pgSimload v.1.0.0-beta Documentation

Jean-Paul Argudo

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Contents

1	Ove	erview	2						
2	Run	Running, Building and Installing binary							
	2.1	Running with Go	. 2						
	2.2	Using binaries provided	. 2						
	2.3	Building binaries	. 3						
	2.4	DEB and RPM packages	. 3						
3	Usa	ages	3						
4	Exai	amples given	4						
	4.1	examples/simple/	. 5						
	4.2	examples/PG_15_Merge_command/	. 5						
	4.3	examples/testdb/	. 6						
5	Ove	Overview of flags and parameters 7							
	5.1	All modes : flags	. 7						
	5.2	SQL-loop mode: parameters							
	5.3	Patroni-monitoring mode: parameters	. 8						
	5.4	Session parameters template file creation	. 8						
6	Refe	Reference: parameters and flags							
	6.1	1 Common flags and parameters							
		6.1.1 config (JSON file) [MANDATORY]	. 9						
		6.1.2 contact (flag) [OPTIONAL]	. 10						
		6.1.3 help (flag) [OPTIONAL]	. 10						
		6.1.4 license (flag) [OPTIONAL]	. 10						
		6.1.5 version (flag) [OPTIONAL]	. 10						
	6.2	SQL-loop mode parameters	. 11						
		6.2.1 create (JSON text file) [OPTIONAL]	. 11						
		6.2.2 script (SQL text file) [MANDATORY]	. 11						
		6.2.3 session_parameters (JSON text file) [OPTIONAL]	. 11						
	6.3	Patroni-monitoring mode flag and parameters	. 13						
		6.3.1 patroni (value) [MANDATORY]	. 13						
	6.4	4 Session parameters template file creation							
		6.4.1 config (value) [MANDATORY]	. 16						
		6.4.2 create gucs template (value) [MANDATORY]	16						

1 Overview

Welcome to pgSimload!

The actual version of the program is:

pgSimload version 1.0.0-beta - July, 24th 2023

pgSimload is a tool written in Go, and accepts 2 different modes of execution:

- **SQL-loop mode** to execute a script infintely on a given schema of a given database with a given user
- **Patroni-monitoring mode** to execute a monitoring on a given Patroni cluster. So you need one of such for this mode to be useful to you

Given the mode you choose, some parameters are mandatory or not. And the contexts of executions are different. Please refer to the complete documentation in docs/pgSimload.doc.md

2 Running, Building and Installing binary

2.1 Running with Go

This is very straightforward if you have Go installed on your system. You can run the tool with Go from the main directory of the project like:

```
1 $ go run main.go <parameters...>
2 $ go run main.go -h
```

2.2 Using binaries provided

If you don't have Go installed on your system, you can also just use one of the binaries provided in bin/. If you want to build your own binary you can build it too, as described in the next paragraph.

Please read carefully bin/README.md, where we told you wich binary to use depending on your environment, specially in Linux.

Note that Mac and Windows versions aren't fully tested at the moment.

Feedback is welcome in any cases!

2.3 Building binaries

You can use the provided script build.sh:

```
1 $ sh build.sh
```

2.4 DEB and RPM packages

We've started tests to build those packages but at the moment, the work hasn't finish yet. Those packages will be available soon.

3 Usages

This tool can be used in different infrastructures:

- on the localhost, if a PostgreSQL is running on it
- on any distant stand-alone PostgreSQL or PostgreSQL cluster, in bare-metal of VMs
- on any PostgreSQL stand-alone PostgreSQL or PostgreSQL cluster running in a Kubernetes environment.

This tool can be used in different scenarios:

- insert dummy data (mostly randomly if you know about, mostly, generate_series() and random() PostgreSQL functions) any DB with the schema of your choice and the SQL script of your choice
 - if your database doesn't have a schema yet, you can create in a create.json file. Look for examples on how to do that in the examples / directory. It should straightforward. That file is **not** mandatory, as pgSimload need at least a -config <file> and a -script <file> to run.
 - the SQL script of your choice. For that purpose you create a plain SQL file, where you put everything you want in it. Beware the parsing is really simple, it would probably fail when creating complex things like functions in this script.
 - you can set special parameters to the session like SET synchronous_commit TO 'on'or SET work_mem TO '12MB' if you want the SQL script's sessions to be tweaked depending your needs. This is usefull to compare the performances or behaviour in replication or others things. For that you'll have to use the -session_parameters <

session_parameters.json> parameter for pgSimpload. Otherwise, without this, every DEFAULT values will of course apply.

- if you're too lazy to gather those session parameters, you can create a template file you can letter modify and adapt to your needs. For that pgSimload will create a template file in the name you want, based on a given connection. Look for -create_gucs_template in this documentation.
- this "dummy data insertion" is most often used to simulate some write work on a PostgreSQL server (standalone or the primary of a PostgreSQL cluster with a(some) replica(s).
- test failovers, or what happens when a DB is down: pgSimLoad handles those errors. Give it a try: simply shuting down your PostgreSQL server while it runs... You'll see it throwing errors, then restarting to load once the PostgreSQL server ("primary" if you use replication) is back.
- monitor a PostgreSQL cluster that uses Patroni, with the special --patroni <config.json
 parameter, that has to come with a --config <config.json> where the later will use
 mandatorily the postgres user, because, on that mode, we use a special trick to get the primary's name, and this trick can only be done by a superuser in PostgreSQL (so it can be something else than postgres, if you set another superuser)
- so when testing a PostgreSQL cluster using Patroni, with multiple hosts (a primary and a given number of replicas, synchronous or not), usually, pgSimload is run in 2 separate terminals, one to load data, and the other, to monitor things in Patroni
 - note the Patroni-monitoring mode can have added information thanks to the Replication_info set to nogucs or <list of gucs separated by a comma (e.g "synchronous_standby_names, synchronous_commit, work_mem") in the patroni.json config file passed as an argument to -patroni patroni.json> parameter. If set to nogucs, no extra GUCs are shown, only the info from pg_stat_replication will be
- demo Crunchy Postgres, a fully Open Source based PostgreSQL distribution using extensively Ansible
- demo Crunchy Postgres for Kubernetes, a fully Open Source based PostgreSQL distribution to run production workloads in Kubernetes

4 Examples given

You can find several examples of usage in the main directory of the project under examples/.

4.1 examples/simple/

This example is the simpliest one:

Prerequisites

- PostgreSQL Server, any version shoud work
- user has LOGIN capabilities

No creation of tables and others are needed, so there's no need to call for -create <create.json > or such. pgSimload accepts the omission of parameter -create.

Here's something for the user to understand:

- script.one_liner.sql contains a simple "select 1;" with a sleep(1) on the same line,
 where
- script.two_liner.sql contains the same, but each command on one line.

When you pass one or the other in the -script <script.sql> parameter, you'll see the difference in counting "statements": actually, pgSimload counts statements as one per line, because the parsing method for the -script <script.sql> thing is very basic.

We did tests with a JSON version of it like -script <script.json>: it adds an overhead and makes everything unreadable in the end. And pgSimload is not designed to have "exact" parameters and results, the thing is having "some data", and watch differences in between different configurations of PG infrastructure, or parameters (GUCS), etc.

Once pgSimload has been compiled **and** the config.json adapted to suit your needs, this could be used as simple as:

```
1 $ pgSimload -config config.json -script script.one_liner.sql
2 $ pgSimload -config config.json -script script.two_liner.sql
```

4.2 examples/PG_15_Merge_command/

This example is to test new MERGE command in PG 15.x

Prerequisites

- PostgreSQL Server version 15+ stand-alone or not
- a user called owning a schema name "test" (please adapt config.json file to match your needs here)
- user has LOGIN capabilities

Creates a schema with 3 tables in create.json.

script.sql will:

- create sample data in test.station_data_new
- merge that data in test.station_data_actual
- merge test.station_data_actualinto test.station_data_history

As per jpa's blog article on MERGE

Once pgSimload has been compiled and the binary placed in some dir your \$PATH points to, this could be used as simple as:

```
1 $ pgSimload -config config.json -create create.json -script script.sql
```

The watcher. sh is a plain psql into watch to get some live stats on the database. You may have to adapt it to match your usage. We've added 2 flavours.

The first show some data, nice to have in a separate terminal (use tilix while you demo!):

```
1 $ sh watcher.sh query
```

The second shows a nice histogram of the data, the query is slightly more complex and heaven tho:

```
1 $ sh watcher.sh histogram
```

4.3 examples/testdb/

This is another example that shows one can:

- create multiple different create.json files to match different scenarios, adding different things like in create.json, create.delete.json, create.delete.vacuum.json, etc. to pass to the paramter create
- create multiple different script.sql, insert.sql, etc.. to pass to the parameter create

If you have a PostgreSQL cluster where you want to test as an example:

- write activity to the primary and
- read activity to the secondary

Then you'll need 2 different files for credentials one to your primary, on let's say port 5432, another one to your secondary (or pool of secondaries, if you're using pgBouncer on a different port, or just HAProxy or anything else to balance to different PostgreSQL replicas, on let's say, port 5433).

You'll need also 2 different SQL script files to run read/write operations on the primary, and obviously, read/only operations to the secondary (or group of secondaries).

Finaly, you will have to run twice pgSimload, in 2 different terminals, to handle boths scenarios at the same time.

We give here a special example of the file session_parameters.json (you can name that like you want), as for you to use the special -session_parameters <session_parameter.json> if you want to modify the parameters of the session in which the script.sql queries will exectute. You can use this to set special values to a lot of configuration parameters that PostgreSQL allows to change within a session. As an example: work_mem, synchronous_commit, etc.

5 Overview of flags and parameters

Those tabulars show basic information about flags and parameters. For full documentation, please read next chapter "Reference: parameters and flags".

5.1 All modes: flags

All flags are optional and intended to run alone.

Name	Description
contact	Shows author name and email
help	Shows some help
license	Shows license
version	Shows current version of pgSimload

5.2 SQL-loop mode: parameters

Name	Mandatory	Optional	Value expected	Description
config	X		JSON file	Sets the PG connexion string (any user)
create		X	JSON file	Sets the SQL DDL to run once prior main loop

Name	Mandatory	Optional	Value expected	Description
script	Х		SQL text file	Sets the script to run inside the loop
session_parameters		X	JSON file	Sets special session configuration parameters

5.3 Patroni-monitoring mode: parameters

Name	Mandatory	Optional	Value expected	Description
config	X (if Replication_info is not empty)	Х	JSON file	Sets the PG connexion string (superuser)
patroni	X		JSON file	Sets parameters for this special mode

5.4 Session parameters template file creation

Name	Mandatory	Optional	Value expected	Description
config	X		JSON file	Sets the PG connexion string
create_gucs_templat * e			output name file	Sets the template file to create

6 Reference: parameters and flags

All flags and parameters can be listed executing pgSimload -h.

There are 2 different modes when executing pgSimload:

- SQL-loop mode to execute a script infintely on a given schema of a given database
- **Patroni-monitoring mode** to execute a monitoring on a given Patroni cluster. So you need one of such for this mode to be useful to you

Given the mode you choose, some parameters are mandatory or not. And the contexts of executions are different.

Before listing each, there are common parameters that can be used. Let's see those first.

6.1 Common flags and parameters

6.1.1 config (JSON file) [MANDATORY]

In the **SQL-loop mode** the "Username" set in the config.json can be any PostgreSQL user.

In the **Patroni-monitoring mode** the "Username" set in the config.json **has to be a superuser** in PostgreSQL, typically "postgres". Because we use special tricks to get the hostname of the PostgreSQL primary server.

"ApplicationName" is used to put a special "pgSimload" there, so the user can ps aux | grep [p]gSimload on any of the PostgreSQL server to isolate the process pgSimload uses... Or for any other SQL / bash command.

As per version 0.6 (June 2023), a valid config.json looks like this:

```
1 {
2    "Hostname": "localhost",
3    "Port" : "5432",
4    "Database": "mydbname",
5    "Username": "myusername",
6    "Password": "123456",
7    "Sslmode": "disable",
8    "ApplicationName": "pgSimload"
9 }
```

Most common values would be there either disable for non-SSL connexion or require for SSL ones.

[&]quot;Sslmode" has to be one among those described in Table 34.1. SSL Mode Descriptions.

6.1.2 contact (flag) [OPTIONAL]

Executing with only -contact will show you where you can contact the programmer of the tool.

This flag is not supposed to be run with other parameters or flags.

6.1.3 help (flag) [OPTIONAL]

Originally, "heredocs" were used in the main program to show this help, but it became too big to do such, it's better to have that doc in the current format you're reading, makes the source code lighter and that's cleaner IMHO.

So as per now, the execution of that -help is only to show where the current documentation is located. Actually, if you are reading this, that means wether you executed that flag...or that you find it by yourself. Kudos:-)

This flag is not supposed to be run with other parameters or flags.

6.1.4 license (flag) [OPTIONAL]

Executing with only -license will show you the license of this tool, currently licensed under The PostgreSQL License.

A full copy of the licence should be present aside the tool, in the main directory, in a file named LICENCE.md.

This flag is not supposed to be run with other parameters or flags.

6.1.5 version (flag) [OPTIONAL]

Executing with only -version will show you the current version of pgSimload. This is intended for general information of the users and also for any further packager of the tool in various systems.

Not supposed to be run with other parameters. No need to add a value to that flag.

6.2 SQL-loop mode parameters

The config flag is not listed down there, but is still **mandatory** to run in this mode, please read carefully informations upper in this documentation. On this mode, no particular "Username" has to be set in the config.json file.

6.2.1 create (JSON text file) [OPTIONAL]

If you need to create tables, or do anything prior to the execution of the main loop, you have to put your SQL commands in this JSON text file.

This script will be run only once prior the main loop on the script described above.

If you're want to execute pgSimload in SQL-loop mode on an existing database, on which you've adapted the SQL present in the script, then you don't need this feature. That's why it is optional.

To have a better idea of what's expected here, please refer to examples/PG_15_Merge_command /create.json or examples/testdb/create.json files.

6.2.2 script (SQL text file) [MANDATORY]

This file is in plain text and contains SQL statements to run, in the main loop of pgSimpload in the "SQL-loop mode".

It can be as simple as a "SELECT 1;". Or much more complex with SQL SQL statements of your choice separated by newlines. As an example of a more complicated example see examples/PG_15_Merge_command/script.sql.

Warning, as per version 0.6, the parsing is very basic for this script: each SQL statement is separated with; \n , so, this doesn't fit complex usages.

So consider limiting the content of those files with simple SQL commands, and not creating functions or other more complex things. If you need to create prior functions, do that with psql ... < create.sql prior to run pgSimload.

6.2.3 session_parameters (JSON text file) [OPTIONAL]

This parameter lets you tweak the PostgreSQL configuration that can be specified in a session. This

can be everything your PostgreSQL version allows, and we let you define proper values for proper parameters.

Every parameter you specify here will be passed at the beginning of the session when the SQL-loop is executed. So everything will be executed accordingly to those parameters in that session.

As an example, you can tweak work_mem in a session, or synchronous_commit, depending your PostgreSQL configuration and version.

The format of the JSON file has to be the following:

```
1
  {
    "sessionparameters": [
2
3
        "parameter" : "synchronous_commit"
4
       ,"value" : "remote_apply"
      },
6
     { "parameter" : "work_mem"
7
       ,"value" : "12MB"
8
9
      }
10
   ]
11 }
```

You can add as many parameters you want in that file, from one to many.

At the moment, we don't check if the parameter and values are OK. As an example, if you set a value for an unknown parameter, you will have this output when running pgSimload:

```
1 The following Session Parameters are set:
2    SET synchronous_commit TO 'remote_apply';
3    SET work_mem TO '12MB';
4    SET connections TO 'on';
5
6    2023/06/27 14:24:38 ERROR: unrecognized configuration parameter "connections" (SQLSTATE 42704)
```

Or if you set a right name, but in the wrong context you could have this too:

```
1 The following Session Parameters are set:
2    SET synchronous_commit TO 'remote_apply';
3    SET work_mem TO '12MB';
4    SET log_connections TO 'on';
5    2023/06/27 14:41:20 ERROR: parameter "log_connections" cannot be set after connection start (SQLSTATE 55P02)
```

And finaly, if you think you set proper values but it seems that nothing is read from the brand new session_parameters.json you just created, like in:

```
1 The following Session Parameters are set:
```

```
2
3
4 Now entering the main loop, executing script "./examples/testdb/script.
sql"
5 Script statements succeeded : |00000060|
```

... that's because you have probably a error in the JSON file, or maybe you changed the keyword "sessionparameters": : don't do that, it's expected in pgSimload to have such keyword there. It is also expected that your JSON file here is valid, like given in the example file given in examples/testdb/session_parameters.json

6.3 Patroni-monitoring mode flag and parameters

To use have pgSimload act as a small Patroni-monitoring tool in a side terminal, all you have to do is to create a patroni.json file in the following format. Note that the name doesn't matter much, you can name the way you want.

6.3.1 patroni (value) [MANDATORY]

When this paramter is set (-patroni <patroni.json>), you're asking pgSimload to run in Patroni-monitoring mode. This parameter is used to give to the tool the relative or complete path to a JSON file formated like the following (note: you can find a copy of this file in examples/patroni monitoring/:

```
$ cat patroni.json
2 {
        "Cluster" : "mycluster",
"Remote_host" : "u20-pg1",
"Remote_user" : "postgres",
"Use_sudo" : "no",
3
4
5
6
        "Ssh_private_key" : "/home/jpargudo/.ssh/id_patroni",
7
        "Replication_info" : "server_version, synchronous_standby_names,
8
             synchronous_commit,work_mem",
        "Watch_timer" : 5,
"Format" : "l
9
         "K8s_selector" : "list",
11
12 }
```

Cluster

You must specify here the Patroni's clustername. You can generally find it where your have Patroni installed in /etc/patroni/<cluster_name>.yml or inside the postgresql.yml.

Remote_host

You have to set here the ip (or hostname) where pgSimload will ssh to issue the remote command patronictl as user patroni_user (see up there).

Remote_user

This is an user on one of the PG boxes where Patroni is installed. That one you use to launch Patroni's patronictl. Depending the security configuration of your PostgreSQL box, Patroni could run with the system account PostgreSQL is running with, or another user. This one may have need to use sudo or not. Again, that all depends on your setup.

Use sudo

If the previous user set in **Remote_user** needs to use sudo before issuing the patronictl command, then set this value to "yes".

Ssh_private_key

Since ssh-ing to the Remote_host IP or (hostname if it's enabled in your DNS) need an SSH pair of keys to connect, we're asking where is that private key. It can be as simple as /home/youruser/. ssh/id_patroni. Beware not to set here the public key, because we need the private one.

Also, we assume you did the necessary thing on SSH so that user can SSH from the box where pgSimload is running to the target host, specifically, that the public key of your user is present in the ~/.ssh/authorized_keys of the taget system and with the matching **Remote_user**.

Replication_info

Thanks to this feature, pgSimload can show extra information about replication. This is usefull if Patroni doesn't do "everything in HA", like the SYNChronous replication, that can be handled by PostgreSQL itself, thanks to the synchronous_commit and synchronous_standby_names parameters. It can also adapt in other scenarios, or just to show the server_version, whatever you want!

If you don't need this extra information, to disable it, just set it to an empty string in the JSON like:

```
1 [...]
2 "Replication_info": "",
3 [...]
```

If disabled, the "Replication information" no extra information will be shown after the output of the patronictl ... list command.

If you want to activate it, like Replication_info is anything different to an empty string, **be sure you also provide** a -config <config.json> parameter, pointing to a file where superuser postgres connection string is defined. So that in this config file, "Username" should be set to "postgres", and the PG box name and port should be directly set.

So there's 2 ways to activate this feature described above.

If you want to activate it, but want pgSimload to only show othe output some extra information from pg_stat_replication system table, then you set the special value "nogucs" like:

```
1 [...]
2 "Replication_info": "nogucs",
3 [...]
```

The other way to activate it is to ask pgSimload to show also settings from the PostgreSQL Primary the whole query will be sent to. In this case, you have to set there all the GUCs you want to be shown, you just have to name those settings separated by a comma in the value of that JSON's fied.

This can be something like:

You can look at examples given at examples/patroni_monitoring/.

Watch_timer

You can ask for the output in the Patroni-monitoring mode to be like a bash "watch" command: it will run every x seconds you define here.

If you want the tool to issue commands each 5 seconds, then set this parameter to simply 5. Since patronictl command can take several seconds to run, the value you set here will be computed by the program to match your request, with timers to take into account the time of execution. So then the tool will iterate a bit before going the closest possible to your match your request.

If the value is less than 1, pgSimload will assume you only want to run it once in the Patroni watcher mode.

Format

The patronictl command offers two modes to list the nodes:

- list will order nodes output by name while
- topology will show the Primary first, so the order may change if you do a switchover of a failover

K8s_selector

This parameter has to be set only if your PostgreSQL Patroni cluster is in Kubernetes.

The value of this field must be what you'd put in the "selector" chain of that particular kubectl command, if you want to get the name of the pod where the current PostgreSQL primary is executing into:

So pgSimload knows the pod where the Primary PostgreSQL server is running.

The usage of pgSimload in "Patroni monitoring mode" in Kubernetes **has requirements**, we urge you to read carrefully the documentation you can access at examples/patroni_monitoring/README.md!

In short, if the Patroni monitoring mode has to be executed on a cluster of PostgreSQL servers in Patroni, the only relevant paramters in the patroni.json file would then be:

- Replication_info: can be set to an empty string (""), if you dont need it, nogucs or list of GUCs separated by a coma> if you want those informations to be shown. In the later case, you'll need then to run mandatorily with the -config config.json parameter too. In than file you'll set a superuser connection (e.g. "postgres" username)
- Watch_timer has to be set to a value >1 otherwise it will only runs once
- Format has to be set either to list or topology. In list, nodes will be ordered by name, while in topology, the Primary will be shown first
- K8s_selector has we already seen up there
- all others parameters won't apply, so you can leave them empty ("")

6.4 Session parameters template file creation

6.4.1 config (value) [MANDATORY]

Same as before, you define in that config.json file (or whatever the name, but it has to be a valid JSON here: see previous examples) the connection that will be used to query the pg_settings system

You can use whatever user here (i.e. superuser or not), because we only gather the parameters in the user context as per pg_settings PostgreSQL documentation.

6.4.2 create_gucs_template (value) [MANDATORY]

This parameter should have been named create_session_parameters_template_file to understand what it does...

Here, pgSimload will connect to a given PostgreSQL server as described in the mandatory -config <config.json> parameter you have to use too. Then, it will query the system view pg_settings to gather the name and the value (aka setting) of each parameter than can be changed in a given session.

Then it will output that in file which format is expected by pgSimload to be passed to the parameter -session_parameter.

Beware that those parameters change from one major PostgreSQL version to another, so likely a file you previously generated, then edited to suit your needs, on a version 15 won't work on a version 12.

Also, since ALL parameters in the context user will be gather (see pg_settings for details), there will be likely many dozens of parameters here. As an example, as per version 15, it's more than 130 parameters...

Since you probably won't need all of these, most likely, you run that command once to have every parameter in the generated template, then you edit it to remove all uncessary parameters. You'll have then your own template you can use in different scenarios, creating as many session_parameters. json you need, to be tested.