

# Change Log for Rate Fitting Code

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25 August 2011

## 1 Notes

- Src is at `~/Work/Rates/Fit_pkg`

## 2 Changes

### 2.1 2011-08-25 Thu (LLS)

- Fixed problems with `adjustrates`.
- Added diagonal regularization for temp and water

### 2.2 2011-08-15 Mon (LLS)

- Jacobians now normalized.  $\max(\text{sum}(\text{abs}(\text{temp}))) = \max(\text{sum}(\text{abs}(\text{water}))) = \max(\text{abs}(\text{gasi}))$  Jacobians. See `get_jac.m`
- `get_jac.m` no longer squeezes the jacobians for the parameters being fit. This is left to `retrieval.m`. This way, the jacobians outputted by `get_jacs.m` are “full” and can be used by `adjustrates.m` to remove the effects of some gas using a known rate.
- Three lambda's are now used. gas/water/temp. The original lambda is now only used if you invoke regularization only on the diagonal.
- Test directory using new covariance matrices (L1 with larger numbers for  $i=1:6$ ) and a true `apriori_zero` = all zeros.
- In `oemlls.m` separated mean square variance of the fitted coefficients in three separate parts: gases, water, temp. Save in `driver.oem.coeffvarxxx` where `xxx` = gas,water,temp. This allows you to do the L-curve separately for each coefficient type.

### 2.3 2011-08-08 Mon

- Removed “cluster” stuff.
- Many many code changes, mostly to simplify.
  - Large speed-up by doing multiple `pinv`'s only once and saving
  - Assume cov input matrix is the regularization matrix!
  - Used `inv` instead of `pinv` for the gain matrix
- Both the Jacobian file and the covariance file now must have same ordering. [gases columns (1-5) `Tsurf` `wv(1:97)`, `t(1:97)`]

- Removed lamba loop, user creates this loop externally if needed
- For now, put gain and ak matrices in driver output
- Fixed qrenorm to reflect that T-Jacobians were scaled by 0.01, not 0.1
- Changed find<sub>covderivoperators</sub> to remove qrenorm scaling on L matrices
- lambda input is now not scaled by  $10^{(\text{lambda})}$
- NOTE: testing now done with a L1 matrix (unity lambda2)

## 2.4 2011-08-05 08:00

- the Cluster directory allows you to put the below job on the cluster, using run\_tara

## 2.5 2011-08-04 21:00

- set\_default\_struct is used to set defaults
- these are then overwritten by calling driver = override\_defaults(driver,ix); where “ix” is the latbin you want to run. So for example a loop to do all latbins would be

```
for ix=1:36
    clear driver;
    run_retrieval;
end
```

## 2.6 2011-08-02 05:05

- jacobian file now has the qrenormalization information

## 2.7 2011-08-01 21:50

- driver.map\_jac2cov allows you to map indices from Jacobian to Covariance

## 2.8 2011-08-01 14:14

- Covariance matrix format changed; no top level structure name. oem\_11s.m changed to reflect this, see b0.
- Apriori format changed; no top level structure name. oem\_11s.m changed to reflect this, see zset, and zstd.
- Changed debugging approach, using Matlab code like:

```
if driver.debug
    addpath Debug
    debug_cov
    rmpath Debug
end
```

These scripts are in the Debug directory.

- Created directory Changelog to hold this file and its various output formats.