

Comparative GC-MS Analysis Of Gomutra Sadhita Triphala Kwatha And Gomutra Sadhita Triphala Arka

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ABSTRACT:

Ayurveda is the science of life, which is dealing with life for 5000 years without any change in basic principles. Panchavidha Kashaya Kalpana are primary preparations that are considered underneath fundamental principles of Bhaishajya Kalpana. They are Swarasa, Kalka, Kwatha, Hima and Phanta and are recommended to consume immediately (sadyosevana). Kwatha Kalpana being one of them has low shelf life, hence same Kalpana can be converted into Arka Kalpana to increase its shelf life up to a year. Arka Prakasha, the classical Ayurveda book mentions the distillation techniques for separation of the volatile bioactive compounds from a drug by using Arka Yantra. In the present study Gomutra sadhita triphala kwatha formulation mentioned in Shotharogadhikara in Bhaishajya Ratnavali textbook is modified into Arka Kalpana taking media as Gomutra. Furthermore, traditional Gomutra sadhita triphala kwatha (GTK) and Gomutra sadhita triphala Arka (GTA) both were subjected to GC-MS analysis.

Method: The present study mainly focuses on GC-MS evaluation of *Gomutra sadhita triphala kwatha* and *Gomutra sadhita triphala Arka* for detection of different components present in them. The comparative analysis will define the efficacy of respective dosage form. The preparation of both the forms of medicines was conducted at GMP certified Parul Ayurveda Pharmacy, Parul university.

Conclusion: Present study reveals that both dosage forms GTK and GTA contains entirely different components. *Arka* contains more volatile compounds having Anti-inflammatory effect. on contrary the *Kwatha* has the components which can exert protein denaturation and thereby act as anti-inflammatory.

Keywords- Kwatha Kalpana, Arka Kalpana, *Gomutra sadhita triphala kwatha* (GTK), *Gomutra sadhita triphala Arka* (GTA), GC-MS, Anti-inflammatory action.

INTRODUCTION:

Ayurveda is the science of life, which is dealing with life for 5000 years without any change in basic principles. Ayurveda gives us the freedom to modify system without changing its basic concepts. Panchavidha Kashaya Kalpana that are Swarasa, Kwatha, Kalka, Hima and Phanta are the basic preparations explained under Bhaishajya Kalpana (1). All these Kalpanas has got its own importance but has shelf life of only 24 hours. Kwatha Kalpana is also having low shelf life hence same Kalpana can be converted into Arka Kalpana to increase its shelf life upto 1 year. The general dose of the Kwatha Kalpana mentioned is 2 Pala i.e., 96ml as per AFI (2) whereas the dose of Arka depends on mode of administration. The AFI (3) mentioned 12-24ml of dose in most of the Arka. In Arka Prakasha, Arka Kalpana is also considered as Panchavidha Kashaya Kalpana (4). Ayurveda Sara Sangraha (5) also mentioned the Arka Kalpana in Arka Prakarana as it explains the quick therapeutic action of Arka due to its Sara, Rupa, Laghu, Sheegra Paka and Drava Rupa. Shotha in Ayurveda can be compared with inflammation in modern aspect. Inflammation is the tissue reaction, succession of changes that occur in living tissues when it is injured. Inflammation is the vascular response of living tissue to injury (6) whereas Shopha is defined as a swelling which is Pruthu (wide spread), Grathita presence of granthi), Sama/Vishama (regular or irregular) and is due to the accumulation of Doshas between the Twak and Mamsa in a particular site of the body (7) Gomutra sadhita triphala kwatha is a preparation given in classical textbook Bhaishajya Ratnavali in Shotha Rogadhikara.(8) Most herbal medicines prepared from crude plant extracts comprise a complex mixture of different phytochemical constituents (plant secondary metabolites). The chemical features of these constituents differ

considerably among different *Kalpana*. GC-MS method used for the analysis of the obtained extracts can be an interesting tool for testing the amount of some active principles in herbs used in therapeutic purpose. The aim of this study is to convert the *Kwatha Kalpana* into a new dosage form i.e., *Arka* with equivalent therapeutic effect. Also identify and compare the bioactive compounds of GTK and GTA by subjecting both traditional *Gomutra sadhita triphala kwatha* (GTK) and *Gomutra sadhita triphala Arka* (GTA) for GC-MS analysis.

MATERIALS AND METHODS:

The drugs *Haritaki*, *Bibhitaki* and *Amalaki* were procured from AIMIL pharmaceuticals Pvt. Ltd. Nalagarh (H.P) and *Gomutra* was arranged from local *Gaushala*. The preparation of both the forms of medicines was conducted at GMP certified Parul Ayurveda Pharmacy, Parul university.

Identification and Authentication of drugs was done at AIMIL Pharmaceuticals Pvt. Ltd. Nalagarh (H.P).

INGREDIENTS:

S.NO.	DRUGS	LATIN NAME	FAMILY	PART USED
1.	Amalaki (9)	Emblica officinalis Gaertn	Euphorbiaceae	Dried fruit pericarp
2.	Bibhitaki (10)	Terminalia bellerica Roxb.	Combretaceae	Dried fruit pericarp
3.	Haritaki (11)	Terminalia chebula Retz.	Combretaceae	Dried fruit pericarp
4.	GO-MUTRA	Cow's urine		

Kwatha preparation-: All the three drugs *Haritaki*, *Bibhitaki* and *Amalaki* were coarsely powdered and each weighed 100gm and mixed homogenously. *Gomutra* was collected from local *gaushala* around Parul Ayurved pharmacy and filtered through cotton cloth. Sixteen times of *Gomutra* was used in 300gm of *Triphala* i.e., 4.8 litres. The mixture is boiled until it reduced to 1/8th part to obtain *Gomutra sadhita Triphala Kwatha* (12) and collected in glass beaker for further study.

Arka preparation: Each drug of *Triphala* is coarsely powdered and weighed in equal quantity and mixed homogenously. This coarse powder was soaked in 10 times of *Gomutra* as per AFI (13) for 4 hours. After that the soaked material was taken into the round bottom flask of distillation apparatus and continues controlled heat of about 60 degrees Celsius was given until 60% of *Arka* was collected in the glass bottle. The first 5 ml and last 5ml of Arka was discarded. The prepared Kwatha and Arka was collected in a glass beaker and used for the GCMS study.

GC-MS (GAS CHROMATOGRAPHY MASS SPECTROSCOPY)

The GC-MS of *Gomutra Sadhita Triphala Kwatha* and *Gomutra Sadhita Triphala Arka* was carried out at Sophisticated Instrumentation Centre for Applied Research and Testing (SICART) Anand, Gujarat.

Preparation of both Sample Extract for GC-MS:

MODEL- Perkin Elmer AUTO 6M TURBOMASS. Sample injected at 250°C,

Oven temperature: 75°C.

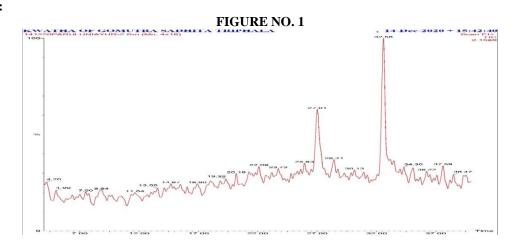
Hold for 5min at the rate of 10°C/min up to 270°C/min; then holded for 10minutes.

Capillary column: PE-5AP. Length of column: 30 meters

ID-0.250 microns Thickness- 0.25mm

Injector temperature: 250°C EI source temperature: 220°C Mass range:20-610AMU. Carrier Gas: Helium gas.

RESULTS:



Area Percent Report

D:\GCMS1\SICART2013.PRO\PeakDB\141220PARULUNIAYUI Monday, December 14, 2020 4:25:46 PM Monday, December 14, 2020 4:25:57 PM Peak List File : Last Modified : Printed :

#	Name	RT	Area	Height	BL	Conc	Units	Area/Conc	m/z	Area %
1	1	27.013	267,186,640.0	679,472,320	MM	0.00		0.00	TIC	30.07
2	2 2	28.414	47,307,052.0	165,531,008	MM	0.00		0.00	TIC	5.32
3	3	32.575	574,001,152.0	513,781,248	MM	0.00		0.00	TIC	64.60

10(1) 2876-2887

Monday, December 14, 2020 4:25:57 PM

FIGURE NO:2

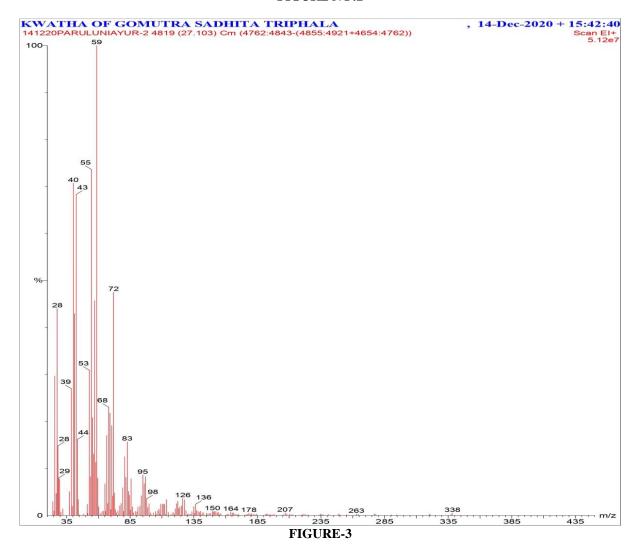
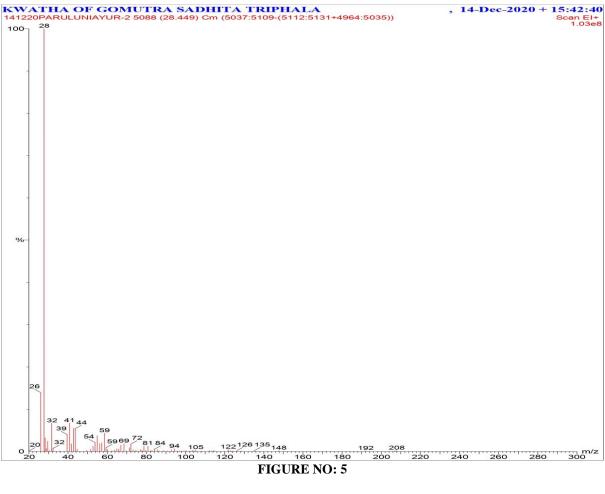




FIGURE NO: 4

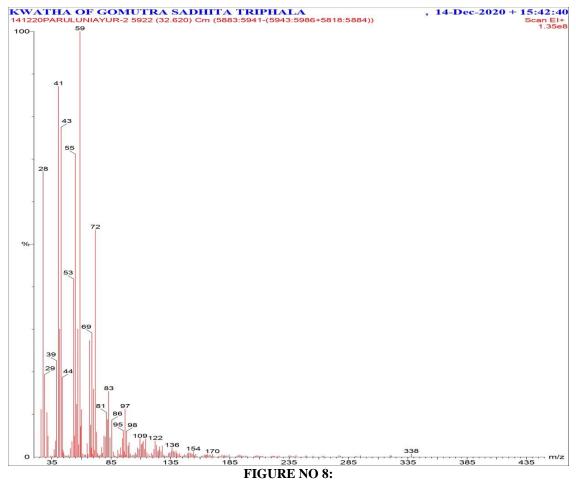


Hit REV	KWATI		MUTRA	A SADHITA TRIPHALA			141220PARULUNIAYUR-2
2 618 504 ÄZETIDINE, 1-NITROSO- 86 C3H6ONZ 15216-10-1 3 577 478 SUCCINIC ANHYDRIDE 100 C4H4O3 108-30-5 4 523 447 1-BUTANAMINE, N-(1-METHYLETHYL)- 5 521 396 ACETIC ACID, [AMINOCARBONYL)AMINO)OXO- 6 520 416 1H-TETRAZOLE, 1-METHYL- 7 514 452 2-ROPENIOL ACID, ETHENYL ESTER 98 C2H4N4 16681-77-9 8 512 436 BUTANAMINE, N-(1-METHYL- 8 512 436 BUTANEDIOL ACID, ETHENYL ESTER 98 C5H6OZ 2177-18-6 8 512 436 BUTANEDIOL ACID, CYCLIC HYDRAZIDE 114 C4H6OZN2 502-95-4 9 498 410 HEXANENITRILE, 6-AMINO- 112 C6H12N12 2432-74-8 10 498 417 SUCCINIMIDE 99 C4H5OZN 123-56-8 11 495 404 2-BUTENE OZONIDE 104 C4H8O3 765-57-1 12 490 426 IRON, TRICARBONYL[(0,1,2,3-ETA.)-METHYL 2-PROPENOATE]- 13 486 361 2-3-DIOXABIOYCLO](2-2-1]HEPTANE 43 C2H5N 151-56-4 14 485 372 ETHYLENIMINE 43 C2H5N 151-56-4 15 485 426 1,3-DIOXABIOYCLO](2-2-1]HEPTANE 43 C2H5N 151-56-4 16 465 388 CYCLOPENTANONE 84 C5H8O 120-92-3 16 465 388 CYCLOPENTANONE 84 C5H8O 120-92-3 17 469 459 391 12,3-TENRENINDE	Hit	REV	for	Compound Name	M.W.	Formula	CAS
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7 514 452 2-PROPENDIC ACID, ETHENYL ESTER 98 C5H6O2 2177-18-6 8 512 436 BUTANEDIDIC ACID, CYCLIC HYDRAZIDE 114 C4H6O2N2 502-95-4 9 498 410 HEXANENITRILE, 6-AMINO- 112 C6H12N2 2432-74-8 10 498 417 SUCCINIMIDE 99 C4H5O2N 123-56-8 11 495 404 2-BUTENE OZONIDE 104 C4H6O3 765-57-1 12 490 426 IRON, TRICARBONYL[(0,1,2,3-ETA,)-METHYL 2-PROPENOATE]- 226 C7H6O5Fe 51922-76-0 13 496 361 2,3-DIOXABICY CLO[2,2,1]-HEPTANE 100 C5H8O2 279-35-6 14 485 372 ETHYLENIMINE 43 C2H5N 151-56-4 15 485 426 13-DIOXABICY CLO[2,2,1]-HEPTANE 144 C8H12N2S 5269-85-2 16 465 388 CYCLOPENTANONE 84 C5H8O 120-92-3 17 483 376 PROPIOLACTONE 72 C3H4O2 57-57-8 18 462 407 2,5-PYRROLIDINEDIONE, 1-HYDROXY- 115 C4H5O3N 6666-82-6 19 459 391 12,3-TRIMETHYLDIAZIRDINE 86 C4H10N2 13604-56-1	6	520	416	1H-TETRAZOLE, 1-METHYL-	84	C2H4N4	16681-77-9
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9 498 410 HEXANENTRILE, 6-AMINO- 10 498 410 SUCCININIDE 99 C41502N 123-56-8 11 495 404 2-BUTENE OZONIDE 104 C41803 765-57-1 12 490 426 IRON, TRICARBONYLI(0,1,2,3-ETA,)-METHYL 2-PROPENOATE]- 226 C714605F6 51922-76-0 13 486 361 2,3-DIOXABICYCLO[2,2,1]HEPTANE 100 C51802 279-35-6 14 495 372 ETHYLENIMINE 43 C215N 151-56-4 15 485 427 1,3-DIAZACYCLOCOTANE-2-THIONE 144 C6112N2S 5289-85-2 16 465 388 CYCLOPENTANONE 84 C5180 120-92-3 17 483 376 PROPIOLACTONE 72 C314002 57-57-8 18 462 407 2,5-PYRROLIDINEDIONE, 1-HYDROXY-19 459 391 12,3-TRIMETHYLDIAZIRDINE 86 C4H10N2 13604-56-1	8	512	436	BUTANEDIOIC ACID, CYCLIC HYDRAZIDE	114	C4H6O2N2	502-95-4
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19 459 391 1,2,3-TRIMETHYLDIAZIRIDINE 86 C4H10N2 113604-56-1							
	19		391				
	20						

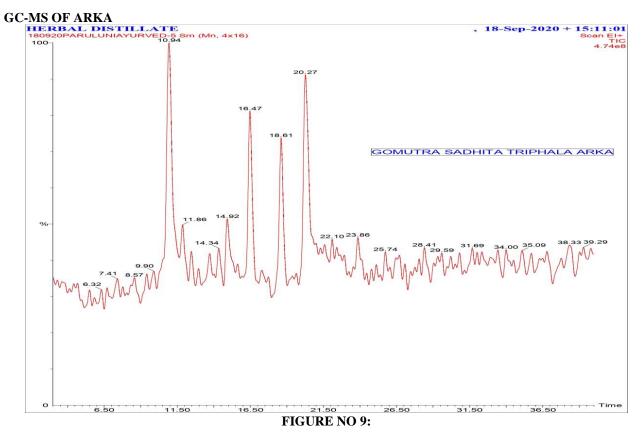
FIGURE NO 6:

KWATE	IA OF GO	MUTRA	A SADHITA TRIPHALA	***		141220PARULUNIAYUR-2
Hit	REV	for	Compound Name	M.W.	Formula	CAS
1	876	661	9-OCTADECENAMIDE, (Z)-	281	C18H35ON	301-02-0
2	839	629	13-DOCOSENAMIDE, (Z)-	337	C22H43ON	112-84-5
3	837	623	13-DOCOSENAMIDE, (Z)-	337	C22H43ON	112-84-5
4	833	597	9-OCTADECENAMIDE	281	C18H35ON	3322-62-1
5	829	680	8-METHYL-6-NONENAMIDE	169	C10H19ON	900293-20-9
6	824	573	9-OCTADECENAMIDE, (Z)-	281	C18H35ON	301-02-0
7	805	542	NONADECANAMIDE	297	C19H39ON	58185-32-3
8	803	558	9-OCTADECENAMIDE, (Z)-	281	C18H35ON	301-02-0
9	799	564	9-OCTADECENAMIDE, (Z)-	281	C18H35ON	301-02-0
10	798	563	9-OCTADECENAMIDE, (Z)-	281	C18H35ON	301-02-0
11	787	614	9-OCTADECENAMIDE	281	C18H35ON	3322-62-1
12	786	561	DODECANAMIDE	199	C12H25ON	1120-16-7
13	784	530	OCTADECANAMIDE	283	C18H37ON	124-26-5
14	767	538	9-OCTADECENAMIDE, (Z)-	281	C18H35ON	301-02-0
15	766	548	N'-(4-NITROBENZYLIDENE)-2-PENTYL-1-CYCLOPROPANECARBOHYDRAZIDE	303	C16H21O3N3	339027-09-7
16	763	624	TRANS-3,4-EPOXYNONANE	142	C9H18O	56769-23-4
17	750	582	CITRONELLOL EPOXIDE (R OR S)	172	C10H20O2	900163-92-8
18	747	490	PENTADECANAMIDE, 15-BROMO-	319	C15H30ONBr	900163-86-1
19	737	606	OXETANE, 2,2,3-TRIMETHYL-	100	C6H12O	23120-43-6
20	735	502	BIS(CIS-13-DOCOSENAMIDO)METHANE	686	C45H86O2N2	10436-19-8

FIGURE NO 7:







Area Percent Report

Peak List File:

D:\GCMS1\SICART2013.PRO\PeakDB\180920APARULUNIAYU Saturday, September 19, 2020 11:25:55 AM Saturday, September 19, 2020 11:26:08 AM Last Modified : Printed :

#	Name	RT	Area	Height	BL	Conc	Units	Area/Conc	m/z	Area %
1	1	10.937	130,712,192.0	298,577,184	MM	0.00		0.00	TIC	31.88
2	2	11.862	14,442,335.0	52,666,196	MM	0.00		0.00	TIC	3,52
3	3	14.919	27,864,182.0	81,944,056	MM	0.00		0.00	TIC	6,80
4	4	16.474	64,466,308.0	211,901,968	MM	0.00		0.00	TIC	15.72
5	5	18.610	63,381,280.0	196,033,248	MM	0.00		0.00	TIC	15.46
6	6	20.271	109,103,392.0	257,105,008	MM	0.00		0.00	TIC	26.61

Saturday, September 19, 2020 11:26:08 AM Area percent report Page 1

FIGURE NO 10:

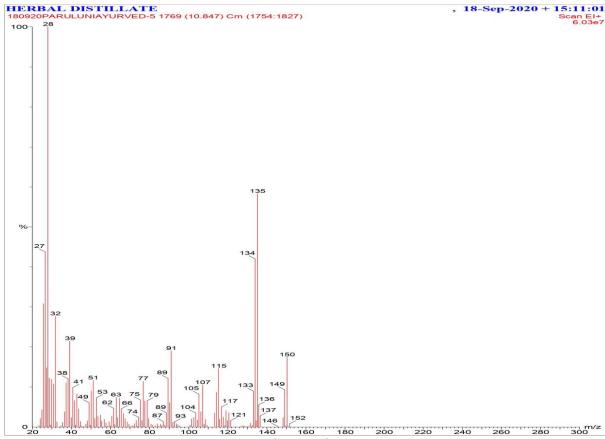


FIGURE NO 11:



FIGURE NO 12:

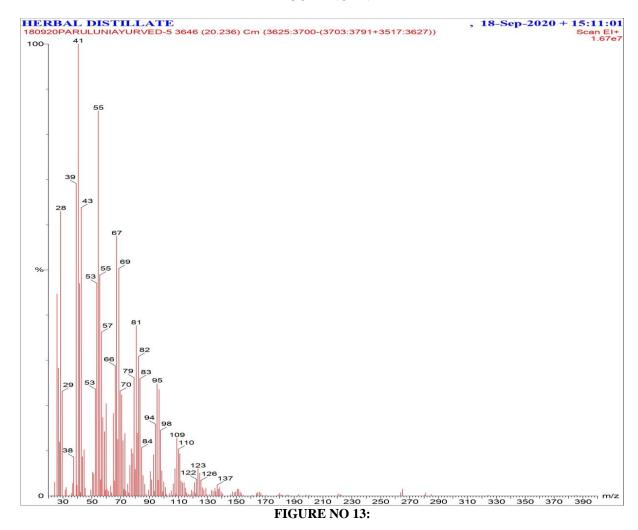




FIGURE NO 14:

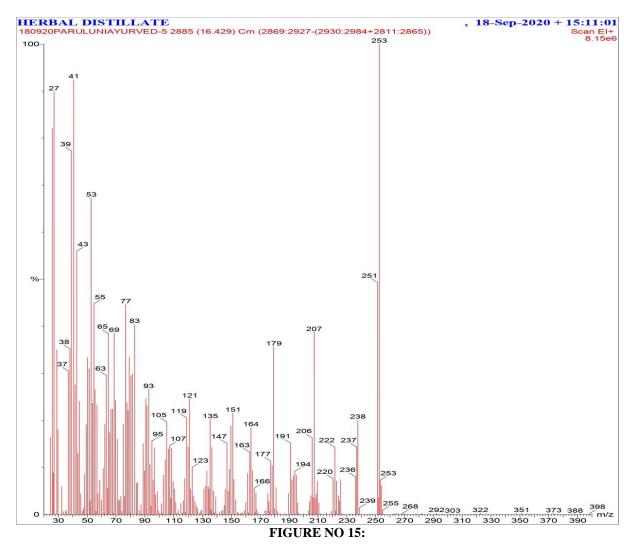
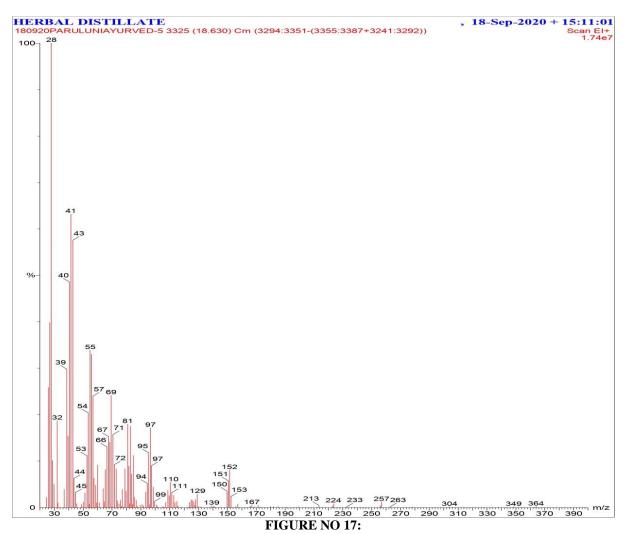




FIGURE NO 16:



Hell	HERBAL DISTILLATE 180920PARULUNIAYO								
1 833 423 2-PIÈERIDINONE, NI-LABROMO-N-BUTYLI- 233 C9H16ONBr 195194-80-0 2 833 408 CVCLOBUTANONE, 2-METHYL-2-OXIRANYL- 126 C7H1002 75314-19-1 3 815 333 11-TRICOSENE 322 C23446 52078-56-5 4 804 446 DODECARE, 12-DIBROMO- 326 C12424Br2 56334-42-4 5 800 415 DISPARLURE 282 C19438O 29804-22-6 6 788 435 2-METHYL-F-7-HEXADECENE 238 C17434 6418-35-24 7 798 392 17-PENTATRIACONTENE 490 C35H70 6971-40-0 8 785 436 7-TETRADECENE 198 C14428 10374-74-0 9 785 436 7-TETRADECENE 240 C19432O 900131-10-1 10 784 437 Z-10-TETRADECENE 254 C18430O 900130-99-3 12 779 429 2-METHYL-Z-18-EPOXYHEXADECENE				Compound Name	M.W.				
2 833 408 CYCLOBUTANONE, 2-METHYL-2-OXIRAÑYL- 3 815 333 11-TRICOSENE 3 820 23446 52078-58-6.5 4 804 446 DODECANE, 1,2-DIBROMO- 5 800 415 DISPARLURE 6 798 435 2-METHYL-E-7-HEXADECENE 7 796 392 17-PENTATRIACONTENE 8 785 436 7-TETRADECENE 9 785 436 E-2-HEXADECENE 9 785 436 E-2-HEXADECENE 10 784 437 2-10-TETRADECEN-1-OL ACETATE 11 781 426 2-METHYL-Z-7-HEXADECENE 12 252 C194380 2980-422-6 13 778 407 SILABLE, TRICHLORODCOSYL- 14 2779 429 2-METHYL-Z-7-HEXADECANE 15 777 422 E-7-OCTADECENE 26 C19438 3535-39-3 15 7774 422 E-7-OCTADECENE 27 900130-99-3 18 789 411 2-METHYL-Z-OCNEENE 28 C7H1002 75314-19-1 29 795 436 E-2-HEXADECANE 29 00130-99-3 20 00130-99-3 20 00130-99-3 20 00130-99-3 20 00130-99-3 21 779 429 2-METHYL-Z-7-HEXADECANE 25 C17H34 900130-87-2 26 C17H34-90130-88-3 27 778 407 SILABLE, TRICHLORODCOSYL- 27 491 3-TRICHLORODCOSYL- 28 C19438 35354-39-3 29 00130-99-3 20 C17H34 22 E-7-OCTADECENE 29 C19438 35354-39-3 20 C19438 35									
4 804 446 DODECANE, 1,2-DIBROMO- 5 800 415 DISPARLURE 5 10;SPARLURE 5 10	2								
4 804 446 DODECANE, 1,2-DIBROMO- 5 800 415 DISPARLURE 5 10;SPARLURE 5 10	3								
5 800 415 DISPARLURE 282 C19438O 29804-22-6 6 798 435 2-METHYLE-7-HEXADECENE 238 C17H34 64183-52-4 7 796 392 17-PENTATRIACONTENE 490 C35H70 6971-40-0 8 785 436 E-2-HEXADECENE 196 C14H28 10374-74-0 9 785 436 E-2-HEXADECACEN-1-OL 240 C16H32O 900131-10-1 10 784 437 Z-10-TETRADECEN-I-OL ACETATE 254 C16H32O 900130-99-3 11 781 426 2-METHYL-Z-1-HEXADECENE 238 C17H34 900130-99-3 12 779 429 2-METHYL-Z-7,8-EPOX/HEXADECANE 254 C17H34O 900130-98-3 13 778 407 SILANE, TRICHLORDODCOSYL- 442 C22H4SCISIS 7325-84-0 14 775 493 GIS-Z-METHYL-7-OCTADECENE 266 C19H38 35354-39-3 774 422 E-7-OCTADECENE 324 <td>4</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>	4								
6 798 435 2-METHYL-E-7-HEXADECENE 238 C17H34 64183-52-4 7969 392 17-PENTATRIACONTENE 490 C35H70 6971-40-0 871-40-0 900131-10-1 871-40-0 900130-93-3 871-40-0 871-40-0 900130-93-3 871-40-0 900130-93-3 871-40-0 900130-93-3 871-40-0 900130-87-2 871-40-0 900130-87-2 871-40-0 900130-88-3 871-40-0 900130-88-3 900130-88-3 900130-88-3 900130-89-3 900130-89-3 900130-93-3	5								
7 796 392 17-PENTATRIACONTENE 490 C35H70 6971-40-0 787-787-786 392 17-PENTADECENE 198 C35H70 6971-40-0 787-787-785 436 E-2-HEXADECACEN-1-OL 240 C16H32O 900131-10-1 10 784 437 2-10-TETRADECEN-1-OL ACETATE 254 C16H33OC 900130-90-3 11 781 426 2-METHYL-Z-7-HEXADECENE 238 C17H34 900130-87-2 12 779 429 2-METHYL-Z-7-REPOXYHEXADECANE 254 C17H34O 900130-87-2 13 778 407 3LANE, TRICHLORODOCOSYL 442 C22HSCISS 7325-84-0 14 775 493 CIS-2-METHYL-7-OCTADECENE 266 C19H38 35354-39-3 15 774 422 E-7-OCTADECENE 255 C18H38 900130-92-0 16 772 491 3-TRIFLUOROACETOXYPENTADECANE 324 C17H31OZF3 900130-92-0 17 771 306 CYCLOPROPANE, 1-(1-METHYL-F-NONYL-1-M	6								
9 785 436 E-2-HEXADECACEN-1-OL 10 784 437 Z-10-TETRADECEN-1-OL ACETATE 254 C18H3002 900131-10-1 275 426 2-METHYL-27-HEXADECENE 238 C17H34 900130-98-3 12 779 429 2-METHYL-27-BEPOXYHEXADECANE 254 C17H340 900130-87-2 254 C17H340 900130-88-3 13 778 407 SILANE, TRICHLORODOCOSYL- 442 C22H45CI3Si 7325-84-0 14 775 493 CIS-2-METHYL-7-OCTADECENE 266 C19H38 35354-39-3 15 774 422 E7-OCTADECENE 275 491 3-TRIFLUOROACETOXYPENTADECANE 276 C18H30 900130-92-0 177 771 306 CYCLOPROPANE, 1-(1-METHYL-E7-NONYL- 277 771 306 CYCLOPROPANE, 1-(1-METHYL-E7-NONYL- 278 66 C19H38 61888-20-0 18 769 411 2-METHYL-1-OCTADECENE 286 C19H38 61888-20-0 189 769 416 2-METHYL-1-OCTADECENE 286 C19H38 61888-20-0	7	796			490		6971-40-0		
10 784 437 Z-10-TETRADECEN-I-OL ACETATE 254 C18H30O2 900130-99-3 11 781 426 2-METHYL-Z-7-REXADECENE 238 C17H34 900130-87-2 12 779 429 2-METHYL-Z-7, 8-EPOXYHEXADECANE 254 C17H34O 900130-88-3 13 778 407 SILANE, TRICHLORODOCOSYL- 442 C22H45CISS 7325-84-0 14 775 493 CIS-2-METHYL-7-OCTADECENE 268 C19H38 35354-39-3 15 774 422 E-7-OCTADECENE 252 C18H36 900130-92-0 16 772 491 3-TRIFLUOROACETOXYPENTADECANE 324 C17H3102F3 900245-47-8 17 771 306 CVCLOPROPANE, 1-(1-METHYLETHYL)-2-NONYL- 210 C15H30 41977-39-3 18 769 411 2-METHYL-DODEC-3-EN-I-OL 266 C19H38 6188-20-0 19 769 416 -METHYL-DODEC-3-EN-I-OL 198 C13H26O 156992-84-6	8		436	7-TETRADECENE	196	C14H28	10374-74-0		
11 781 426 2-METHYL-Z-7-HEXADECENE 238 C17H34 900130-87-2 12 779 429 2-METHYL-Z-7-BEPOXYHEXADECANE 254 C17H340 900130-88-3 13 778 407 SILANE, TRICHLORODOCOSYL- 442 C22H45CI3Si 7325-84-0 14 775 493 CIS-Z-METHYL-7-OCTADECENE 266 C19H38 35354-39-3 15 774 422 E-7-OCTADECENE 252 C18H36 900130-92-0 16 772 491 3-TRIFLUOROACETOXYPENTADECANE 324 C17H31OZF3 900245-47-8 17 771 306 CVCLOPROPANE, 1-(1-METHYLETHYL)-2-NONYL- 210 C15H30 41977-39-3 18 769 411 2-METHYL-1-OCTADECENE 266 C19H38 6188e-20-0 19 769 416 -4METHYL-1-DODEC-3-EN-1-OL 198 C13H26O 156992-84-6									
12 779 429 2-METHYL-Z-7,8-EPOXYHEXADECANE 254 C17H34O 900130-86-3 13 778 407 SILANE, TRICHLORODOCOSYL- 442 C22H45CISSI 7325-84-0 14 775 493 CIS-2-METHYL-7-OCTADECENE 266 C19H38 35354-39-3 15 774 422 E-7-OCTADECENE 252 C18H36 900130-92-0 16 772 491 3-TRIFLUOROACETOXYPENTADECANE 324 C17H3102F3 900245-47-8 17 771 306 CYCLOPROPANE, 1-(1-METHYLETHYL)-2-NONYL- 210 C15H30 41977-39-3 18 769 411 2-METHYL-1-OCTADECENE 266 C19H38 6188-20-0 9 769 416 4-METHYL-1-DODEC-3-EN-1-OL 198 C13H28O 156992-84-6									
13 778 407 SILANE, TRICHLORODOCOSYL- 442 C22H45CISSi 7325-84-0 14 775 493 CIS-2-METHYL7-OCTADECENE 268 C19H38 3535-39-3 15 774 422 E-7-OCTADECENE 252 C18H36 900130-92-0 16 772 491 3-TRIFLUOROACETOXYPENTADECANE 324 C17H31'OCF3 900245-47-8 17 771 306 CVCLOPROPANE, 1-(1-METHYLETHYL)-2-NONYL- 210 C15H30 41977-39-3 18 769 411 2-METHYL-1-OCTADECENE 266 C19H38 61868-20-0 9 769 416 -METHYL-1-DODEC-3-EN-1-OL 198 C13H28O 156992-84-6									
14 775 493 CIS-2-METHYL-7-OCTADECENE 268 C19H38 35354-39-3 15 774 422 E-7-OCTADECENE 252 C18H36 900130-92-0 16 772 491 3-TRIFLUOROACETOXYPENTADECANE 324 C17H3102F3 900245-47-8 17 771 306 CVCLOPROPANE, 1-(1-METHYLETHYL)-2-NONYL- 210 C15H30 41977-39-3 18 769 411 2-METHYL-1-OCTADECENE 266 C19H38 6186-20-0 19 769 416 4-METHYL-1-DDDEC-3-EN-1-OL 198 C13H26O 156992-84-6						C17H34O			
15 774 422 E-7-OCTADECENE 252 C18H36 900130-92-0 16 772 491 3-TRIFLUOROACETOXYPENTADECANE 324 C17H3102F3 900245-47-8 17 771 306 CVCLOPROPANE, 1-(1-METHYLETHYL)-2-NONYL- 210 C19H30 41977-39-3 18 769 411 2-METHYL-1-OCTADECENE 266 C19H38 61888-20-0 19 769 416 4-METHYL-1-DODEC-3-EN-1-CU 198 C13H28O 156992-84-6									
16 772 491 3-TRIFLUOROACETOXYPENTADECANE 324 C17H31O2F3 900245-47-8 17 771 306 CVCLOPROPANE, 1-(1-METHYLETHYL)-2-NONYL- 210 C15H30 41977-39-3 18 769 411 2-METHYL-1-OCTADECENE 266 C19H38 61868-20-0 19 769 416 4-METHYL-DODEC-3-EN-1-OL 198 C13H26O 156992-84-6									
17 771 306 CYCLOPROPANE, 1-(1-METHYLETHYL)-2-NONYL- 210 C15H30 41977-39-3 18 769 411 2-METHYL-1-OCTADECHNE 266 C19H38 61868-20-0 19 769 416 4-METHYL-DODEC-3-EN-1-OL 198 C13H26O 156992-84-6									
18 769 411 2-METHYL-1-OCTADÉCENE 266 C19H38 61868-20-0 19 769 416 4-METHYL-DODEC-3-EN-1-OL 198 C13H26O 156992-84-6									
19 769 416 4-METHYL-DODEC-3-EN-1-OL 198 C13H26O 156992-84-6		771							
20 769 408 UNDEC-10-YNOIC ACID, DODECYL ESTER 350 C23H42O2 900406-16-5									
	20	769	408	UNDEC-10-YNOIC ACID, DODECYL ESTER	350	C23H42O2	900406-16-5		

FIGURE NO 18:

DISCUSSION:

Gas Chromatography and mass spectroscopy is mainly used to detect and separate the volatile and thermally stable components of drugs. It was done for both Kwatha as well as Arka. In Gas Chromatography of Kwatha, three values of RT i.e., 27.013, 28.414, 32.575 were found that covers the 100% area. These three peaks were followed to run on mass spectroscopy for which 20 chemical compounds were found in each peak. The compound 9-octadecenamide is same for RT 32.575 and RT 27.103 and repeated 8 and 4 times respectively in both the RT. This compound has sedative effect (14). The compound dodecanamide is found which is fatty amide of lauric acid and has a role of metabolite. Octadecanamide is another compound having metabolite role (14). In RT 27.103 Octadecanamide and Dodecanamide were also obtained (14). In RT 28.449, compounds don't have any pharmacological use regarding inflammation. In Gas Chromatography of Arka, six RT values i.e., 10.937, 11.862, 14.919, 16.474, 18.610, 20.271 were found. Four RT values out of six were run for mass spectroscopy because maximum area was covered by these four peaks. For 10.937, 18.619 and 20.271, twenty compounds for each were found. For 16.474. total three compounds were found. Thymol, Dihydrobenzo pyran, Carvacrol, Isothymol and Acetic acid compounds were found which are having Anti- inflammatory actions (14). In probable compound list, the anti-inflammatory action was not found in Kwatha whereas in Arka compounds having anti-inflammatory action were found. Whereas, Anti-inflammatory In-vitro study of Gomutra sadhita triphala kwatha (GTK), Gomutra sadhita triphala Arka (GTA) by protein denaturation inhibition method states that IC 50 value of Kwatha is less that Arka. Denaturation of proteins is the well documented cause of inflammation. The compounds that inhibit the denaturation of protein may be used as anti-inflammatory agents. Lower value of inhibition

CONCLUSION:

concentration shows stronger inhibition of Kwatha than Arka. (15)

Present study reveals that both dosage forms GTK and GTA contains entirely different components. *Arka* contains more volatile compounds having Anti-inflammatory effect. on contrary the *Kwatha* has the components which can exert protein denaturation and thereby act as anti-inflammatory.

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