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## F-Polynomials and Newton Polytopes

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# F-POLYNOMIALS AND NEWTON POLYTOPES 

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#### Abstract

We provide an effective algorithmic method for computation of Gross-Keel-Hacking-Kontsevich potential, F-polynomials and Bernstein-Kazhdan decoration function and it's complexity bounds. For simply laced Lie algebras we make conjecture and provide experimental evidence that Newton polytopes for Gross-Keel-Hacking-Kontsevich potential do not contain any interior lattice points.


## 1. F-polynomials and Gross- <br> Hacking-Keel-Kontsevich POTENTIALS

Let $G$ be a group with the Lie algebra of simply-laced type, $B_{+}$and $B_{-}$be its Borel subgroups, with the set of simple roots $\alpha_{a}, a \in I, W$ the Weyl group. The Gross-Hacking-Keel-Kontsevich potential (GHKK for short) $W_{G H K K}$ is a function on the double Bruhat cell $G^{w_{0}, e}=B_{-} \cap B_{+} \overline{w_{0}} B_{+}$, defined using cluster algebra's tools [6]. Because of validity the Fock-Goncharov conjecture in such cases [5], we get the polyhedral parametrization of canonical bases of the ring of regular functions on $G / B$ arising from the tropicalizations of the potential.

Specifically, the ring of regular functions on the double Bruhat cell is endowed with the cluster algebra structure. Namely, for a reduced decomposition $\mathbf{i}$ of the longest element $w_{0} \in W$ with length $N$, let $\Sigma_{\mathbf{i}}$ be a corresponding X-cluster seed and $Q_{\mathbf{i}}$ be the corresponding quiver (due to [1]). Then $W_{G H K K}$ is a polynomial in the cluster variables $\Sigma_{\mathbf{i}}([10])$.

Namely, a seed $\Sigma$ with underlying quiver $Q$ is optimal for a frozen vertex $a \in I$ if after deleting arrows between frozen vertices and $a$, the vertex $a$ becomes a source of the quiver $Q$.

For the optimal seed $\Sigma$, the $a$ th part of the GHKK-potential is equal to the value of the corresponding frozen,

$$
\begin{equation*}
W_{a}=Y_{a} \tag{1}
\end{equation*}
$$

For a frozen $a \in I$, there exists an appropriate reduced word $\mathbf{i}^{\prime}$, such that $\Sigma_{\mathbf{i}^{\prime}}$ is an optimal seed for $a$.

Because of that, for a given reduced decomposition i, one can compute the half of the GHKK-potential

$$
W_{G H K K}=\sum_{a \in I} W_{a}
$$

using cluster mutations corresponding to 3 -braid moves between the reduced decompositions of $w_{0}$ (for $l$ and $k$, such that $\left.a_{l k}=-1, s_{k} s_{l} s_{k}=s_{l} s_{k} s_{l}\right)$. Namely, for computing $W_{a}$, we apply a sequence of cluster mutations corresponding to 3braid moves which transform $\Sigma_{\mathbf{i}}$ into an

[^0]optimal seeds for $a$, then $W_{a}$ is the $X$ cluster variable at the frozen vertex labeled by $a$ in the optimal seed computed in the variables of the seed $\Sigma_{\mathbf{i}}$. In variables of the seed $\Sigma_{\mathbf{i}}$, such an $X$-cluster variable is equal to the specification of the $F$-polynomial (see 4, 9) and takes the form
\[

$$
\begin{align*}
W_{a}= & Y_{1}^{c_{1 a}(t)} \cdots Y_{N}^{c_{N a}(t)} \\
& \prod_{i} F_{i}(t)\left(Y_{1}, \cdots, Y_{N}\right)^{b_{i a}(t)} . \tag{2}
\end{align*}
$$
\]

In the above formula we take notations of [9], where $t$ means the end vertex of the path in the mutation graph from the optimal seed for $a$ to $\Sigma_{\mathbf{i}}$ and $Y_{j}$ 's are cluster variables of $\Sigma_{\mathbf{i}}$.

Precisely (see [10]) $W_{a}$ is of the form product of the frozen $Y_{a}(t)$ and an $F$ polynomial.

## 2. Newton polytopes

We are interested of properties of the Newton polytopes of the individual terms $W_{a}, a \in I$, of the half-potential $W_{G H H K}$, as well as the Newton polytope of $W_{G H H K}$.

Fei in [3] conjectured that the Newton polytope of an $F$-polynomial has no interior integer points.

For minuscule weight $a \in I$, the validity of this conjecture for $F$-polynomials corresponding to terms $W_{a}$ follows from Remark 5.17 [7]. Namely in such a case, the Newton polytope is a geometric realisation of a distributive lattice of the corresponding decorated graph $D G_{a}$. Since such a polytope is a convex hull of a subset of the vertices of a unit cube the claim follows.

An integer polytope without interior integer points is void.

We state the following
Conjecture. For a simply-laced group $G$, and any reduced decomposition i of $w_{0}$, the Newton polytope of $W_{G H K K}$ is void.

Note that validity of this conjecture implies that the Newton polytopes of $F$ polynomials for frozens are void.

We made computer verification of Conjecture for the following cases
$A_{n}, n=3,4,5,6,7, D_{n}, n=4,5,6,7$, $E_{6}, E_{7}$

Note that the cluster computation of $W_{G H K K}$ is time consuming, because the division of Laurent polynomials in many variables is time consuming.

We use another approach. Namely, because of Theorem 1 in [5], we compute $W_{G H K K}$ by applying the algorithm of [8] for computing the Berenstein-Kazhdan decoration function $\Phi_{B K}$.

Theorem 1. For simply-laced $G$, and $a$ given reduced decomposition $\mathbf{i}$, the Newton polytopes $\Phi_{B K}$ and $W_{G H K K}$ are isomorphic under a unimodular transformation.

Corollary 1. The Newton polytopes $\Phi_{B K}$ is void if and only if the Newton polytopes $W_{G H K K}$ is void.

Thus, for the numeric verification of Conjecture we compute the Newton polytope $\Phi_{B K}$ using the algorithm [8] and Polymake.

## 3. Algorithm Description

The algorithm for computation Berenstein-Kazhdan decoration function $\Phi_{B K}$ is based on Theorem 4.4 in [8]. For the input data consisting of a group $G$ and reduced word $w_{0}=\mathbf{i}$. Let $W$ be a corresponding Weyl group, $I$ be the set of simple roots, $\Lambda_{i}$ be $i$-th fundamental weight.

We compute $\Phi_{B K}$ as sum:

$$
\Phi_{B K}=\sum_{j \in I} \Delta_{w_{0} \Lambda_{j}, s_{j} \Lambda_{j}},
$$

where $\Delta$ is generalized minor function (see Defenition 2.2 in [8]). By Theorem 4.4. [8] it is possible to compute all monomials in $\Delta_{w_{0} \Lambda_{i}, s_{i} \Lambda_{i}}$ by

```
set initial algebraic structures:
    Weyl group,
    Cartan matrix,
    ring of Laurent polynomials
set initial value of word w
compute source monomial
    compute weyl_action(i}\mp@subsup{\mathbf{i}}{1}{},\mp@subsup{\mathbf{i}}{2}{},\ldots\mp@subsup{\mathbf{i}}{j}{})(\mp@subsup{\alpha}{j}{})\mathrm{ for all j
    find index }h(j)\mathrm{ of one equal to exactly }\mp@subsup{a}{j}{
    set th(j) as starting monomial
start graph enumeration
set new variables
    St=[[source monomial]] //layers of associahedron
    Gs=[] //graph structure
until new_nodes=[1]
    //until we get only one new monomial equal to 1)
    mutate all monomials in last layer of St
    add new monomials to new layer of St
    add arrows to Gs for all mutations
        with marks corresponding to }\mp@subsup{A}{k}{-1}\mathrm{ factors
return sum(flatten(St))
```


## Listing 1: main loop of the algorithm

consequently applying multiplication by monomials $A_{k}^{-1}$, where

$$
A_{k}=t_{j} t_{j^{+}} \prod_{j<l<j^{+}} t^{c_{i_{l}, i_{j}}}
$$

and
$k+:=\min \left\{l \in[1, N] \mid w_{l}=w_{k}, l>k\right\} \cup\{\infty\}$
$k-:=\max \left\{l \in[1, N] \mid w_{l}=w_{k}, l<k\right\} \cup\{0\}$
(3.7 [8]) starting from predefined source monomial. The algorithm performs graph enumeration for graph with vertices being monomials of $\Delta_{w_{0} \Lambda_{j}, s_{j} \Lambda_{j}}$ and edges are relations of monomials being different by multiplication by $A_{k}^{-1}$.

Computation of each $\Delta_{w_{0} \Lambda_{i}, s_{i} \Lambda_{i}}$ in pseudo-code using SageMath computer algebras system is shown in listing 1.

The mutation procedure in listing 2 checks which mutations are applicable to a given monomial due to Theorem 4.4 [8 and computes new monomials. To speed up computation we compute
the b-vector $b^{0}$ of the source monomial $t_{h(i)}$ and then we apply Lemma 5.2 [8]. Computed $b$-vectors are stored for later use. For +-notation see (3.5) [8]. The pseudo-code for $b$-vectors implementation with reference to SageMath functions is shown in listing 3 .

After $\Delta_{w_{0} \Lambda_{i}, s_{i} \Lambda_{i}}$ are computed for all simple weights all is left is sum them to obtain $\Phi_{B K}$. Computation of it's Newton polytope using Polymake is straightforward. We check it's interior integer lattice points using Polymake method N_INTERIOR_LATTICE_POINTS.

For computation of monomials in $W_{G H K K}$ it is necessary to compute each set of monomials $W_{a}$ from $\Phi_{B K}$ computation data. Using

$$
\begin{aligned}
W_{a}= & Y_{1}^{c_{1 a}(t)} \cdots Y_{N}^{c_{N a}(t)} \\
& \prod_{i} F_{i}(t)\left(Y_{1}, \cdots, Y_{N}\right)^{b_{i a}(t)} .
\end{aligned}
$$

```
def mutate_node(M)
```



```
    compute bj for node if it was not computed previously
    set initial result as empty list
    for all simple roots }\mp@subsup{\alpha}{l}{
        for all k in Rl
            if k+<\infty
                if }\mp@subsup{d}{k}{}<2\mathrm{ and }\mp@subsup{b}{k+}{}>
                    new_monomial = M* A - = 
                    compute new_monomial }\mp@subsup{b}{j}{}\mathrm{ :
                            Lemma 5.2 [8
                                new_monomial }\mp@subsup{b}{j}{}=\mathrm{ old_monomial }\mp@subsup{b}{j}{
                                new_monomial bk +=1
                                new_monomial b}\mp@subsup{b}{k+}{}-=
                    cache new_monomial bj
                    add new_monomial and graph edge to result
                if }\mp@subsup{d}{k}{}==\mp@subsup{d}{k+}{
                    set lookup depth h=2
                        while }\mp@subsup{k}{}{+h}<\infty\mathrm{ and }\mp@subsup{d}{\mp@subsup{k}{}{+h}}{}=0\mathrm{ and }\mp@subsup{b}{\mp@subsup{k}{}{h+}}{}=
                    h++
                    if d}\mp@subsup{d}{\mp@subsup{k}{}{+h}}{}=-1\mathrm{ and }\mp@subsup{b}{\mp@subsup{k}{}{+h}}{}=
                    new_monomial = M* A
                    compute new_monomial }\mp@subsup{b}{i}{}\mathrm{ :
                            [Lemma 3.4]
                            new_monomial }\mp@subsup{b}{j}{}=\mathrm{ old_monomial }\mp@subsup{b}{j}{
                    new_monomial }\mp@subsup{b}{k}{}+=
                    new_monomial b}\mp@subsup{b}{k+}{}-=
                    cache new_monomial }\mp@subsup{b}{j}{
                    add new_monomial and graph edge to result
    return result
```

Listing 2: mutation procedure

```
def b(M)
    //M= \prod m=1 吕 m
    set b}\mp@subsup{b}{N}{}=\mp@subsup{d}{N}{}
            fundamental_weights[j].weyl_action( }\mp@subsup{\alpha}{j}{}\mathrm{ ).
                scalar(simple_coroots(i}\mp@subsup{|}{N}{})
    for t from N-1 to 1
        b}\mp@subsup{b}{t}{}=\mp@subsup{d}{t}{}
            fundamental_weights[j].weyl_action( }\mp@subsup{\alpha}{j}{}\mathrm{ ).
                        scalar(simple_coroots(it))
        for l from t to N
        bt -= b[l]*a\mp@subsup{a}{\mp@subsup{\mathbf{i}}{t}{},\mp@subsup{\mathbf{i}}{l+1}{}}{}
    return b
```

Listing 3: b-vector computation

```
compute GHKK support
    b_start = get b}\mp@subsup{b}{i}{}\mathrm{ vector of source monomial in }\mp@subsup{\Delta}{\mp@subsup{w}{0}{}\mp@subsup{\Lambda}{i}{},\mp@subsup{s}{i}{}\mp@subsup{\Lambda}{i}{}}{
    b_stop = get b}\mp@subsup{b}{i}{}\mathrm{ vector of stop monomial in }\mp@subsup{\Delta}{\mp@subsup{w}{0}{}\mp@subsup{\Lambda}{i}{},\mp@subsup{s}{i}{}\mp@subsup{\Lambda}{i}{}}{
    ee = basis spanned by }\mp@subsup{e}{k}{}-\mp@subsup{e}{k+}{
    GHKK_support = coordinates of b_stop-b_start in ee
compute Y_frozen
    Y_frozen=Y[number last occurrence of j in i]
compute }\mp@subsup{W}{j}{
    compute start monomial
        Y_start = Y_frozen * }\mp@subsup{\prod}{0<k<=len(i)}{}\mp@subsup{Y}{k}{
    W = dictionary with keys in monomials in }\mp@subsup{\Delta}{\mp@subsup{w}{0}{}\mp@subsup{\Lambda}{i}{},\mp@subsup{s}{i}{}\mp@subsup{\Lambda}{i}{}}{
    W[start monomial of }\mp@subsup{\Delta}{\mp@subsup{w}{0}{}\mp@subsup{\Lambda}{i}{},\mp@subsup{s}{i}{}\mp@subsup{\Lambda}{i}{\prime}}{}]=\mathrm{ Y_start
    enumerate edges in graph Gs starting from source monomial:
        v_b = start of the edge
        v_e = end of the edge
        k = mark on the edge
        W[v_e]=W[e_b]*Y[k]
    return set(values(W))
```


## Listing 4: computation of $W_{G H K K}$ monomials

we Because in $A$-cluster variables $t_{m}=$ $\frac{X_{m}}{X_{m^{-}}}$([2] Corollary 4.10). Recall, that cluster variables are generalized minors computed at the twisted matrix corresponding to $\left.\theta_{\mathbf{i}}^{-}\left(t_{1}, \cdots, t_{N}\right)\right)$. Namely, at the vertex $m$

$$
X_{m}:=\Delta_{\Lambda_{i_{m}}, s_{N} \cdots s_{m+1} \Lambda_{i_{m}}}\left(\psi^{w_{0}, e}(X)\right)
$$

where $X=\theta_{\mathbf{i}}^{-}\left(t_{1}, \cdots, t_{N}\right)$, and $\psi^{w_{0}, e}(X)$ is the twist, $\psi^{w_{0}, e}(X) \in U^{e, w_{0}}=U \cap$ $B_{-} \bar{w}_{0} B_{-}$. Because of that relation to $A$ cluster variables $X$ 's, we get by direct computations, using $p$-map, the following relations to $X$-cluster variables, $Y$ 's

$$
\begin{gathered}
A_{j}=t_{j} t_{j^{+}} \prod_{j<l<j^{+}} t_{l}^{a_{i_{l}, i_{j}}}= \\
\frac{X_{j}}{X_{j^{-}}} \frac{X_{j^{+}}}{X_{j}} \prod_{j<l<j^{+}}\left(\frac{X_{l}}{X_{l^{-}}}\right)^{a_{i_{l}, i_{j}}} \\
\quad=\frac{\prod_{l^{\prime} \rightarrow j} X_{l^{\prime}}}{\prod_{j \rightarrow l^{\prime \prime}} X_{l^{\prime \prime}}}=Y_{j} .
\end{gathered}
$$

Similarly, we get
$t^{a^{a_{J+1}, i}} t^{a_{i_{N, i}}} Y^{c_{1 a}(t)} Y^{c_{N a}(t)} \quad$ edges in Gs graph starting from frozen


To determine starting monomial of $F$ polynomial part it is needed to compute GHKK support - coordinates of $b_{\text {source }}-b_{\text {stop }}$ vector in basis $e_{k}-e_{k+}$ ((3.5) [8]), where $b_{\text {source }}$ is $b$-vector of source monomial and $b_{\text {stop }}$ is $b$-vector of last computed monomial in $\Delta_{w_{0} \Lambda_{i}, s_{i} \Lambda_{i}}$ (listing 4).

Last step is to compute monomials in $W_{G H K K}$ as $\bigcup_{j \in I} W_{j}$. Again, checking if Newton polytope of $W_{G H K K}$ is void is done via Polymake using same routine.

The implementation using SageMath is provided in [13].

## 4. Algorithm complexity

We establish a bound on the complexity of the algorithm computing $\Delta_{w_{0} \Lambda_{i}, s_{i} \Lambda_{i}}$ and $W_{j}$ with respect to the number of monomials in $\Delta_{w_{0} \Lambda_{i}, s_{i} \Lambda_{i}}$. Let $K$ be the number of such monomials, and $r$ be the rank of the Lie algebra.

In the mutation function, each node can produce at most $N=\operatorname{length}(\mathbf{i})$ mutations, $N \sim O\left(r^{2}\right)$. Note that mutation branch with lookup in i need no more than $r$ cycles and $k+/ k$ - operations can also be computed in $r$ operations if $\mathbf{i}$ splitting into slices of same root number is cached. With monomial multiplication complexity $(O(N))$, we get $O\left(r^{4} K\right)$.

The main tree search can be done in linear time using prefix trees for string representation of monomials in alphabet $t_{j}$ is of complexity $O\left(r^{2} K\right)$, because monomials can have only $N$ variables. Therefore total complexity of generating the monomials of $\Delta_{w_{0} \Lambda_{i}, s_{i} \Lambda_{i}}$ is of complexity

$$
O\left(r^{4} K\right) \sim O\left(r^{2} \star\right.
$$

length of string representation).
$W_{j}$ computation is bounded by multiplication complexity and number of edges in Gs: $O\left(r^{2}\right) * O\left(K * r^{2}\right) \sim O\left(r^{4} K\right)$. For a fixed $r$, this complexity is the lowest possible complexity being linear with respect to actual complexity to print out
the answer. Complexity of lattice point counting in polytopes has theoretical exponential upper bound with respect of number of inequalities defining polytope (11).

Actual computing speed is mostly determined by speed of Polymake operations. Average time of $\Delta_{w_{0} \Lambda_{i}, s_{i} \Lambda_{i}}$ and $W_{j}$ computation for single simple root of $D_{6}$ with polytope checking is around 2 seconds, for $E_{7}-8$ seconds. For comparison, one computation for single simple root of $D_{6}$ without any Polymake operations takes around 70 ms . For all computations we used a PC with dual 3.8 Ghz Intel ${ }^{\circledR}$ Xeon ${ }^{\circledR}$ Gold 5222 CPU running Ubuntu Linux.

Computing all half-potentials and $W_{G H K K}$ for given Dynkin type relies on enumeration of all reduced decompositions of longest element in Weyl group. This is done using SageMath library with BraidOrbit [12] function. Original BraidOrbit function provides full list of reduced decompositions, which can be enormous and not feasible to compute in real applications (even for $D_{5}$ ). This can be resolved making it generator function, returning a new reduced decomposition each call. Second difficulty is related to algorithm computing all reduced decompositions. Function loops trough all braid moves and applies them to given reduced decomposition, then compares new decompositions to ones that were produced and stored before. This leads to extensive memory consumption when computing all halfpotentials for given Lie group. As can be seen from algorithm, if two reduced decompositions can be transformed one into another using only 2-braid moves, then the results of computation differ only by variable exchange. If it is possible to enumerate all classes of reduced decompositions up to equivalence by 2 braid moves, it can drastically reduce number of cases for computation (cf. for
$D_{4}$ number of classes is 182 and total number of reduced words is 2316, for $D_{5}-13198$ and 12985968). (Number of classes of equivalence fot type $D$ can be found in A180607 OEIS sequence.) Thus we propose an open problem to find fast algorithm to produce only relevant reduced decompositions which give different half-potentials.

Problem. Find algorithm to enumerate quotient by 2-braid moves equivalence set of set of all reduced decompositions of longest element of Weyl group and produce one reduced decomposition from each class with respect of equivalence by far commuting braid moves. Complexity of one step of generation should be polynomial with respect of rank of Lie-group.

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