1-CHLORO	C14H29Cl	232	0.32	······
7	2,5-Hexanediamine, 2,5- dimethyl-	$C_8H_{20}N_2$	144	0.30	H2N NH2
8	Cyclononasiloxane, octadecamethyl-	C18H54O9Si9	666	0.11	
9	Eicosanal-	C ₂₀ H ₄₀ O	296	0.47	~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~
10	1-Heptacosanol	C27H56O	396	0.89	~~~~~он

		a 11 No		0.04	
11	1-Methyl-7- azabicyclo[4.1.0]hepta-2,4- diene-7-carboxylic acid, 3,17-diacetoxy-4,4,10,13- tetramethylhexadecahydroc yclopenta[a]phenan threne	C33H47NO6	553	0.21	
12	Hexacosanal	C ₂₆ H ₅₂ O	380	0.23	
13	Tetracosanoic acid, methyl ester	$C_{25}H_{50}O_2$	382	0.18	
14	2,6,10,14,18,22- TETRACOSAHEXAENE, 2,6,10,15,19,23- EXAMETHYL	C ₃₀ H ₅₀	410	0.15	for the second s
15	Stigmasta-5,22-dien-3-ol, acetate, (3.beta.)-	C ₃₁ H50O2	454	0.61	H H H H
16	STIGMAST-5-EN-3-OL, OLEAT	C47H82O2	678	0.82	
17	Stigmasterol	C ₂₉ H ₄₈ O	412	0.70	
18	Lup-20(29)-en-3-one	$C_{30}H_{48}O$	424	7.96	
19	.gammaSitosterol	C29H50O	414	0.79	
20	.betaAmyrone	$C_{30}H_{48}O$	424	3.48	
21	Lupeol	C ₃₀ H ₅₀ O	426	5.47	
22	4-Campestene-3-one	C28H46O	398	2.76	

	1 20(20) 2	0 11 0	42.4	00.40	1
23	Lup-20(29)-en-3-one	C30H48O	424	33.40	
24	4,22-Stigmastadiene-3-one	C29H46O	410	4.00	
25	9,19-Cyclolanost-24-en-3- ol, (3.beta.)-	C ₃₀ H ₅₀ O	426	0.96	H OV
26	METHYL COMMATE B	C31H50O3	470	4.86	
27	.gammaSitostenone	C29H48O	412	13.66	
28	hexadecanamide, N-methyl- N-(1-oxohexadecyl)-	C ₃₃ H ₆₅ NO ₂	507	1.57	
29	17-[5-(1-HYDROXY-1- METHYL-ETHYL)-2- METHYL-TETRAHYDRO- FURAN-2-YL]-4,4,10,13,14- PENTAMETHYL-2,3,4,5,6,7, 8,10,12,13,14,15,16,17- TETRADECAHYDRO-1H- CYCLOPENTA[A]PHENANT HRENE-3,6,16-TRIOL	C30H50O5	490	1.07	
30	Stigmasta-7,22-dien-3-ol, acetate, (3.beta.,5.alpha.,22E)-	C31H50O2	454	1.02	
31	Stigmastane-3,6-dione, (5.alpha.)-	C29H48O2	428	1.33	

CONCLUSION:

Presence of Mome inositol, Lupeol and Lupen-3-one in barks of Butea monosperma makes reliable to proceed further for pharmacology to design drugs.

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Conflicts of Interest: The authors declare no conflict of interest.

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