

NAME

fi ndBgpAtoms – compute BGP policy atoms

SYNOPSIS

fi ndBgpAtoms [-d *debug-level*] [-k] [-T *dir*] fi le-prefi x

PARAMETERS

file-prefix

The *prefi x* of fi lenames generated by straightenRV.

OPTIONS

-d *debug-level*

Verbosity of debugging output, 0 for no debugging output, 1 (default) for normal debugging output, 2 for all debugging output.

-k

Whether to keep intermediate fi les or not after fi ndBgpAtoms has run. All fi les generated by fi ndBgpAtoms except fi le-prefi x.atoms.ccdf, fi le-prefi x.atoms.asp.gz, fi le-prefi x.atoms.p2a and fi le-prefi x.atoms.full.gz are intermediate fi les.

-T *dir*

fi ndBgpAtoms uses UNIX sort. This option provides sort with a temporary directory. Handy when sort runs out of space.

DESCRIPTION

fi ndBgpAtoms computes BGP policy atoms. A BGP policy atom, or 'atom' for short, is defi ned relative to a system of BGP routers. This collection of scripts uses RouteViews peers as the system of BGP routers. Each atom consists of *prefi xes* that are treated equivalently by the chosen system of BGP routers.

A precise defi nition of an atom is as follows. Two *prefi xes* are said to be *path equivalent* if no BGP router can be found among the considered BGP routers that sees them with different AS paths. An equivalence class of this relation is called a *BGP policy atom*. It follows from this defi nition that *prefi xes* in the same atom share a collection of AS paths.

Atoms are computed from a number of fi les generated by straightenRV, which must be located in the working directory:

fi le-prefi x.atoms.p2a.gz

The information needed to compute atoms: *prefi xes*, peers and AS paths.

fi le-prefi x.peer

Used for *prefi x* selection. In this version, atoms are computed from 'global *prefi xes*' only. (Non-global *prefi xes* are ignored.) A global *prefi x* is a *prefi x* that is carried by all the peers listed in the fi le-prefi x.peer fi le. Use straightenRV's -c option to create a selection of peers (and therefore global *prefi xes*) based on peer size.

Many intermediate and auxiliary fi les are produced. The main output fi les for atoms are:

fi le-prefi x.atoms.asp.gz

Maps each atom name to the collection of AS paths for the atom.

fi le-prefi x.atoms.p2a

Maps each *prefi x* to the atom that the *prefi x* belongs to.

The following summarises the input and output fi les of straightenRV and fi ndBgpAtoms: To summarise the

above:

```
straightenRV -> .peer + .pfastp.gz + ...
.peer + .pfastp.gz -> fi ndBgpAtoms -> .atoms.asp.gz + .atoms.p2a + ...
```

ATOM COMPUTATION

This section describes the computation of atoms by `fi ndBgpAtoms` in several steps. Each step takes one or more `fi les` produced by the previous step and generates `fi les` to be used in the next step.

Sorting `.pfastp.gz` to `.pfastpsorted`

The `.pfastp.gz` `fi le` is sorted and written to a `.pfastpsorted` `fi le`. It sorts on `prefi x` `fi rst`, then on `peer IP address`, and `fi nally` on `AS path`.

Joining AS paths in `.pfastpsorted` to `.pfastpcoll`

For each `prefi x`, groups all AS paths for that `prefi x` in `.pfastpsorted` together, forming a collection of AS paths. Writes the `prefi x` and the AS path collection to `.pfastpcoll`. Each AS path is `fi rst` prepended with the IP address of the peer through which the AS path was learnt.

Only global `prefi xes` are considered; non-global `prefi xes` are dropped in this step, and are not written to `.pfastpcoll`. Nor are they considered in the remaining steps. See the '`fi le-prefi x.peer`' description above for a definition of global `prefi xes`.

Sorting `.pfastpcoll` by AS path collection to `.pfastpcollsorted`

Sorts the `.pfastpcoll` `fi le` by the collections of AS paths. Note that this effectively places the `prefi xes` that are in the same atom together.

Grouping `prefi xes` in `.pfastpcollsorted` by atoms to `.atoms`

From the `.pfastpcollsorted` `fi le`, reads `prefi xes` and their collections of AS paths, groups them into atoms, sorts the atoms by the number of `prefi xes` per atom (in reverse order), and writes the result to the `.atoms` `fi le`. Also writes a distribution of `prefi x` counts per atom to the `.atoms.ccdf` `fi le`.

The format of the `.atoms` `fi le` consists of four lines per atom:

1. The number of `prefi xes` in the atom.
2. The sorted list of `prefi xes` in the atom.
3. The collection of AS paths for the atom (as copied from `.pfastpcollsorted`).
4. An empty separator line.

Naming atoms in `.atoms` to `.atoms.asp.gz`, `.atoms.p2a` and `.atoms.full.gz`

Uniquely names the atoms found in `.atoms`. The atom name is a string composed as follows:

```
as<origin>np<#prefi xes>at<sequence>
```

where `<origin>` is the origin AS of the atom (or multiple origin ASes separated by '`_`'), `<#prefi xes>` is the number of `prefi xes` in the atom, and `<sequence>` is a sequence number (starting with 1) ranging over all atom names that have the same `<origin>` and `<#prefi xes>` parts. Examples:

```
as701np1133at1
as5676np125at1
as5676np125at2
as6140_16528np34at1
```

The following files are produced:

.atoms.asp.gz

Maps each atom name to the collection of AS paths for the atom. Note that in an earlier step, each AS path was prefixed with the IP address of the peer from which the AS path had been learnt.

.atoms.p2a

Maps each prefix to the atom that the prefix belongs to.

.atoms.full.gz

Contains the data of .atoms, adding an atom counter (counter ranging over all atoms found) and the atom name.

FILES

All output files generated by `findBgpAtoms` start with the *file-prefix* that was passed on the command line. The output files are described above.

SEE ALSO

`straightenRV(1)`

BUGS and TODO

Atoms definition

This script computes BGP policy atoms according to the definition of atoms given earlier. We should also compute atoms using a modified definition more appropriate for use within the 'atoms' project:

In determining path equivalence of two prefixes, we ignore the part of a prefix's AS path that does not include any of the RV peers' ASes. From this definition, it follows that prefixes in the same atom share a set of truncated AS paths, where each AS path is truncated to exclude the part that falls outside the set of peer ASes. Note that the number of atoms under this definition is smaller than the number of atoms under the original definition.

Origin-declared atoms

A third kind of atom exists, the 'origin-declared atom'. An origin-declared atom is defined by the AS that originates the prefixes in the atom. The prefixes in an origin-declared atom share the same origin link (the rightmost link in the AS path), i.e. are all announced by the same origin AS to the same neighbour AS. Origin-declared atoms are 'declared', in that they are 'imposed' by the origin AS; any policy applied by ASes other than the origin AS is ignored. Per prefix origin links are listed in the .pref file produced by `straightenRV` in the 'oriLkstats' column, but are not further grouped into atoms.

Statistics

Andre Broido's original `findBgpAtoms` script produced many interesting statistics on atoms that will one day be incorporated.

AUTHORS

Patrick Verkaik (patrick@caida.org)

Andre Broido: algorithms and scripts before rewrite

Young Hyun: code review

REFERENCES

Atoms web page:

<http://www.caida.org/projects/routing/atoms/>

Andre Broido, kc claffy, 'Analysis of RouteViews BGP data: policy atoms', Proceedings of the Network-

Related Data Management workshop, Santa Barbara, May 23, 2001.
<http://www.caida.org/outreach/papers/2001/NdrmBgp/>

Andre Broido, Evi Nemeth, kc claffy, 'Internet Expansion, Refinement, and Churn', European Transactions on Telecommunications, January 2002.
<http://www.caida.org/outreach/papers/2002/EGR/>

Andre Broido, kc claffy, 'Complexity of global routing policies'.
<http://www.caida.org/outreach/papers/2001/CGR/>