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CORRECTION

Correction to: GC-MS Analysis, Molecular Docking and Pharmocokinetic Properties of Phytocompounds from *Solanum torvum* Unripe Fruits and Its Effect on Breast Cancer Target Protein

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The original article has been corrected. The following errors in the Table 1 & 2 are listed below.

The original article can be found online at https://doi.org/10.1007/s12010-021-03698-3.

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Table 1: The phytocompounds identified from GC-MS analysis of aqueous extract of *S. torvum* unripe fruits against breast cancer

S. no.	RT	Compound name	Molecular formula	Molecular weight (g/mol)	% area	PubChem (CID)
1	23.065	Ergost-25-ene-3,6-dione, 5,12-dihydroxy-, (5alpha,12beta)-	$C_{28}H_{44}O_4$	444.6	0.26	91692405
2	17.722	Aspidospermidin-17-Ol, 1-acetyl-16-methoxy-	$C_{22}H_{30}N_2O_3$	370.5	0.45	632854
3	15.286	2-[3,4-Dichlorophenyl]-4-[[2-[1-methyl-2-pyrrolidinyl]ethyl]amino]-6-[trichloromethyl]-S-triazine	$C_{17}H_{18}C_{15}N_5$	467	1.10	558706
4	46.491	Dihydroartemisinin, 10-O-(T-butyloxy)-	$C_{19}H_32O_6$	356.5	0.07	537898
5	43.664	3-[(2-Fluoroanilino)methyl]-5-(2-methoxyphenyl)-1,3,4-oxadiazole-2(3h)-thione	$C_{16}H_{14}FN_3O_2S$	331.4	0.27	578971
6	35.204	2(1H)-naphthalenone, 5-[2-(3-furanyl)ethyl]octahydro-1,5,6,8a-tetramethyl-,	$C_{20}H_{30}O_2$	302.5	0.01	565269
7	43.270	4H-1-Benzopyran-4-one, 2-(3,4-dimethoxyphenyl)-3,5-dihydroxy-7-methoxy-	$C_{18}H_{16}O_7$	344.3	0.31	5748558
8	28.017	2-butenoic acid, 2-methyl-, 4,4a,5,6,7,8,8a,9-octahydro-8a-hydroxy-3,4a,5-trimethylnaphtho[2,3-B]furan-6-yl ester	$C_{20}H_{28}O_4$	332.4	0.14	5367763
9	18.747	Benzonitrile, 4-(4-ethylcyclohexyl)-, trans-	$C_{15}H_{19}N$	213.32	0.27	175307
10	31.404	1-(2-adamantylidene)semicarbazide	$C_{11}H_{17}N_3O$	207.27	0.04	541478

Table 2: The 2D structure of the best identified phytocompounds from GC-MS analysis

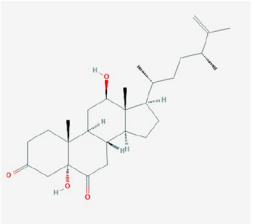
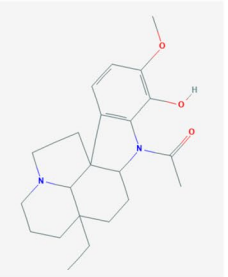
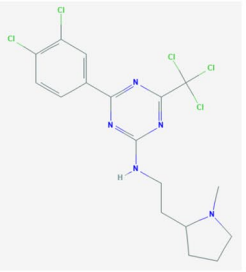
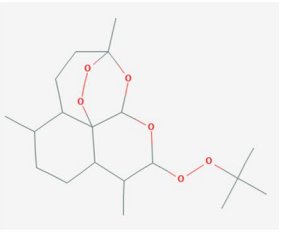
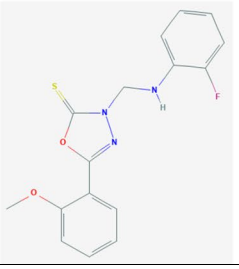
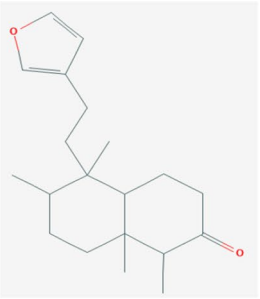
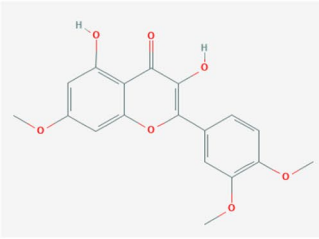
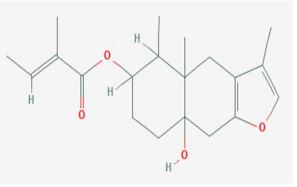
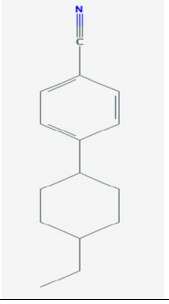
S. No.	PubChem (CID)	Compound Name	2D Structure of Phytocompounds
1	91692405	Ergost-25-Ene-3,6-Dione, 5,12-Dihydroxy-, (5.Alpha.,12.Beta.)-	
2	632854	Aspidospermidin-17-Ol, 1-Acetyl-16-Methoxy-	
3	558706	2-[3,4-Dichlorophenyl]-4-[[2-[1-Methyl-2-Pyrrolidinyl]Ethyl]Amino]-6-[Trichloromethyl]-S-Triazine	
4	537898	Dihydroartemisinin, 10-O-(T-Butyloxy)-	
5	578971	3-[(2-Fluoroanilino)Methyl]-5-(2-Methoxyphenyl)-1,3,4-Oxadiazole-2(3h)-Thione	

Table 2: (continued)

6	565269	2(1H)-Naphthalenone, 5-[2-(3-Furanyl)Ethyl]Octahydro-1,5,6,8a-Tetramethyl-,	
7	5748558	4H-1-Benzopyran-4-one, 2-(3,4-Dimethoxyphenyl)-3,5-Dihydroxy-7-Methoxy-	
8	5367763	2-Butenoic Acid, 2-Methyl-, 4,4a,5,6,7,8,8a,9-Octahydro-8a-Hydroxy-3,4a,5-Trimethylnaphtho[2,3-B]Furan-6-yl Ester	
9	175307	Benzonitrile, 4-(4-Ethylcyclohexyl)-, Trans-	
10	541478	1-(2-Adamantylidene)Semicarbazide