**SUPPLEMENTARY TABLE:2 BIOACITVE COMPOUNDS TARGETS LIST**

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| S.No. | Bioactive compounds | Targets |
| 1. | 3-Cyclohexene-1-methanol, .alpha.,.alpha.,4-trimethyl-, acetate | HSD11B1, HSD11B2, AR, CYP19A1, F10, CES2, DHCR24, PLA2G1B, PTGES, PTGS2, CYP17A1, TRPM2, PTPN1 |
| 2. | 2(1H)-Isoquinolinecarboximidamide, 3,4-dihydro- | MAOB |
| 3. | Caryophyllene oxide | HSP90AA1 |
| 4. | 4,8-Divinylbicyclo[3.3.1]nonane-2,6-dione | DRD1, NR1I2, DLAT |
| 5. | 12-Hydroxy-14-methyl-oxa-cyclotetradec-6-en-2-one | PKD1, PTGS2, PRKCE |
| 6. | 1,2,3,4-Tetrahydroisoquinolin-6-ol,1-[3-hydroxybenzyl] | ADRB1, ADRB3, DRD1, DRD5, DRD2, DRD3, DRD4, TBXA2R |
| 7. | Ursodeoxycholic acid | AKR1B10, AR, NR1H4, FXR1, HSD11B2, ALB, CYP19A1, F10, ESR2, GPBAR1, GABRA4, NR3C1, PTGFR, SHBG, F3, UGT2B7 |
| 8. | 3-Cyclohexene-1-methanol, .alpha.,.alpha.,4-trimethyl-, acetate+ 12-Hydroxy-14-methyl-oxa-cyclotetradec-6-en-2-one | PTGS2 |
| 9. | 3-Cyclohexene-1-methanol, .alpha.,.alpha.,4-trimethyl-, acetate + Ursodeoxycholic acid | HSD11B2,AR,CYP19A1,F10 |
| 10. | 4,8-Divinylbicyclo[3.3.1]nonane-2,6-dione + 1,2,3,4-Tetrahydroisoquinolin-6-ol,1-[3-hydroxybenzyl] | DRD1 |

In this above table it was found that bioactive compound 3-Cyclohexene-1-methanol, .alpha.,.alpha.,4-trimethyl-, acetate, 2(1H)-Isoquinolinecarboximidamide, 3,4-dihydro, Caryophyllene oxide, 4,8-Divinylbicyclo[3.3.1]nonane-2,6-dione, 12-Hydroxy-14-methyl-oxa-cyclotetradec-6-en-2-one, 1,2,3,4-Tetrahydroisoquinolin-6-ol,1-[3-hydroxybenzyl, Ursodeoxycholic acid have 13, 01, 01, 03, 03, 08, 16 targets respectively. Further we checked for the compound’s having more than one targets and it was found that 3-Cyclohexene-1-methanol, .alpha.,.alpha and 12-Hydroxy-14-methyl-oxa-cyclotetradec-6-en-2-one have one common target 3-Cyclohexene-1-methanol, .alpha.,.alpha.,4-trimethyl and Ursodeoxycholic acid have four common targets. Further we have also looked for more than three targets and four targets but there were no any common targets were found.