

```

gap> g:= SymmetricGroup( 4 );
Sym( [ 1 .. 4 ] )
gap> tbl:= CharacterTable( g );; HasIrr( tbl );
i5 : betti(t,Weights=>{1,0})
false
      0 1 2 3 4 gap> tblmod2:= CharacterTable( tbl, 2 );
o5 = total: 1 4 13 14 4 BrauerTable( Sym( [ 1 .. 4 ] ), 2 )
      0: 1 . . . .
      1: . 2 2 4 2 gap> tblmod2 = CharacterTable( tbl, 2 );
      2: . 2 5 6 . true
      3: . . 4 . 2
      4: . . . 4 . gap> tblmod2 = BrauerTable( tbl, 2 );
      5: . . 2 . . true
      6: . . . . . gap> tblmod2 = BrauerTable( tbl, 2 );
o5 : BettiTally
i6 : betti(t,Weights=>{0,1})
true
      0 1 2 3 4 gap> libtbl:= CharacterTable( "M" );
o6 = total: 1 4 13 14 4 CharacterTable( "M" )
      0: 1 . . . . gap> CharacterTableRegular( libtbl, 2 );
      1: . 2 2 4 2 BrauerTable( "M", 2 );
      2: . 2 5 6 . BrauerTable( libtbl, 2 );
      3: . . 4 . 2 gap> BrauerTable( libtbl, 2 );
      4: . . . 4 . fail
      5: . . 2 . .
gap> CharacterTable( "Symmetric", 4 );
o6 : BettiTally
i7 : t1 = betti(t,Weights=>{1,1})
CharacterTable( "Sym(4)" )
gap> ComputedBrauerTables( tbl );
      0 1 2 3 4 [ , BrauerTable( Sym( [ 1 .. 4 ] ), 2 ), ]
      0: 1 . . . . option(noprot);
      1: . . . . . timer=1;
      2: . . . . . ring r1 = 32003,(x,y,z),ds;
      3: . 2 . . . int a,b,c,t=11,5,3,0;
      4: . . . . . poly f = x^a+y^b+z^(3*c)+x^(c+2)*y^(c-1)+x^(
      5: . 2 . . . x^(c-2)*y^c*(y^2+t*x)^2;
      6: . . 1 . . option(noprot);
      7: . . 8 6 . timer=1;
      8: . . 4 8 4 ring r2 = 32003,(x,y,z),dp;
      poly f=imap(r1,f);
      ideal j=jacob(f);
      vdim(std(j));
      ==> 536
      vdim(std(j+f));
      ==> 195
      timer=0; // reset timer

o7 : BettiTally
i8 : peek t1

o8 = BettiTally{(0, {0, 0}, 0) => 1 }
      (1, {2, 2}, 4) => 2
      (1, {3, 3}, 6) => 2
      (2, {3, 7}, 10) => 2
      (2, {4, 4}, 8) => 1
      (2, {4, 5}, 9) => 4
      (2, {5, 4}, 9) => 4
      (2, {7, 3}, 10) => 2
      (3, {4, 7}, 11) => 4
      (3, {5, 5}, 10) => 6
      (3, {7, 4}, 11) => 4
      (4, {4, 4}, 8) => 1
      (4, {5, 5}, 9) => 4
      (4, {7, 5}, 12) => 2

```

Journal of Software for Algebra and Geometry

Phylogenetic trees

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ABSTRACT: We introduce the package *PhylogeneticTrees* for *Macaulay2*, which allows users to compute phylogenetic invariants for group-based tree models. We provide some background information on phylogenetic algebraic geometry and show how the package *PhylogeneticTrees* can be used to calculate a generating set for a phylogenetic ideal as well as a lower bound for its dimension. Finally, we show how methods within the package can be used to compute a generating set for the join of any two ideals.

MOTIVATION. A central problem in phylogenetics is to describe the evolutionary history of n species from their aligned DNA sequences. One way to do this is through a model-based approach. A phylogenetic model is a statistical model, specified parametrically, of molecular evolution at a single DNA site. We can regard the aligned sequences as a collection of n -tuples of the four DNA bases, one from each site. Each choice of parameters results in a probability distribution on the 4^n possible n -tuples. The goal of model-based reconstruction is to find a choice of parameters that yields a distribution close to the empirical distribution. If we are able to do so then it is reasonable to assume that the model is an accurate reflection of the underlying evolutionary process. Most significantly, we can infer that the underlying tree parameter of the phylogenetic model is the evolutionary tree of the species under consideration.

MATHEMATICAL BACKGROUND. In phylogenetic algebraic geometry, the statistical models under consideration are tree-based Markov models. This means that we assume a Markov process proceeds along a tree with a transition matrix associated to each edge. A κ -state phylogenetic model on an n -leaf tree \mathcal{T} induces a polynomial map from the parameter space $\Theta_{\mathcal{T}} \subseteq \mathbb{R}^m$ to the probability simplex Δ^{κ^n-1} ,

$$\psi_{\mathcal{T}} : \Theta_{\mathcal{T}} \rightarrow \Delta^{\kappa^n-1} \subset \mathbb{C}^{\kappa^n}.$$

The image of this map is the set of all probability distributions we obtain by varying the entries of the transition matrices; we refer to the image as the model $\mathcal{M}_{\mathcal{T}}$. For phylogenetic applications, usually $\kappa = 2$ or $\kappa = 4$.

The Zariski closure of the model $V_{\mathcal{T}} := \overline{\mathcal{M}_{\mathcal{T}}} \subseteq \mathbb{C}^{\kappa^n}$ is an affine algebraic variety. For the models we consider, the entries of $\psi_{\mathcal{T}}$ are homogeneous polynomials of uniform degree. Thus, $V_{\mathcal{T}}$ can be viewed

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PhylogeneticTrees version 2.0

as the affine cone of a projective variety in \mathbb{P}^{κ^n-1} . The ideal $\mathcal{I}_{\mathcal{T}} := I(V_{\mathcal{T}}) \subseteq \mathbb{C}[x_1, \dots, x_{\kappa^n}]$ of all polynomials vanishing on $V_{\mathcal{T}}$ is a homogeneous ideal called the *ideal of phylogenetic invariants*; this ideal carries useful information about the model that can be used for determining model identifiability and performing model selection [E. S. Allman and Sullivant 2011; Allman et al. 2012; Cavender and Felsenstein 1987; Casanellas and Fernández-Sánchez 2006; A. 1987; Long and Sullivant 2015; Matsen et al. 2008; Matsen and Steel 2007; Rhodes and Sullivant 2012; Rusinko and Hipp 2012].

Given two models, $\mathcal{M}_{\mathcal{T}_1}$ and $\mathcal{M}_{\mathcal{T}_2}$, we define the *mixture model* $\mathcal{M}_{\mathcal{T}_1} * \mathcal{M}_{\mathcal{T}_2}$ to be the image of the map

$$\begin{aligned} \psi_{\mathcal{T}_1, \mathcal{T}_2} : \Theta_{\mathcal{T}_1} \times \Theta_{\mathcal{T}_2} \times [0, 1] &\rightarrow \Delta^{\kappa^n-1} \subset \mathbb{C}^{\kappa^n}; \\ \psi_{\mathcal{T}_1, \mathcal{T}_2}(\theta_1, \theta_2, \pi) &= \pi \psi_{\mathcal{T}_1}(\theta_1) + (1 - \pi) \psi_{\mathcal{T}_2}(\theta_2). \end{aligned} \quad (1)$$

As before, we take the Zariski closure $\overline{\mathcal{M}_{\mathcal{T}_1} * \mathcal{M}_{\mathcal{T}_2}} \subseteq \mathbb{C}^{\kappa^n}$, and now we obtain the algebraic variety $V_{\mathcal{T}_1 * \mathcal{T}_2} = V_{\mathcal{T}_1} * V_{\mathcal{T}_2}$, the join of $V_{\mathcal{T}_1}$ and $V_{\mathcal{T}_2}$. The *join* of two algebraic varieties V and W embedded in a common ambient space is the variety

$$V * W := \overline{\{\lambda v + (1 - \lambda)w \mid \lambda \in \mathbb{C}, v \in V, w \in W\}}.$$

In the special case when $V = W$, the join variety $V * V$ is called the *secant variety of V* . Similarly, given two ideals $I_1, I_2 \subset \mathbb{C}[x_1, \dots, x_n]$, the ideal $I_1 * I_2 \subset \mathbb{C}[x_1, \dots, x_n]$ is the *join ideal*. As for varieties, if $I_1 = I_2$, the ideal $I_1 * I_2$ is the *secant ideal*. We refer to [Sturmfels and Sullivant 2006, Section 2] for the definition of $I_1 * I_2$, but note the following important property:

$$\mathcal{I}_{\mathcal{T}_1} * \mathcal{I}_{\mathcal{T}_2} = I(V_{\mathcal{T}_1} * V_{\mathcal{T}_2}). \quad (2)$$

Our package provides a means of computing invariants for those working in phylogenetic algebraic geometry. We handle a class of commonly used models called *group-based models* that have special restrictions on the entries of the transition matrices. These entries are indexed by elements of a group and thus are subject to the Fourier–Hadamard coordinate transformation, which makes the parametrization monomial and the ideals toric [Evans and Speed 1993; Székely et al. 1993]. We will refer to the original coordinates, which represent leaf probabilities, as *probability coordinates* and the transformed coordinates as *Fourier coordinates*; furthermore, following the literature, we will use p for probability coordinates and q for Fourier coordinates. For these group-based models, we implement a theoretical construction for inductively determining the ideal of phylogenetic invariants for any tree from the invariants for claw trees [Sturmfels and Sullivant 2005]. We also handle the join and secant ideals formed from these ideals, which allows for computations involving mixture models.

FUNCTIONALITY FOR TORIC PHYLOGENETIC VARIETIES. As an example, let \mathcal{T} be the four-leaf tree illustrated in Figure 1 and consider the Cavender–Farris–Neyman (CFN) model, a two-state group-based model, on \mathcal{T} . Then the toric ideal $\mathcal{I}_{\mathcal{T}}$ is generated in degree 2. Using our package, we can compute a generating set for $\mathcal{I}_{\mathcal{T}}$ using two different methods, `phyloToric42` and `phyloToricFP`.

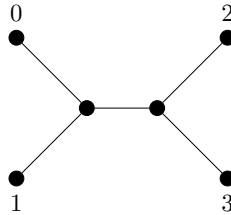


Figure 1. Four leaf tree

The first method `phyloToric42` calls `FourTiTwo.m2`, the [Macaulay2] interface to [4ti2], a software package with functionality for computing generating sets of toric ideals. We input \mathcal{T} by its set of non-trivial splits. In this instance, $01|23$ is the only nontrivial split of \mathcal{T} , which we can enter as either $\{0, 1\}$ or $\{2, 3\}$. The indices on the q 's correspond to two-state labelings of the four leaves of \mathcal{T} :

Macaulay2, version 1.7

```
i1: load "PhylogeneticTrees.m2"
```

```
i2: n = 4; T = {{0,1}}; M = CFNmodel;
```

```
i5: toString phyloToric42(n,T,M)
```

```
o5: = ideal(-q_(0,1,1,0)*q_(1,0,0,1)+q_(0,1,0,1)*q_(1,0,1,0),
           -q_(0,0,1,1)*q_(1,1,0,0)+q_(0,0,0,0)*q_(1,1,1,1))
```

The second method `phyloToricFP` computes generators of $\mathcal{I}_{\mathcal{T}}$ using Theorem 24 in [Sturmfels and Sullivant 2005]; the “FP” in the method name stands for fiber product [Sullivant 2007]. For this example, this theorem allows us to explicitly construct a generating set of $\mathcal{I}_{\mathcal{T}}$ from generators of $\mathcal{I}_{K_{1,3}}$, the ideal associated to the CFN model on the claw tree $K_{1,3}$. While our example here is binary, we note that this method is implemented for all trees, binary or not.

```
i6: = toString phyloToricFP(n,T,M)
```

```
o6: = ideal(-q_(0,0,1,1)*q_(1,1,0,0)+q_(0,0,0,0)*q_(1,1,1,1),
           q_(0,0,1,1)*q_(1,1,0,0)-q_(0,0,0,0)*q_(1,1,1,1),
           q_(0,0,1,1)*q_(1,1,0,0)-q_(0,0,0,0)*q_(1,1,1,1),
           -q_(0,0,1,1)*q_(1,1,0,0)+q_(0,0,0,0)*q_(1,1,1,1),
           -q_(0,1,1,0)*q_(1,0,0,1)+q_(0,1,0,1)*q_(1,0,1,0),
           q_(0,1,1,0)*q_(1,0,0,1)-q_(0,1,0,1)*q_(1,0,1,0),
           q_(0,1,1,0)*q_(1,0,0,1)-q_(0,1,0,1)*q_(1,0,1,0),
           -q_(0,1,1,0)*q_(1,0,0,1)+q_(0,1,0,1)*q_(1,0,1,0))
```

The algorithm used by `phyloToricFP` returns more polynomials than are required to generate the ideal. If we wish to directly compare this ideal to that returned by `phyloToric42` we must reconstruct both ideals in the same ring. Thus, we use the function `qRing` to define the ring of Fourier coordinates and use the option of specifying the ring for our ideals:

```
i7: R = qRing(n,M)
```

```
i8: = phyloToric42(n,T,M,QRing=>R) == phyloToricFP(n,T,M,QRing=>R)
```

```
o8: = true
```

In our experiments, for most cases, `phyloToric42` runs much faster than `phyloToricFP`. This is likely because we have implemented a naive version of the toric fiber product algorithm from [Sturmfels and Sullivant 2005] with no attempt to avoid producing redundant polynomials. It would be worth investigating if there is a faster implementation of this algorithm. Still, one advantage offered by the fiber product method is the ability to inductively construct a single invariant when computing the entire ideal is infeasible. The method `phyloToricRandom` returns such a randomly constructed invariant.

The polynomials that are returned by both methods are in Fourier coordinates, however, they can be converted to probability coordinates using the function `fourierToProbability`. To do so, we must first construct the ring of probability coordinates using `pRing`. Then the method `fourierToProbability` returns a ring map that converts polynomials in Fourier coordinates to probability coordinates:

```
i9: = S = pRing(n,M);
i10: = phi = fourierToProbability(S,R,4,M);
i11: = f = (vars R)_(0,0)
o11: = q_(0,0,0,0)
i12: = phi(f)
o12: = (1/2)*p_(0,0,0,0)+(1/2)*p_(0,0,0,1)+(1/2)*p_(0,0,1,0)+(1/2)*p_(0,0,1,1)
      +(1/2)*p_(0,1,0,0)+(1/2)*p_(0,1,0,1)+(1/2)*p_(0,1,1,0)+(1/2)*p_(0,1,1,1)
      +(1/2)*p_(1,0,0,0)+(1/2)*p_(1,0,0,1)+(1/2)*p_(1,0,1,0)+(1/2)*p_(1,0,1,1)
      +(1/2)*p_(1,1,0,0)+(1/2)*p_(1,1,0,1)+(1/2)*p_(1,1,1,0)+(1/2)*p_(1,1,1,1)
```

FUNCTIONALITY FOR SECANT VARIETIES. Mixtures of group-based phylogenetic models correspond to secants and joins of toric ideals, objects that are of interest in combinatorial commutative algebra, but are notoriously hard to compute. In the methods `joinIdeal` and `secant`, we implement the elimination method described in [Sturmfels and Sullivant 2006, Section 2] for computing the join of two homogeneous ideals or the secant of one homogeneous ideal.

Consider now the Jukes–Cantor model on \mathcal{T} from Figure 1. The phylogenetic ideal for the mixture of $\mathcal{M}_{\mathcal{T}}$ with itself is the second secant ideal of the homogeneous ideal $\mathcal{I}_{\mathcal{T}}$, denoted $\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}}$. For secants, the method `secant` takes as input a homogenous ideal and an integer k and returns a generating set for the k -th secant ideal. The method also accepts the optional argument `DegreeLimit=>{1}`, which computes generators of the ideal only up to degree l . Thus, we can obtain generators of degree 3 or less of $\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}}$ with the following commands. The minimal generating set of `SecI3` contains 49 linear invariants; we only print the generators with degree greater than one:

```
i13: = I = phyloToric42(n,T,JCmodel);
i14: SecI3 = secant(I,2,DegreeLimit={3});
i15: toString for i in flatten entries mingens SecI3
      list (if (degree i)#0 == 1 then continue; i)
o15 = {q_(0,3,3,0)*q_(3,0,2,1)*q_(3,2,0,1)-q_(0,3,2,1)*q_(3,0,3,0)*q_(3,2,0,1)
      +q_(0,3,2,1)*q_(3,0,0,3)*q_(3,2,1,0)-q_(0,3,0,3)*q_(3,0,2,1)*q_(3,2,1,0)
      -q_(0,3,3,0)*q_(3,0,0,3)*q_(3,2,3,2)+q_(0,3,0,3)*q_(3,0,3,0)*q_(3,2,3,2)}
```

The degree bound allows for the possibility of obtaining some invariants when computing a generating set for the secant ideal is infeasible. Let $(\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}})_l$ be the ideal generated by the elements of $\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}}$ of degree less than or equal to l . In some instances $(\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}})_l$ may be equal to $\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}}$. To prove this, we must verify that $\dim((\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}})_l) = \dim(\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}})$ and that $(\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}})_l$ is prime. Assuming we are able to compute $(\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}})_l$, we can compute its dimension and verify that it is prime. We then know that $\dim((\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}})_l) \geq \dim(\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}})$, leaving the inequality $\dim((\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}})_l) \leq \dim(\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}})$ to show. The method `toricSecantDim` enables us to do this using a probabilistic method based on Terracini’s lemma [1911] to compute a lower bound on $\dim(\mathcal{I}_{\mathcal{T}} * \mathcal{I}_{\mathcal{T}})$.

Using this method, we can show that the secant of the ideal from the previous example is in fact generated in degree less than three:

```
i16: = dim(SecI3)
o16: = 12
i17: = isPrime(SecI3)
o17: = true
i18: = toricSecantDim(phyloToricAMatrix(4,{{0,1}},JCmodel),2)
o18: = 12
```

In the code above, we used `phyloToricAMatrix(n,T,JCmodel)` to construct the defining integral matrix of the toric ideal. For more details, see the documentation and [Sturmfels 1996]. In this instance, the method outlined is substantially faster than using the secant method without a degree bound.

ADDITIONAL FUNCTIONALITY. Although this package was developed with toric ideals from phylogenetics in mind, the methods `secant` and `joinIdeal` can be used for *any* homogeneous ideals. Thus, these can be employed for computations outside of phylogenetic algebraic geometry.

The following models are loaded with the package: the Cavender–Farris–Neyman model, the Jukes–Cantor model, the Kimura 2-parameter model, and the Kimura 3-parameter model. Additionally, some functionality for working with trees is available in this package, which includes the methods `edgeCut`, `vertexCut`, `edgeContract`, `internalEdges`, `internalVertices`.

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SUPPLEMENT. The online supplement contains version 2.0 of `PhylogeneticTrees`.

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