

From the Weak Bruhat

Order to Crystal Posets

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Perspective & Main Goal:

- Study crystal graphs regarded as posets via poset map to weak Bruhat order, namely via the (right) key map.

Poset Structure for Many Crystals

$$u <_{\text{crystal}} v \iff u \xrightarrow{f_i} v \text{ for some } i$$

- Transfer properties of weak Bruhat order to crystals.

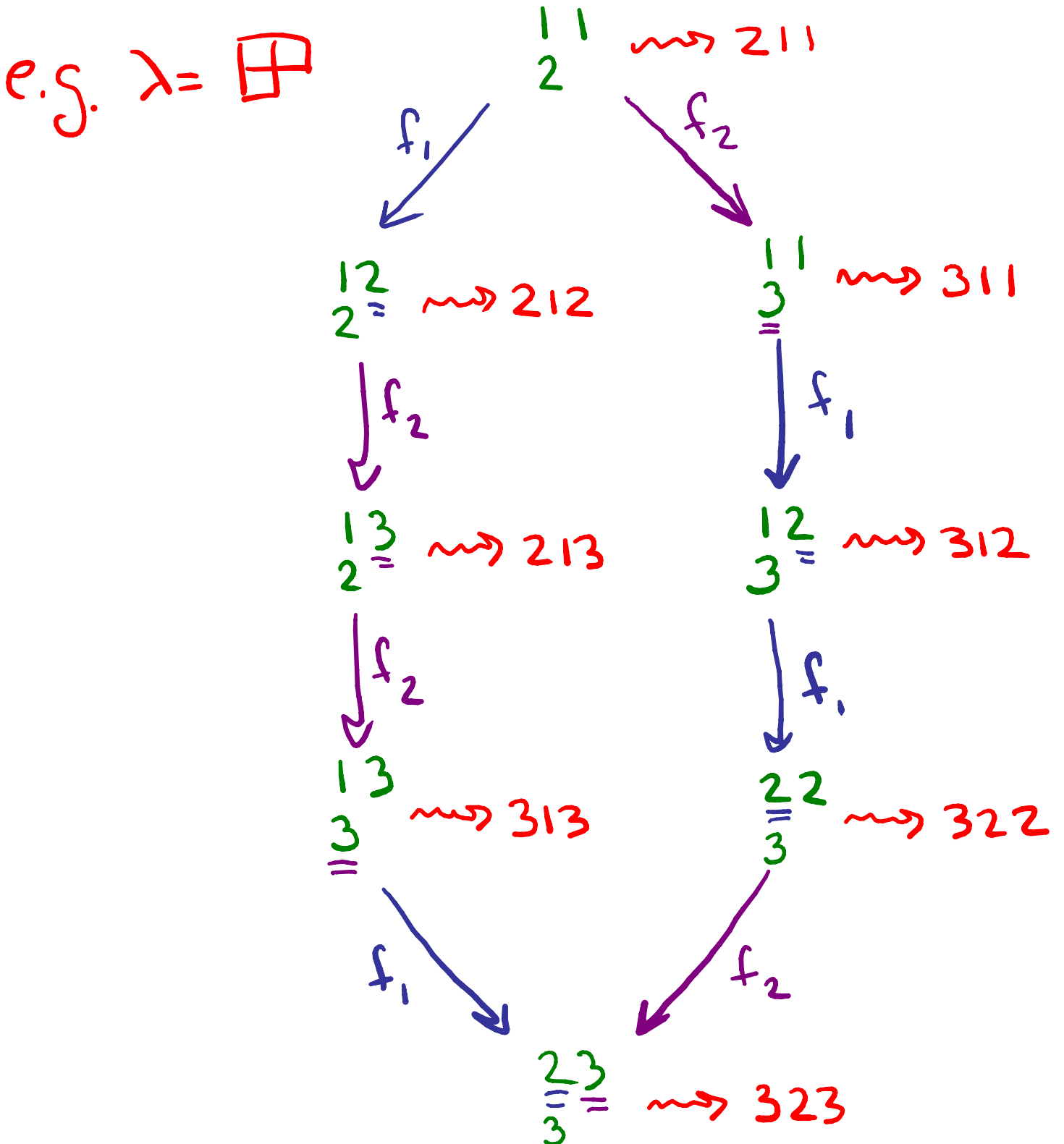
Weak Bruhat order:

$$u <_{\text{weak}} s_i u \text{ if } l(s_i u) > l(u)$$

Motivations for Crystals

- Study representation theory of Kac-Moody algebras (e.g. affine Lie algebras)
- Take universal enveloping algebra, \neq its quantum algebra w/ parameter q
- $q \rightsquigarrow 1$ yields $U(A)$ for Kac-Moody algebra A
- $q \rightsquigarrow 0$ yields algebra with same dimensions of weight spaces, described by combinatorics of "crystal graphs"

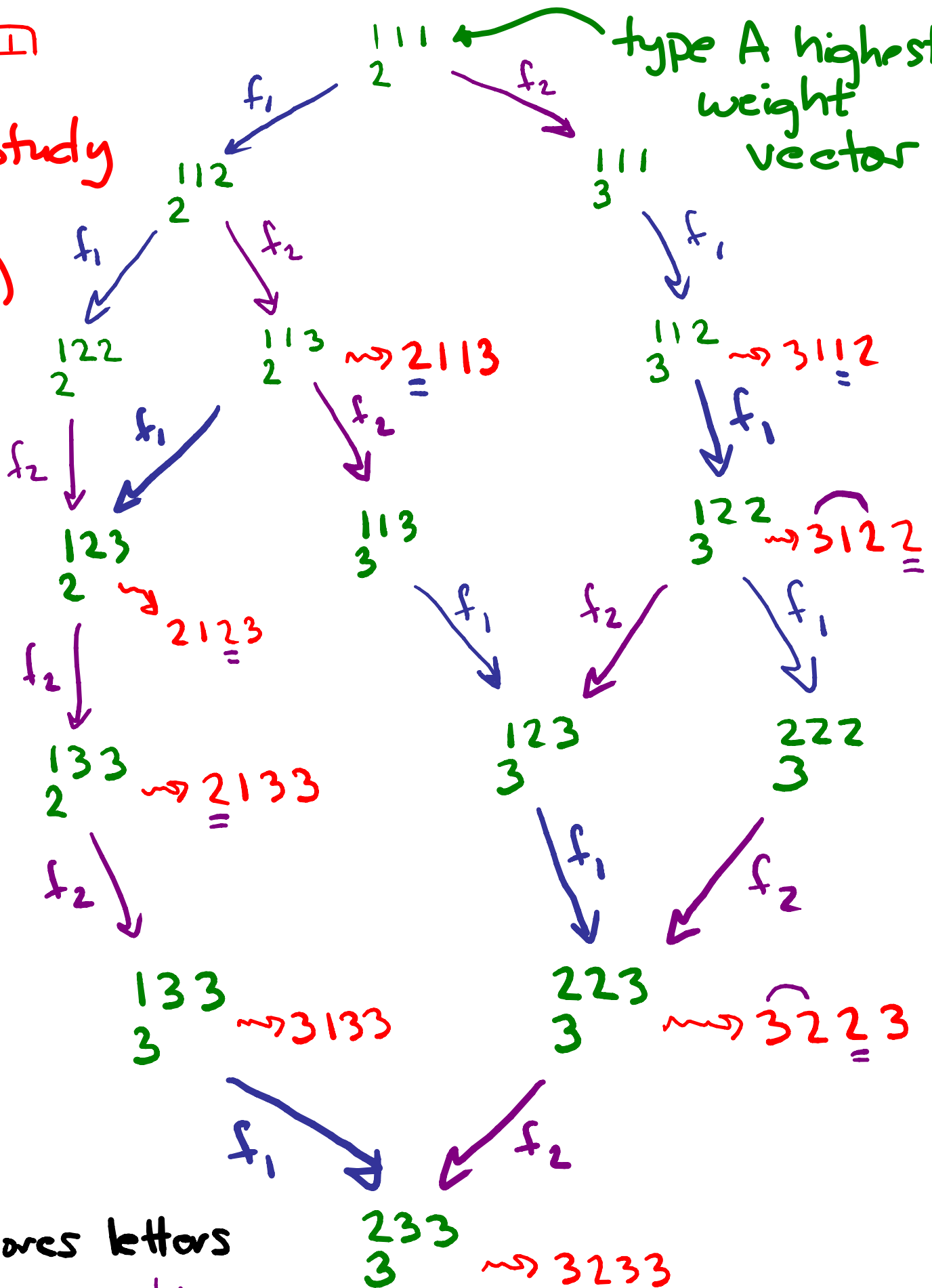
(Type A) Crystals of Highest Weight Representations & their Kashiwara Lowering Operators



$$\lambda = \begin{array}{|c|c|} \hline & \square \\ \hline \square & \square \\ \hline \end{array}$$

(We study dual poset)

type A highest weight vector



f_i ignores letters other than i & $i+1$,

pairs $i+1$ followed by i , then $f_i: i^r (i+1)^s \rightarrow i^{r-1} (i+1)^{s+1}$

Talk Outline:

I. Background Review

II. New Algorithm to Calculate

Right Key of Crystal

- does not depend on choice of model

III. Positive Results for Lower Intervals $[\hat{0}, u]$

- Möbius function \neq homotopy type
- Connectedness of saturated chains under "Stembridge moves" \neq "Sternberg moves"

IV. Negative Results for Arbitrary Type A Intervals $[u, v]$

- Arbitrarily large Möbius functions
- Arbitrarily high degree non-redundant "relations" amongst crystal operators

I. Background

Def'n: The (left) weak Bruhat order

on Coxeter system (W, S) is the partial order with cover relations

$u < \cdot v \iff v = s_i u$ for $u, v \in W$ with

$l(v) > l(u)$, for $l(v) = \min \{ r \mid v = s_{i_1} \dots s_{i_r} \}$
for $s_{i_1}, \dots, s_{i_r} \in S$

e.g. $W = S_n$

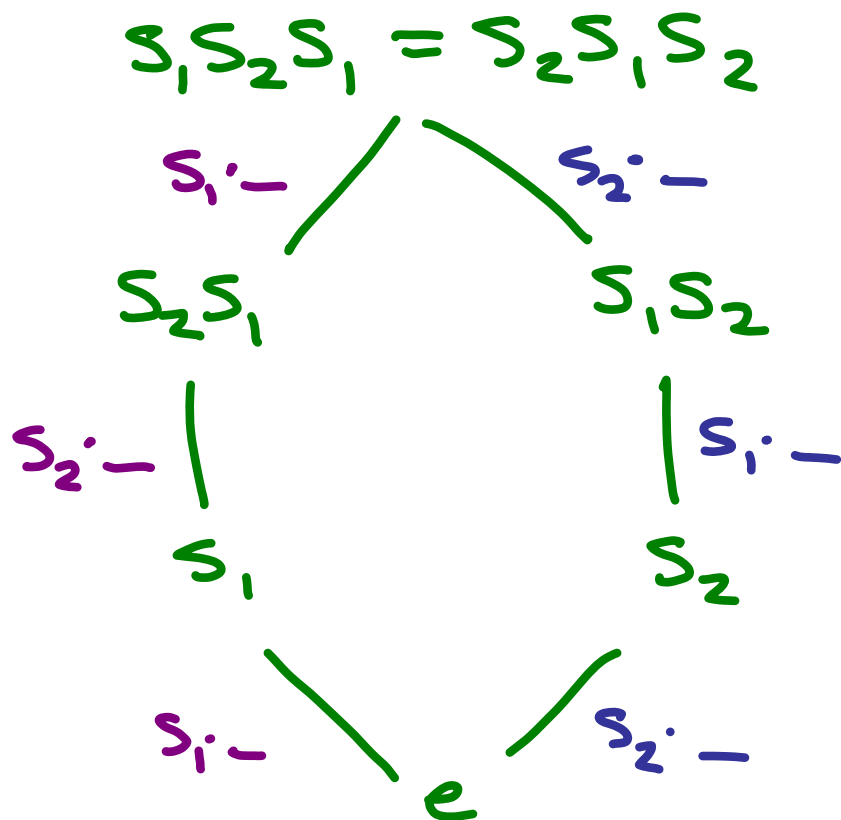
$S = \{s_1, s_2, \dots, s_{n-1}\}$ for $s_i = (i, i+1)$

with relations:

$$s_i^2 = e \quad \& \quad s_i s_{i+1} s_i = s_{i+1} s_i s_{i+1} \quad \& \quad s_i s_j = s_j s_i \\ \text{(for } |j-i| > 1 \text{)}$$

"braid moves"

e.g. Left weak order for S_3



Key Fact: Saturated chains from e to w naturally labeled w/ the "reduced expressions" $s_{i_1} \dots s_{i_{\ell(w)}}$ for w .
Likewise, saturated chains from u to $v \iff$ reduced expressions for vu^{-1} !

Properties of Reduced Expressions & Weak Bruhat Order

(see Björner-Brenti book for details)

Thm 3.2.1: Weak order on W is a meet-semilattice.

Corollary: Each $[u, v]$ is a lattice.

Lemma 3.2.3: For $J \subseteq S$, the join of atoms $\bigvee_{j \in J} a_j$ exists $\Leftrightarrow W_J = \langle j \mid j \in J \rangle$

finite, in which case $\bigvee_{j \in J} a_j = \omega_0(W_J)$

Useful Related

Fact: $a_j \leq u \Leftrightarrow \exists$ red. exp. for u s.t. s_j rightmost

longest element

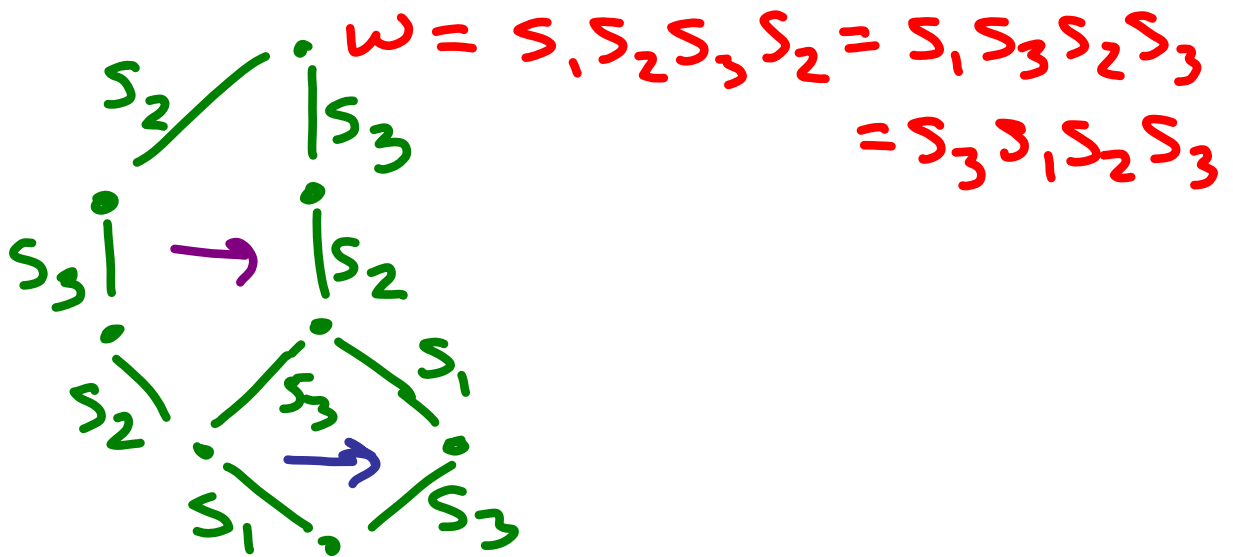
parabolic subgroup gen'd by J

Connectedness under Braid Moves

Thm 3.3.1 (Björner-Brenti) Let (W, S) be a Coxeter group $w/w \in W$. Then every two reduced expressions for w are connected via braid moves.

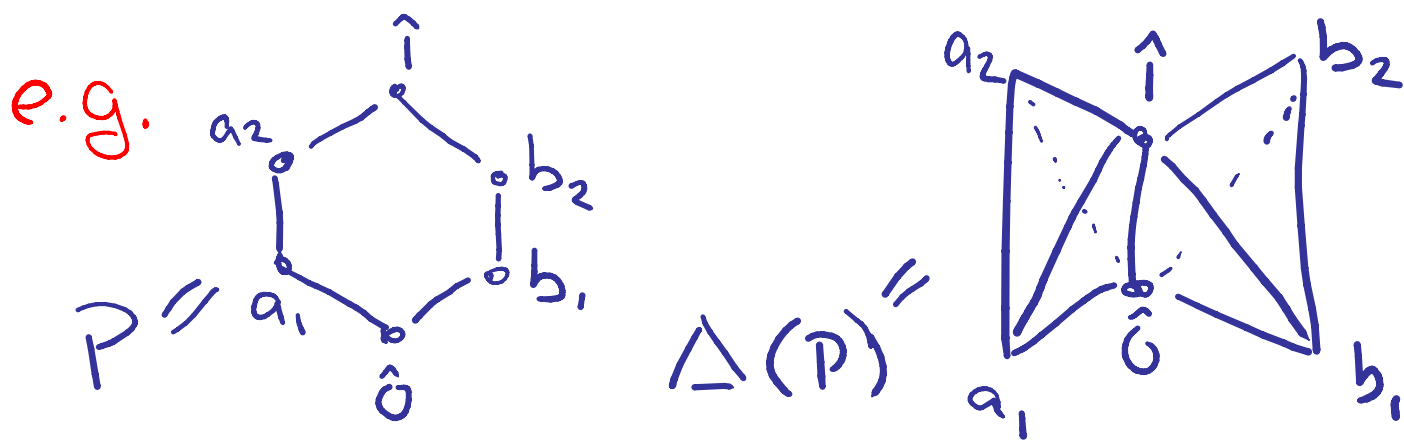
e.g. $s_1 s_2 s_3 s_2 \rightarrow s_1 s_3 s_2 s_3$
 $\rightarrow s_3 s_1 s_2 s_3$

Right weak order:



Note: Proof via lattice property for $[u, v]$

Def'n: The **order complex** (or **nerve**) of a poset P is the simplicial complex $\Delta(P)$ whose i -dimensional faces are the $(i+1)$ -chains $v_0 < \dots < v_i$ in P



Recall: $M_P(u, v) = \tilde{\chi}(\Delta(\underbrace{u, v}_{\{z \in P \mid u < z < v\}}))$

($M_P(u, v) = 0, \pm 1$ suggests ball or sphere)

Crystals + g-Crystals

A **crystal** B of type ϕ is a nonempty set B with raising \uparrow & lowering operators $e_i, f_i \uparrow f_i$

lowering operators $e_i, f_i \uparrow f_i$

$$\varepsilon_i, \varphi_i : B \rightarrow \mathbb{Z} \cup \{-\infty\}$$

$\text{wt} : B \rightarrow \Lambda = \text{weight lattice of type } \phi$

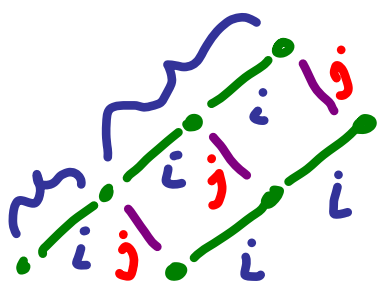
s.t.

(A1) $x, y \in B$, then $e_i(x) = y \Leftrightarrow x = f_i(y)$

both implying $\text{wt}(y) = \text{wt}(x) + \alpha_i$

$$\varepsilon_i(y) = \varepsilon_i(x) - 1$$

$$\varphi_i(y) = \varphi_i(x) + 1$$



(A2) $\varphi_i(x) - \varepsilon_i(x) = \langle \text{wt}(x), \alpha_i^\vee \rangle$

Stanbridge Crystals: "g-crystals"

(Crystals of highest weight
reps in Simply laced case)

• $\chi_{B(\lambda)}(t) = \sum_{b \in B(\lambda)} t^{\text{wt}(b)}$ = character
of irrep $B(\lambda)$

• if $y = e_i(x) \neq 0$ & $z = e_j(x) \neq 0$
for $i \neq j$ then either:

• $e_i e_j(x) = e_j e_i(x) \neq 0$

OR

• $e_i e_j^2 e_i(x) = e_j e_i^2 e_j(x) \neq 0$

• likewise for f_i operators

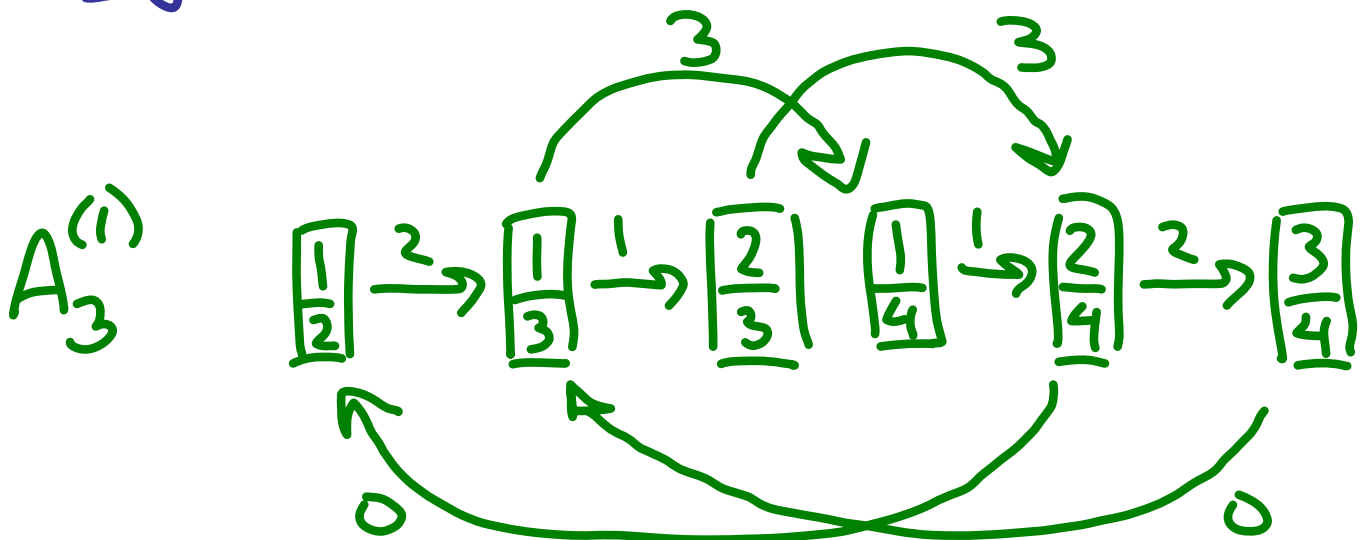
• axioms yield this & characterize
crystals of highest weight reps
in simply laced case

Additional Important Facts

- Stembridge crystals are posets with cover reln's
 $u < \cdot v \iff v = f_i(u)$

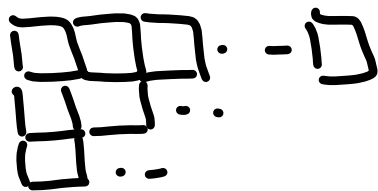
- Stemberg gave analogous "relations" for doubly laced crystals (not characterization)

A Crystal that is not a Poset



Type A SSYT Model for Crystals

A **semistandard Young tableau (SSYT)** of shape λ for $\lambda \vdash n$ is a filling of a "Ferrers diagram"



having λ_i boxes in row i for $\lambda_1 \geq \lambda_2 \geq \dots$

with positive integers $\{a_{i,j} \mid 1 \leq i \leq \ell(\lambda), 1 \leq j \leq \lambda_i\}$

$$\text{s.t. } a_{11} \leq a_{12} \leq a_{13} \leq \dots \leq a_{1,\lambda_1}$$

$$\hat{a}_{21} \leq \hat{a}_{22} \leq \dots \leq \hat{a}_{2,\lambda_2}$$

\vdots

e.g. $\lambda =$ $= (4, 2, 1, 1)$

w/ SSYT

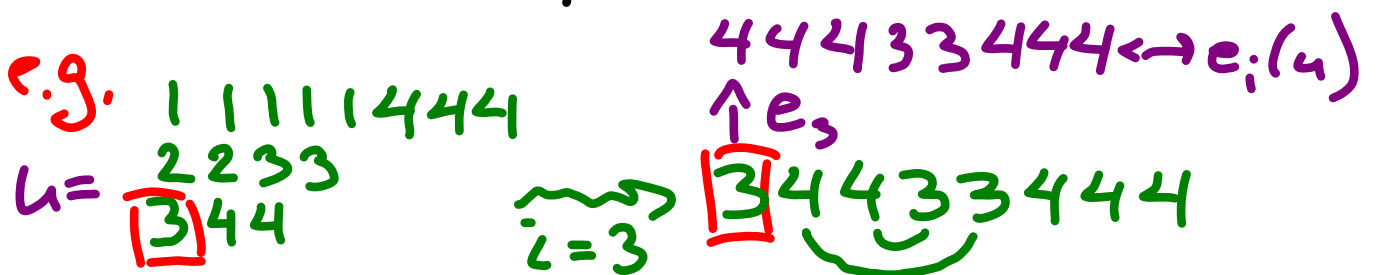
$$\begin{matrix} 1 & 1 & 2 & 5 \\ 3 & 4 & & \\ 4 & & & \\ 9 & & & \end{matrix}$$

Type A highest weight rep'n of type λ

1. $\hat{G} = \begin{matrix} 1 & 1 & 1 & - & 1 \\ 2 & 2 & - & 2 \\ 3 & 3 & - & \\ \vdots & & & \end{matrix}$ of shape λ

2. $u \xrightarrow{i} v$ for v obtained from u by increment rightmost i not in "parenthesization pair" w/ $i+1$ to an $i+1$

Parenthesization Pairs: Read leftmost column bottom to top, then subsequent columns L to R, ignoring all but $i, i+1$; omit consec $i+1, i$ and repeat.



II. Right key "k" of a KM-crystal

$$k \begin{pmatrix} 124 \\ 23 \end{pmatrix} = s_3 s_2 s_1 s_2$$



$$k \begin{pmatrix} 114 \\ 23 \end{pmatrix} = s_3 s_2 s_1$$

$$k \begin{pmatrix} 123 \\ 23 \end{pmatrix} = s_2 s_1 s_2$$



$$\begin{pmatrix} 114 \\ 22 \end{pmatrix}$$

$$k \begin{pmatrix} 113 \\ 23 \end{pmatrix} = s_2 s_1$$

$$k \begin{pmatrix} 122 \\ 23 \end{pmatrix} = s_1 s_2$$



$$R = s_3 s_2 s_1$$

$$k \begin{pmatrix} 113 \\ 22 \end{pmatrix} = s_2 s_1$$

$$k \begin{pmatrix} 112 \\ 23 \end{pmatrix} = s_1 s_2$$

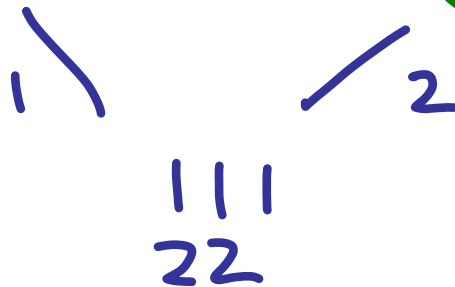


$$k \begin{pmatrix} 112 \\ 22 \end{pmatrix} = s_1$$

$$k \begin{pmatrix} 111 \\ 23 \end{pmatrix} = s_2$$

k: crystal poset \rightarrow weak Bruhat order

$$u \leq v \Rightarrow k(u) \leq k(v)$$



$$\text{Key } k(\hat{0}) = e$$

New Algorithm to Calculate Right Key of a KM-Crystal

- (1) $\text{key}(\hat{\sigma}) = e$
- (2) if $\hat{\sigma} \xrightarrow{i} a$, then $\text{key}(a) = s_i$
(i.e. $\hat{\sigma} \leftarrow a$)
- (3) if v covers 2 or more elements
then $\text{key}(v) = \bigvee_{\{u \mid u \rightarrow v\}} \text{key}(u)$
(for join taken in weak order)
- (4) if $u \xrightarrow{i} v$ and v does not cover
any other elements, then:
 - (a) $\text{key}(v) = \text{key}(u)$ if $\exists u' \xrightarrow{i} u$
 - (b) $\text{key}(v) = s_i \cdot \text{key}(u)$ otherwise

Crucial Properties of Key

Thm (Littlemann): Given any symmetrizable Kac-Moody algebra A , the key of any crystal of type A satisfies:

$$K(f_p(F)) = \begin{cases} K(F) & \text{if } e_p(F) \neq 0 \\ s_p K(F) \text{ or } K(F) & \text{if } e_p(F) = 0 \end{cases}$$

Also, if $e_p(F) = 0$ then $s_p K(F) > K(F)$

Corollary: If $K(F) = s_{i_1} \dots s_{i_r}$ then there exists saturated chain from F to $\hat{0}$ given by applying $e_{i_r}^{d_r} \dots e_{i_1}^{d_1}$ to F for some $d_1, \dots, d_r > 0$.

Key Polynomials \neq right / left key

(see Lascoux-Schutzenberger \neq e.g.
Reiner-Shimozono)

- Motivations:
- (1) Schubert poly. G_w is positive sum of "key polynomials"
 - (2) Key polynomial records character for Demazure module
 - (3) The (closely related) right \neq left key maps determine smallest Demazure modules containing a given crystal element
 - (4) These will give us poset map from g -crystal to weak Bruhat order, transferring properties

Relation to Reiner-Shimozono Viewpoint on Key Polynomials

- $\partial_i = \frac{1-s_i}{x_i - x_{i+1}} \quad \neq \quad \pi_i = \partial_i x_i$
- $K_\alpha = \pi_{i_1} \dots \pi_{i_r} x^{\lambda(\alpha)}$ for α composition of $n \neq s_{i_1} \dots s_{i_r}$ sorting α to $\lambda(\alpha)$

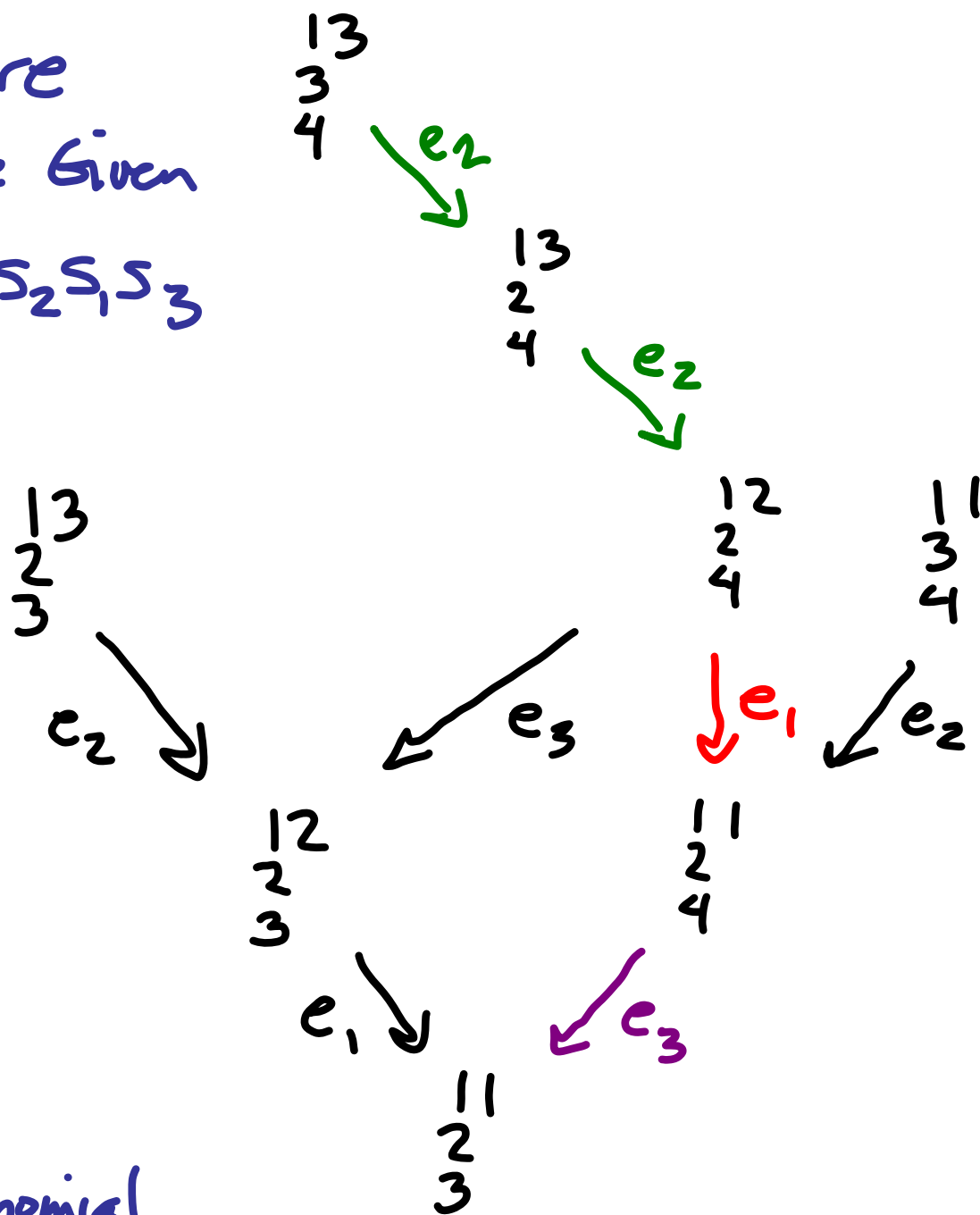
e.g. $K_{(1,0,2,1)} = \pi_2 \pi_1 \pi_3 x^{(2,1,1,0)}$

$= \pi_2 \pi_1 (x_1^2 x_2 (x_3 + x_4))$

$= \pi_2 (x_1 x_2 x_3 (x_1 + x_2) + x_1 x_2 x_4 (x_1 + x_2))$

$= x_1^2 x_2 x_3 + x_1 x_2 x_3 (x_2 + x_3) + x_1^2 x_4 (x_2 + x_3) + x_1 x_4 (x_2^2 + x_2 x_3 + x_3^2)$

Demazure
Module Given
by $w = s_2 s_1 s_3$



Key Polynomial

$$K_{(1,0,2,1)} = \sum_{T' \leq T} x^{T'}$$

componentwise

$$K(T') \leq_{\text{Bruhat}} K(T) \iff \text{no higher } e_i \text{ exponents}$$

III. Positive Results for Lower Intervals $[\hat{0}, u]$

Recall: $M_p(u, u) = 1$
 $M_p(u, v) = -\sum_{u \leq z < v} M_p(u, z)$

Thm 1 (H.-Leuvert): Given u in a symmetric KM-crystal, then $M(\hat{0}, u) = 0, \pm 1$. More specifically, $M(\hat{0}, u) = 0$ unless $\text{key}(u) = \omega_0(J)$ for some parabolic subgroup ω_J with u the unique smallest element in $\text{key}^{-1}(\omega_0(J))$, in which case $M(\hat{0}, u) = (-1)^{|J|}$.

Proof Ingredients:

Thm 2 (H.-Lenart): Given a symmet.

KM-crystal \dagger given any parabolic

W_J , then $\text{key}^{-1}(\omega_0(W_J))$

has a unique minimal element

and a unique maximal element.

(Proof via alcove path model)

Prop'n: Each $w \in W$ has unique maximal element $u \in W$ in weak Bruhat order s.t.

- $u \leq_{\text{weak}} w$

- there exists parabolic subgroup W_J s.t. $u = \omega_0(J) = \text{longest el't}$

Thm 3 (H-Lenart): Each lower interval $(\hat{0}, u)$ in a symmet. KM-crystal has $\Delta(\hat{0}, u) \simeq$ ball or sphere, getting $S^{|\mathcal{J}|-2}$ for $u = \min(k^{-1}(\omega_0(\mathcal{J})))$. Likewise for upper intervals in finite KM-crystals.

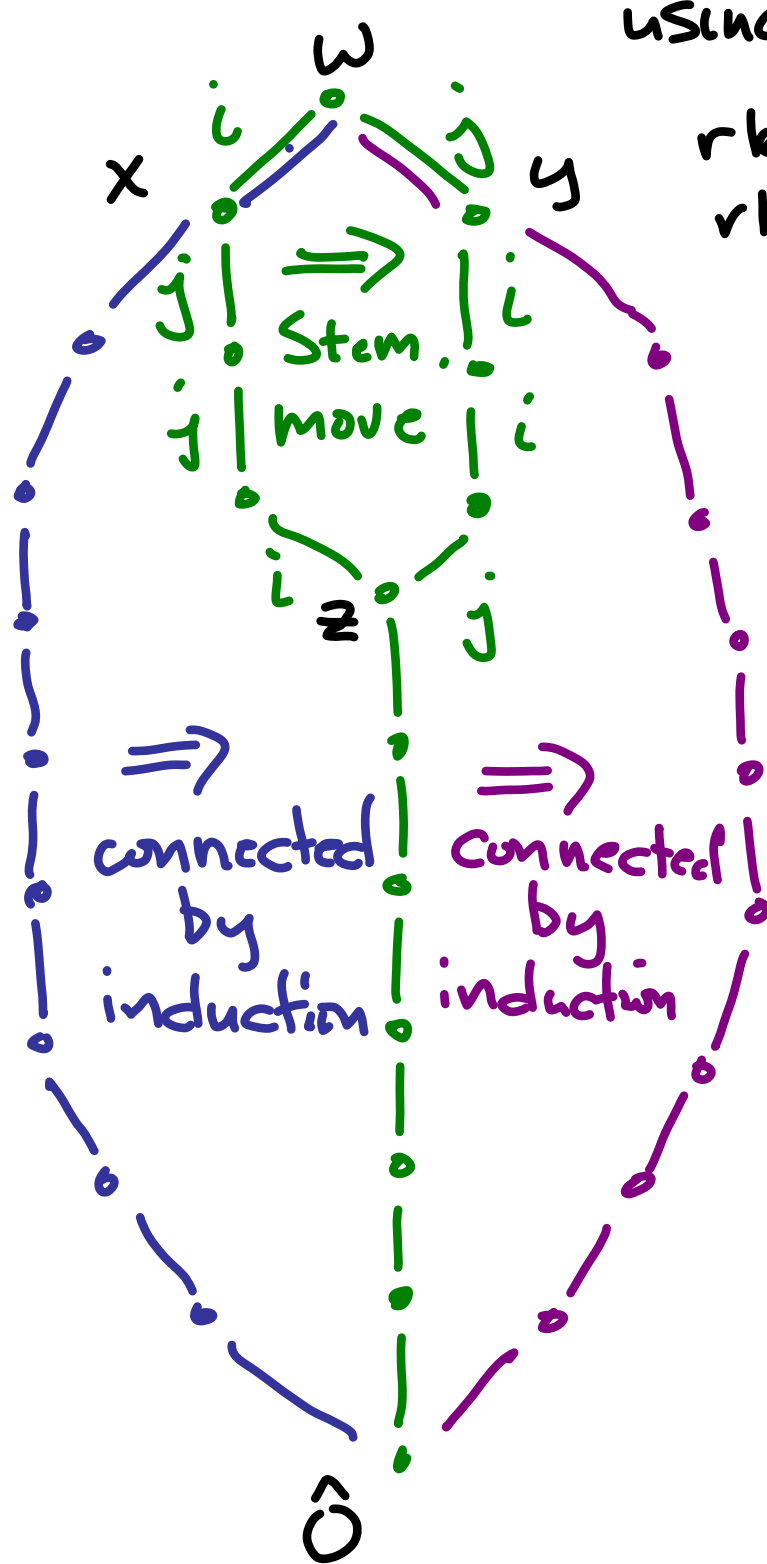
Proof Method: Quillen fiber lemma based upon:

$f: \text{Crystal Poset} \longrightarrow \text{Boolean Algebra } \{\mathcal{J} \mid \mathcal{J} \subseteq \mathcal{I}\}$

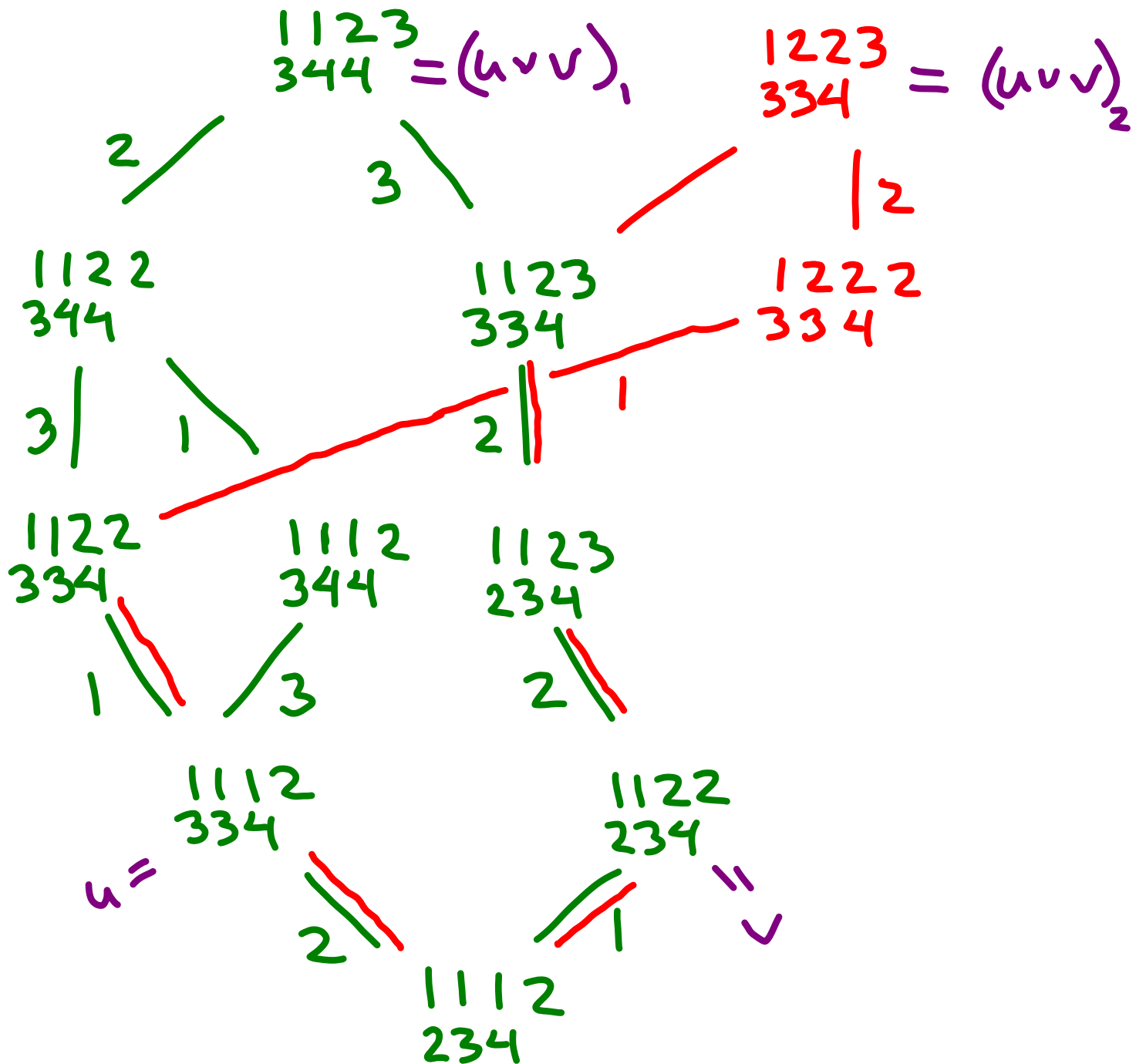
$x \longmapsto \max \{\mathcal{J} \mid \omega_0(\mathcal{J}) \leq k(x)\}$

Quillen Fiber Lemma: Poset map $f: P \rightarrow Q$ s.t. each $\Delta(f^{-1}(q))$ is contractible implies $\Delta(P) \simeq \Delta(Q)$.

Proof Idea: Induction on rank using $z \geq \hat{0}$ \dagger
 $rk(x) < rk(w)$
 $rk(y) < rk(w)$



Non-Lattice Example



(Why proof fails for arbitrary intervals)

IV. Negative Results for Arbitrary (not necessarily lower)

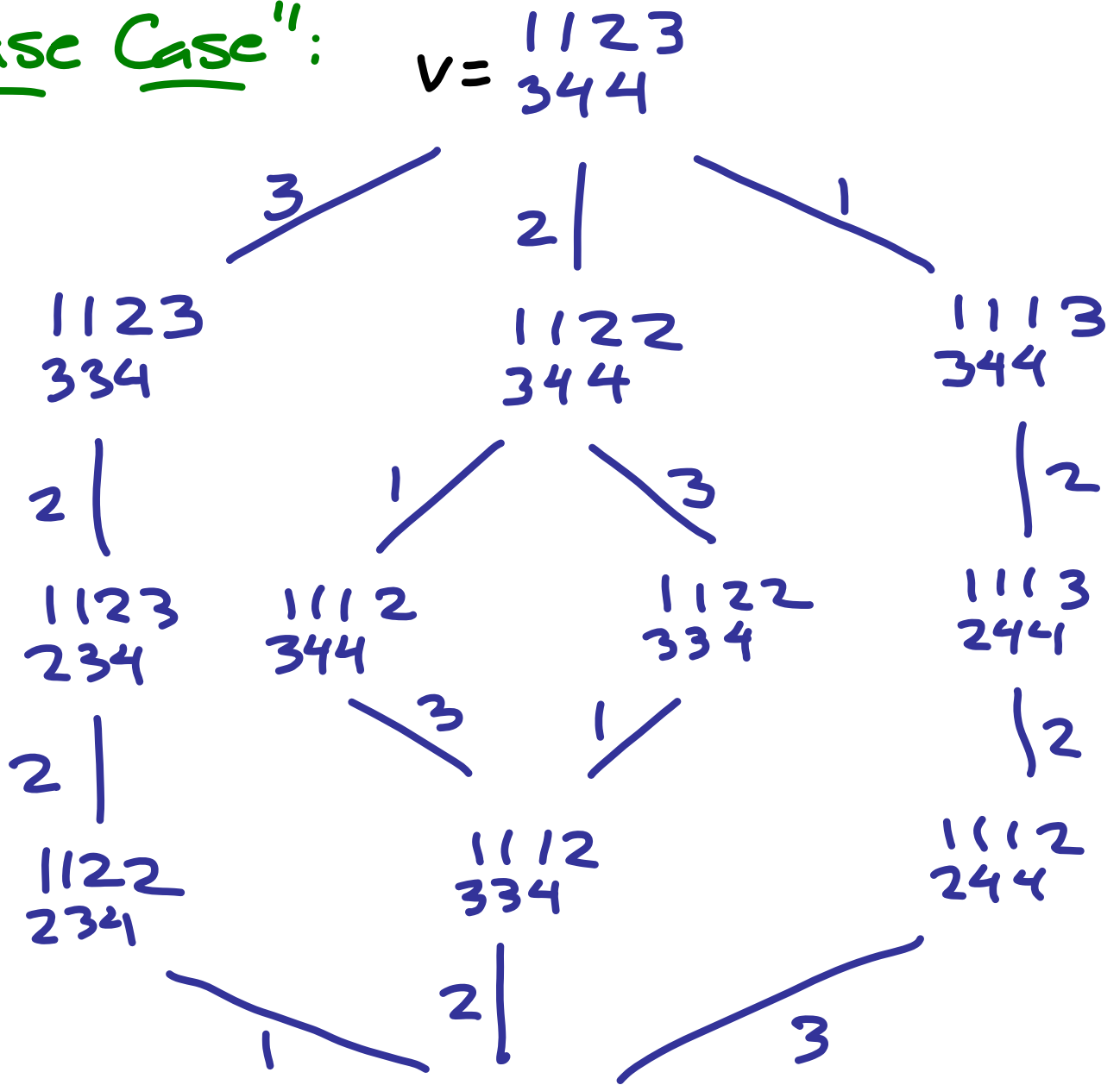
Crystal Poset Intervals (type A)

Thm 5 (H. - Lenart): There exist elements u, v in type A g -crystals with $M(u, v) = 2j$ for every positive integer j .

Thm 6 (H. Lenart): There exist type A intervals $[u, v]$ with $\text{rk}(v) - \text{rk}(u)$ arbitrarily large s.t. (u, v) is disconnected

Infinite Family of Examples

"Base Case":



$M_p(u, v) = 2$ $u = \begin{matrix} 1112 \\ 234 \end{matrix}$

‡ not connected by "Steinbridge moves"

Examples with $M(u,v) = 2^j$

$j=1: u = \begin{matrix} 1112 \\ 234 \end{matrix}$

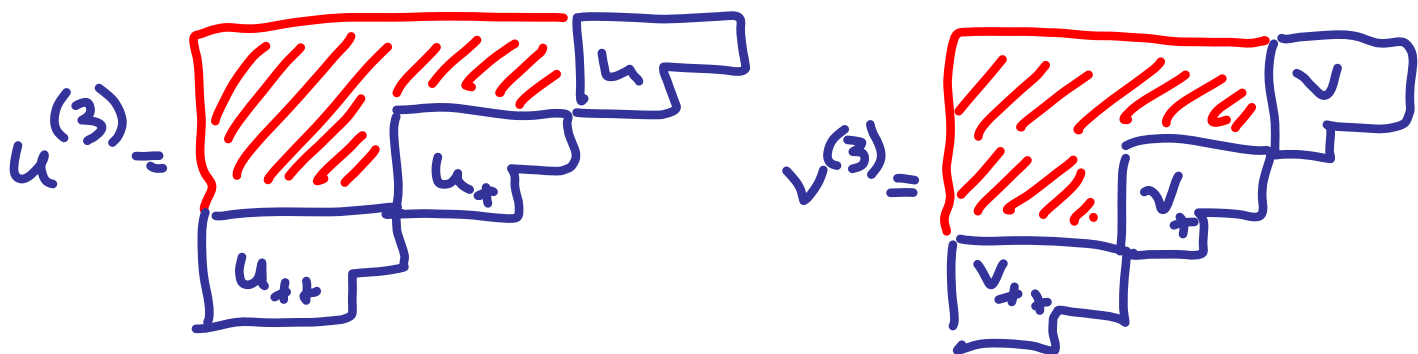
$v = \begin{matrix} 1123 \\ 344 \end{matrix}$

$j=2: u^{(2)} = \begin{matrix} 1111 & \boxed{1112} \\ 2222 & \boxed{234} \\ \boxed{6667} & \\ \boxed{789} & \end{matrix}$ $v^{(2)} = \begin{matrix} 1111 & \boxed{1123} \\ 2222 & \boxed{344} \\ \boxed{6678} & \\ \boxed{899} & \end{matrix}$

$u_+ := u+S = \begin{matrix} \boxed{6667} \\ \boxed{789} \end{matrix}$ $v_+ := v+S = \begin{matrix} \boxed{6678} \\ \boxed{899} \end{matrix}$

$[u^{(2)}, v^{(2)}] \cong [u, v] \times [u, v]$

so $M(u^{(2)}, v^{(2)}) = 2^2$



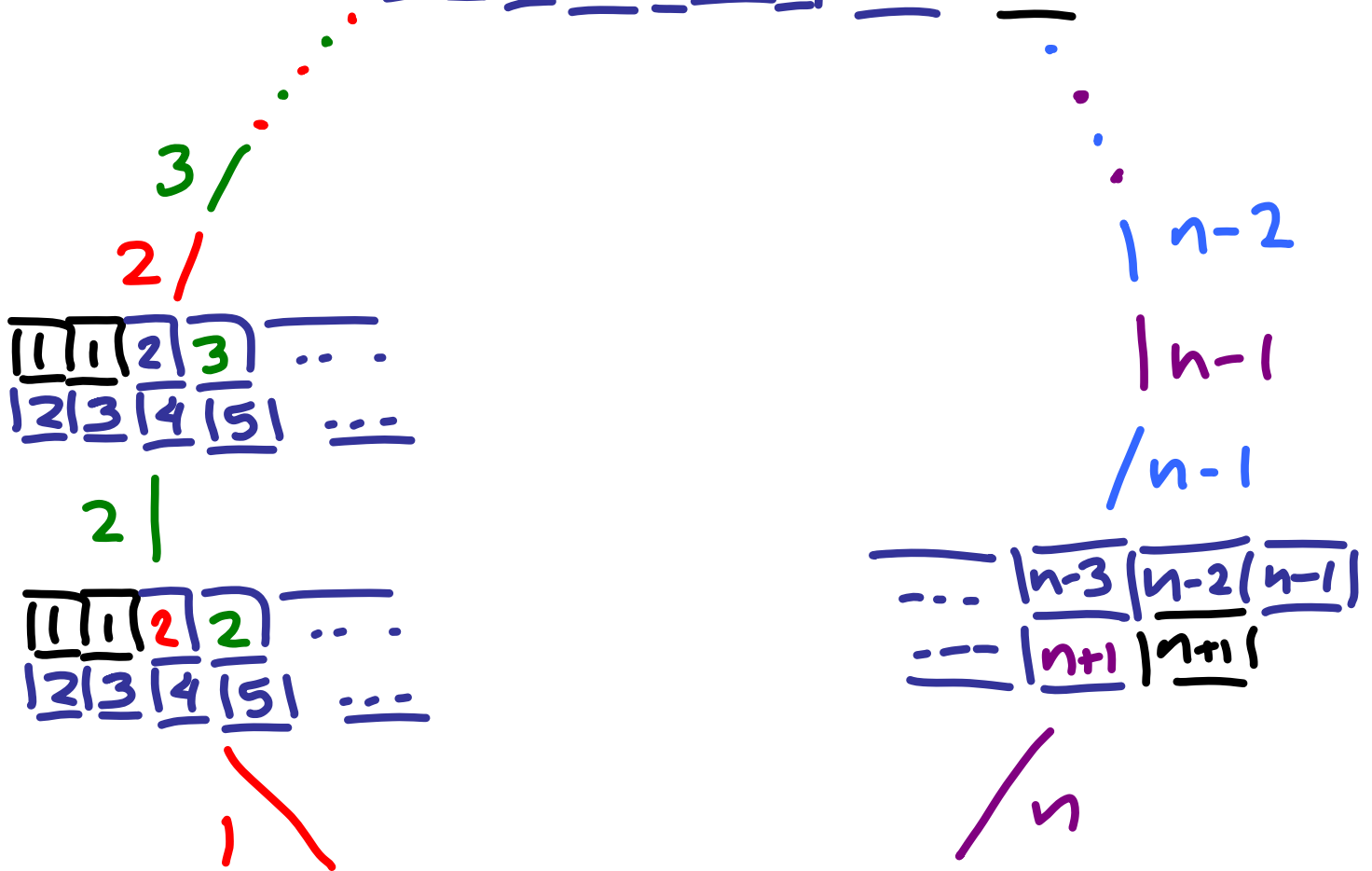
$[u^{(k)}, v^{(k)}] \cong \underbrace{[u, v] \times \dots \times [u, v]}_{k\text{-fold}} \quad M = 2^k$

Arbitrarily High Rank

Disconnected Open Intervals

$$v = \begin{array}{|c|c|c|c|} \hline 1 & 1 & 2 & 3 \\ \hline \end{array} \dots \begin{array}{|c|c|c|} \hline n-2 & n-1 & n \\ \hline \end{array}$$

$$\begin{array}{|c|c|c|c|} \hline 3 & 4 & 5 & 6 \\ \hline \end{array} \dots \begin{array}{|c|c|} \hline n+1 & n+1 \\ \hline \end{array}$$



$$u = \begin{array}{|c|c|c|c|} \hline 1 & 1 & 1 & 2 \\ \hline \end{array} \dots \begin{array}{|c|c|c|} \hline n-3 & n-2 & n-1 \\ \hline \end{array}$$

$$\begin{array}{|c|c|c|c|} \hline 2 & 3 & 4 & 5 \\ \hline \end{array} \dots \begin{array}{|c|c|} \hline n & n+1 \\ \hline \end{array}$$

label sequences: $1, 2, 2, 3, 3, 4, 4, \dots, n-1, n-1, n$
 $\neq n, n, n, \dots, 2, 2, 1$ in distinct components

Consequence: Arbitrarily high degree relations $e_{i_1} \dots e_{i_d}(u) = e_{j_1} \dots e_{j_d}(u)$ amongst crystal operators applied to u not implied by any lower degree relations.

Some Further Questions:

1. Positive results in any additional generality (more general intervals)?
2. Interpretation/applications for Möbius function of a crystal?
3. Where/how exactly can the lattice property fail?