Lazy vs. non lazy

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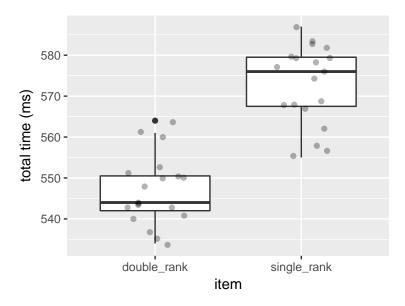
1 Double vs. single rank

Table 1: Time (in ms) of 500K calls to w1() based on single_rank() or double_rank() methods on 100MB random DNA input; Mean/sd over 20 repetitions.

item	avg_time	sd_time
double_rank	546.89	8.49
$single_rank$	572.74	9.65

Table 2: Absolute (double_rank / single_rank) and relative (100 * |double_rank - single_rank| / single_rank) ratios of average times.

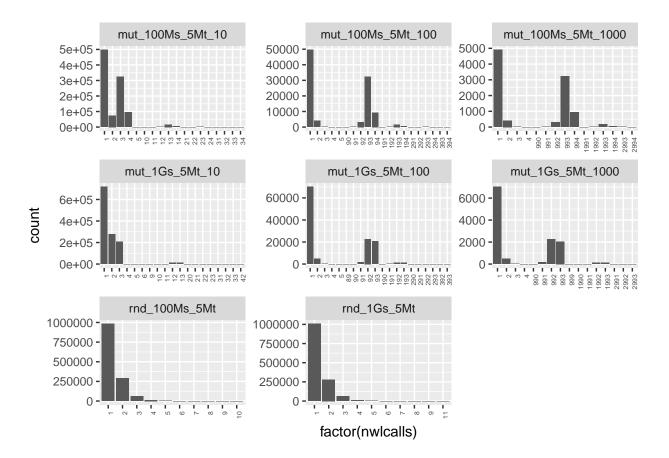
double_rank	single_rank	abs_ratio	rel_ratio
546.89	572.74	0.95	4.51



2 Lazy vs non-lazy

2.1 Input properties

For various types ("mut_XMs_YMt_Z" means s and t are random identical strings of length X, and Y million respectively with mutations inserted every Z characters. "rnd_XMs_YMt" means s and t are random strings of length X, and Y million respectively) of inputs run the MS algorithm and count the number of consecutive lazy_wl() calls.



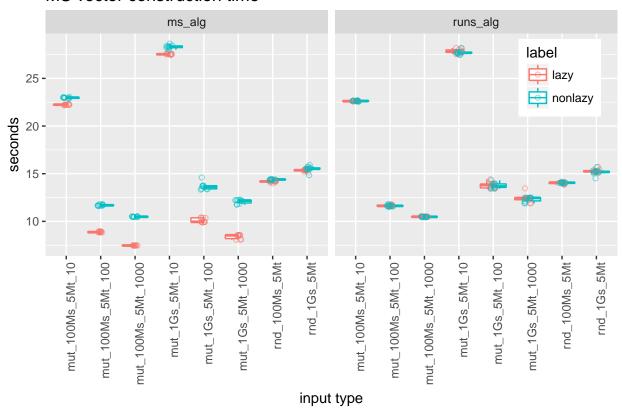
2.2 Code

The lazy and non-lazy versions differ in a couple of lines of code as follows

```
if(flags.lazy){
    for(; I.first <= I.second && h_star < ms_size; ){</pre>
        c = t[h_star];
        I = bstep_interval(st, I, c); //I.bstep(c);
        if(I.first <= I.second){</pre>
            v = st.lazy_wl(v, c);
            h_star++;
        }
    }
    if(h_star > h_star_prev) // // we must have called lazy_wl(). complete the node
        st.lazy_wl_followup(v);
} else { // non-lazy weiner links
    for(; I.first <= I.second && h_star < ms_size; ){</pre>
        c = t[h_star];
        I = bstep_interval(st, I, c); //I.bstep(c);
        if(I.first <= I.second){</pre>
            v = st.lazy_wl(v, c);
            h_star++;
        }
    }
}
```

2.3 Run time

MS vector construction time



The right panel shows the time to construct the **runs** vector. This stage is the same for both versions and is shown as a control. On the left panel it can be seen that speedup correlates positively with both the size of the indexed string and the mutation period.

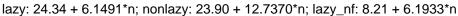
2.4 Sandbox timing

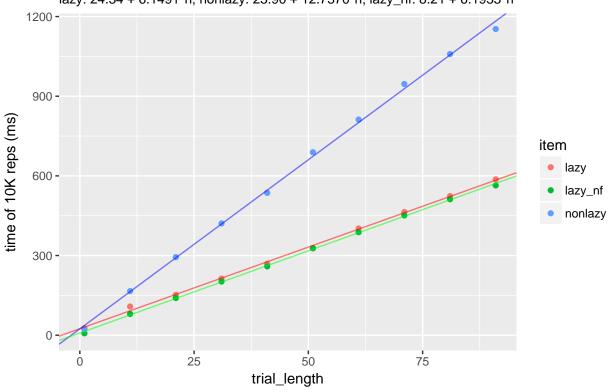
Measure the time of 10k repetitions of

- (lazy) n consecutive lazy_wl() calls followed by a lazy_wl_followup()
- (nonlazy) n consecutive wl() calls
- (lazy_nf) n consecutive lazy_wl() calls

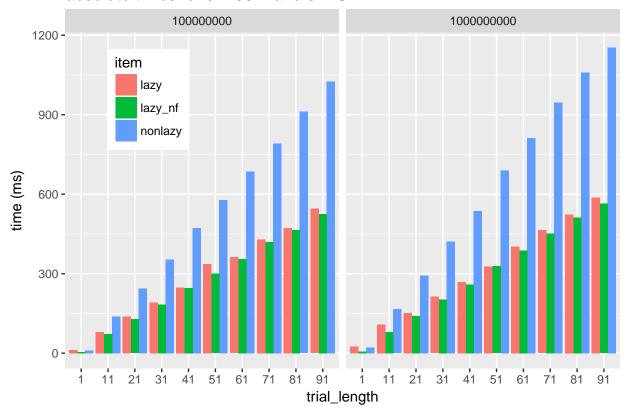
```
// lazy
for(size_type i = 0; i < trial_length; i++)
    v = st.lazy_wl(v, s_rev[k--]);
if(h_star > h_star_prev) // // we must have called lazy_wl(). complete the node
    st.lazy_wl_followup(v);
...
// non-lazy
for(size_type i = 0; i < trial_length; i++)
    v = st.wl(v, s_rev[k--]);
...
// lazy_nf
for(size_type i = 0; i < trial_length; i++)
    v = st.lazy_wl(v, s_rev[k--]);</pre>
```

indexed input size 1G





absolute times for s=100M and s=1G



2.5 Check

In the experiments above we ran the program with the "lazy" or "non-lazy" flag and measured. The total time of each experiment can be written as $t_l = l_l + a$ and $t_n = l_n + a$ for the two versions respectively; only the ts being known. Furthermore, we have \hat{l}_l and \hat{l}_n estimations. Hence we should expect

$$\delta t = t_l - t_n = l_l + a - l_n - a = l_l - l_n \approx \delta \hat{l} = \hat{l}_l - \hat{l}_n$$

1	. 1 1 4
b_path t_l t_n l_l l_n delta_t delta	a_l_hat
mut_100Ms_5Mt_10 22.23 22.97 4.05 3.53 -0.74	0.52
mut_100Ms_5Mt_100 8.89 11.70 2.90 4.07 -2.81	-1.16
mut_100Ms_5Mt_1000 7.47 10.49 2.79 4.12 -3.02	-1.33
mut_1Gs_5Mt_10 27.54 28.32 4.08 4.08 -0.78	-0.01
mut_1Gs_5Mt_100 10.11 13.66 3.15 4.65 -3.55	-1.50
mut_1Gs_5Mt_1000 8.40 12.09 3.05 4.70 -3.68	-1.65
rnd_100Ms_5Mt 14.17 14.38 4.68 3.65 -0.20	1.03
rnd_1Gs_5Mt 15.38 15.51 4.15 3.94 -0.14	0.20

The numbers are not identical (process dependent factors might influence the running time of function calls), but they are correlated $(corr(\delta t, \delta \hat{l}) = 0.96)$.