User guidelines for Advanced Model Diagnostics with ss3diags

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1 Getting started

This vignette explains functions for applying advanced model diagnostics for Stock Synthesis model with the ss3diags R package for model. The ss3diag R package accompanies the paper "A cookbook for using model diagnostics in integrated stock assessments" by Carvalho, Winker et al. (2021). The ss3diags package builds on R package r4ss (Taylor et al. 2021), which is designed to support the use of the Stock Synthesis software (Methot and Wetzel, 2013).

1.1 Installation

Both ss3diags and r4ss can be installed from gihtub using library(devtools):

```
installed.packages("devtools")

devtools::install_github("JABBAmodel/r4ss")

devtools::install_github("JABBAmodel/ss3diags")

library(r4ss)
library(ss3diags)
```

1.2 Loading built-in example data

The package contains two examples of Stock Synthesis assessments as presented in Carvalho, Winker et al. (2021).

1.2.1 The Pacif hake (Merluccius productus) base case model

2020 Pacific hake/whiting Stock Synthesis assessment. Joint Technical Committee of the U.S. and Canada Pacific Hake/Whiting Agreement, National Marine Fisheries Service and Fisheries and Oceans Canada

```
data("pac.hke")
```

Individual list objects generated using R package r4ss:

• ss3phk: output from Stock Synthesis as read by r4ss::SS_output()

```
dir.path = "C:/Users/henni/Dropbox/ss3diags_demo/PacificHake"
ss3pke = r4ss::SS_output(file.path(dir.path, "Reference_Run"))
```

• retro.phk: list of retrospective runs with r4ss:SS_doRetro() as read by r4ss::SSgetoutput()

In this case the retrospective runs, produced with r4ss:SS_doRetro()(see recipe below), were located in the subfolders /Retro_Reference_Run and named "retro0", "retro-1", "retro-2" and so on. To load those at once we use the r4ss::SSgetoutput(). In this case we conducted eight retrospective runs, where "retro0" corresponds to the "Reference_Run" and "retro-1" to "retro-7" are "peels". To assign the model names, we there specify start.retro = 0 and end.retro = 7 below.

```
start.retro = 0
end.retro = 7
retro.runs = "Retro_Reference_Run"
# load models
retro.phk <- r4ss::SSgetoutput(dirvec=file.path(dir.path,
retro.runs,paste0("retro",start.retro:-end.retro)))</pre>
```

Again, note that list object retro.phk[[1]] ("retro0") corresponds to the reference run ss3phk

aspm.phk: Comprises of two runs the "Reference_Run" and the "ASPM", which can be loaded together using r4ss::SSgetoutput() and then summarized with r4ss::SSsummarize()

1.2.2 The ICCAT North Atlantic shortfin make (Isurus oxyrinchus) reference model

```
data("natl.sma")
```

Individual list objects generated using R package r4ss:

- ss3sma: output from Stock Synthesis as read by r4ss::SS_output()
- retro.sma: list of retrospective runs with r4ss:SS_doRetro() as read by r4ss::SSgetoutput()
- aspm.sma: list of Stock Synthesis referece and aspm created with r4ss::SSsummarize()

2 Model Diagnostics with ss3diags

2.1 Plotting residual diagnostic with ss3diags

The plotting options are kept mostly to those provided by r4ss. Just like with r4ss, if, for example, 'SSplotRunstest() called with no further specifications several windows will open, which depends in this case on the number abundance indices.

SSplotRunstest(ss3sma)

The runs test is a nonparametric hypothesis test for randomness in a data sequence that calculates the 2-sided p-value to estimate the number of runs (i.e., sequences of values of the same sign) above and below a reference value. the runs test can diagnose model misspecification using residuals from fits to abundance indices (Carvalho et al. 2017). It can also be applied to other data components in assessment models such as the mean-length residuals and mean-age residuals. In addition, the three-sigma limits can be considered to identify potential outliers as any data point would be unlikely given a random process error in the observed residual distribution if it is further than three standard deviations away from the expected residual process average of zero.

To visicalized the runs test for multiple indices, it is suggested to make use of the function sspar() and the option plot add=TRUE. sspar() facilitates setting the graphic parameters so that they are suitable for ss3diags plots and option add=TRUE prevents the plotting functions from over-writing sspar().

```
sspar(mfrow=c(3,2),plot.cex=0.8)
rt = SSplotRunstest(ss3sma,add=T,verbose=F)
```

It is also possible to select the indices that should be plotted. For example, in this case we may want to exclude CPUE2 as it was not fitted (zero weight to the likelihood). This also creates space to add another plot, such as joint-residual SSplotJABBAres to summarize all selected indices.

```
sspar(mfrow=c(3,2),plot.cex=0.8)
rt= SSplotRunstest(ss3sma,add=T,indexselect = c(1,3:6),legendcex = 0.8,verbose=F)
jr = SSplotJABBAres(ss3sma,add=T,indexselect = c(1,3:6),legendcex = 0.55,verbose=F)
```

The default for SSplotRunstest() and SSplotJABBAres() is plot the residual runs for the abundance indices, but it is also possible to do the plot plot for the composition data by specifying subplots="len" (or "age")

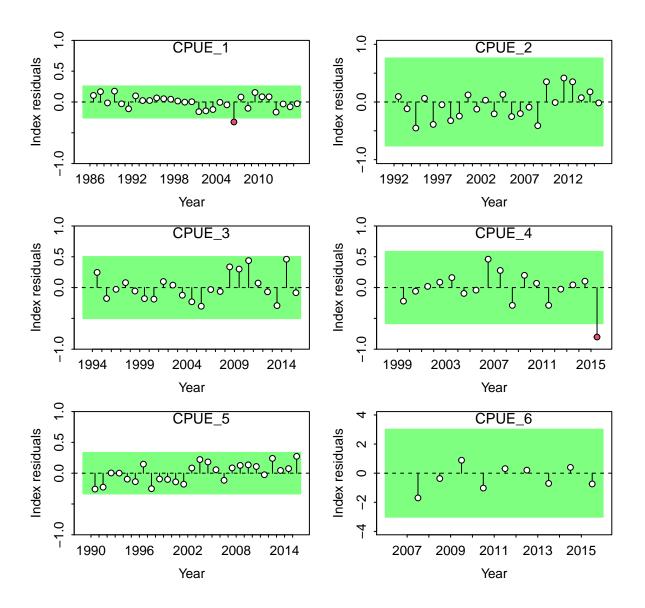


Figure 1: Runs test plots for CPUE index fits. Green shading indicates no evidence (p = 0.05) and red shading evidence (p < 0.05) to reject the hypothesis of a randomly distributed time-series of residuals, respectively. The shaded (green/red) area spans three residual standard deviations to either side from zero, and the red points outside of the shading violate the 'three-sigma limit' for that series.

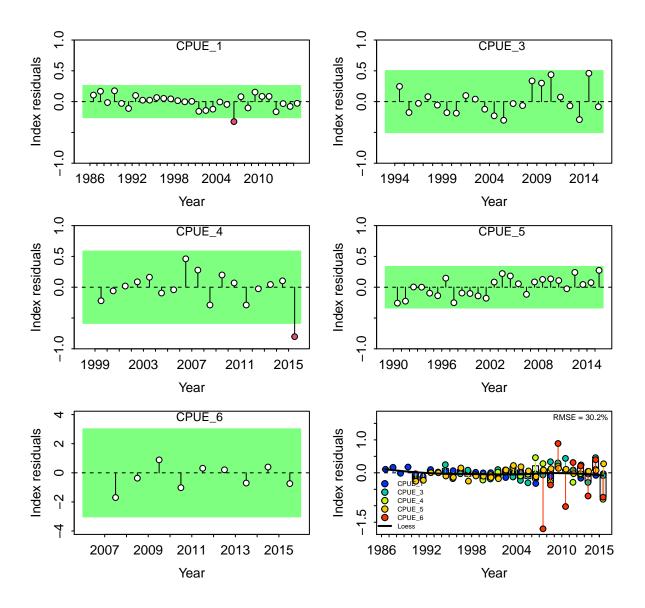


Figure 2: Runs test plot and Joint residual plot for fits to CPUE indices, where the vertical lines with points show the residuals, and solid black lines show loess smoother through all residuals. Boxplots indicate the median and quantiles in cases where residuals from the multiple indices are available for any given year. Root-mean squared errors (RMSE) are included in the upper right-hand corner of each plot.

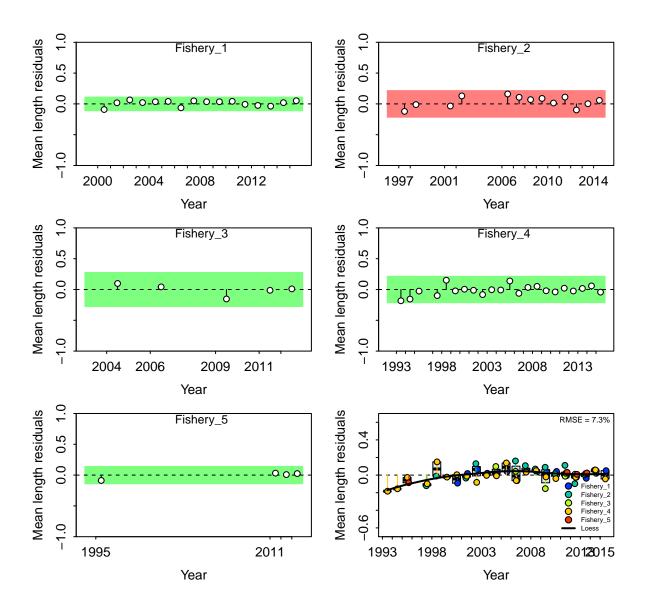


Figure 3: Runs test plot and Joint residual plot for mean lengths from fits length composition data

Several diagnostic tests can also be called without plotting, e.g. to facilitate automated processing

```
rti= SSrunstest(ss3sma,quant="cpue",verbose=F)
rtl= SSrunstest(ss3sma,quant="len",verbose=F)
rbind(rti,rtl)
          Index runs.p
                         test sigma3.lo sigma3.hi type
         CPUE_1 0.069 Passed -0.2605783 0.2605783 cpue
   2
         CPUE_2 0.717 Passed -0.7630777 0.7630777 cpue
   3
                0.229 Passed -0.5033328 0.5033328 cpue
   4
         CPUE 4
                0.406 Passed -0.5868380 0.5868380 cpue
   5
         CPUE 5
                0.065 Passed -0.3369695 0.3369695 cpue
         CPUE 6 0.870 Passed -3.0226356 3.0226356 cpue
   6
   7
     Fishery 1
                0.127 Passed -0.1103741 0.1103741
     Fishery_2
                0.040 Failed -0.2156036 0.2156036
     Fishery_3 0.331 Passed -0.2759992 0.2759992
                                                    len
   10 Fishery_4
                0.806 Passed -0.2141490 0.2141490
                                                    len
   11 Fishery_5
                0.159 Passed -0.1407228 0.1407228
                                                    len
```

The pacific hake assessment provides an example of fits to age composition instead of length composition data, which can plotted by specifying subplots="age"

```
sspar(mfrow=c(2,2),plot.cex=0.8)
rti = SSplotRunstest(ss3phk,add=T,legendcex = 0.8,subplot="cpue",verbose=F)
rta = SSplotRunstest(ss3phk,add=T,legendcex = 0.8,subplot="age",verbose=F)
jra = SSplotJABBAres(ss3phk,add=T,legendcex = 0.7,subplot="age",verbose=F)
```

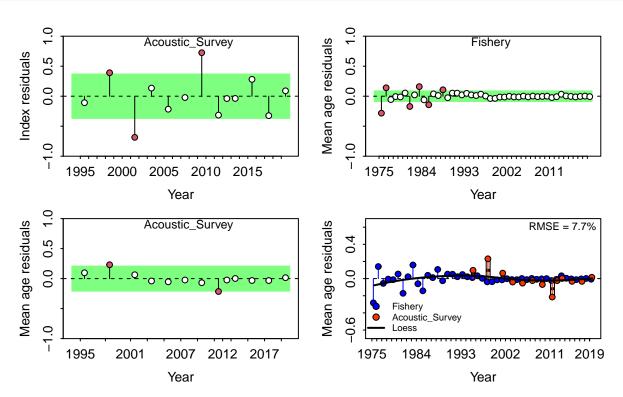


Figure 4: Runs test plot and Joint residual plot for a survey abudance index and mean ages from fits to survey and fisheries dependent age-composition data

2.2 Ploting Retrospective and Forecast bias

Retrospective analysis is commonly used to check the consistency of model estimates, i.e., the invariance in spawning stock biomass (SSB) and fishing mortality (F) as the model is updated with new data in retrospect. The retrospective analysis involves sequentially removing observations from the terminal year (i.e., peels), fitting the model to the truncated series, and then comparing the relative difference between model estimates from the full-time series with the truncated time-series.

In Stock Synthesis, retrospective analysis can be routinely implemented in Stock Synthesis using r4ss:SS_doRetro() available in r4ss as demonstrated in Section 3.1. ss3diags then provides the function SSplotRetro() to visualize the retrospective patterns of SBB and F and to compute the associated Mohn's rho value (i.e. retrospective bias). This would require to first load the retrospective runs (Section 1.2), which are already inbuilt into ss3diags in this case. The next step is summarize the list of retrospective runs using r4ss::SSsummarize().

```
retroI.phk <- r4ss::SSsummarize(retro.phk,verbose=F)</pre>
```

We use notation "retroI" because r4ss::SSsummarize() summarizes the modelled quantaties and abudances indices, but length or age composition data, which becomes relevant again in the next Section. With summarized object retroI.phk it is now possible to produce some basic retrospective plots.

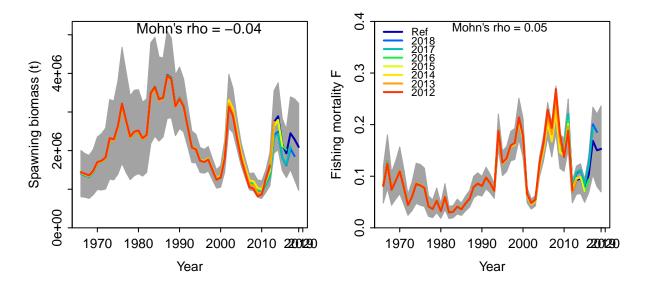


Figure 5: Retrospective analysis of spawning stock biomass (SSB) and fishing mortality estimates for Pacific hake conducted by re-fitting the reference model (Ref) after seven years, one year at a time sequentially. Mohn's rho statistic are denoted on top of the panels. Grey shaded areas are the 95 % confidence intervals from the reference model in cases where the analysis was run with Hessiean

Retrospective analysis is useful to evaluate how consistent the modeled quantities are in retrospect. However, providing fisheries management advice requires predicting a stock's response to management and checking that predictions are consistent when updated by new data in the future. A first, intuitive extension of the retrospective analysis is to assess potential forecast bias by adding the additional step of forward projecting quantities, such as SSB, over the truncated years. A desirable feature of Stock Synthesis is that forecasts are automatically done when using r4ss:SS_doRetro().The forecasts are based on the settings in

'forecast.ss', which are also evoked when conducting future projections with the same model, only that the observed catches are used for the retrospective forecasts. Retrospective forecasts (i.e. hindcasts) with Stock Synthesis is therefore only a matter of visualization, which can be done by setting the SSplotRetro() option forecast=TRUE.

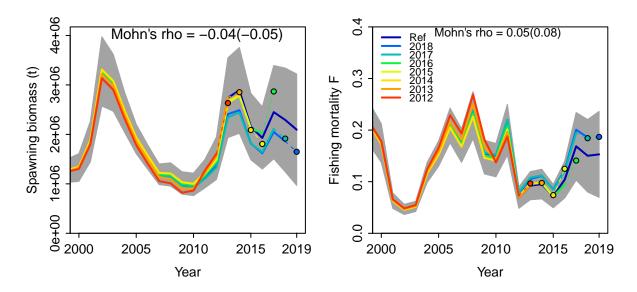


Figure 6: Retrospective results shown for the most recent years only. Mohn's rho statistic and the corresponding 'hindcast rho' values (in brackets) are now printed at the top of the panels. One-year-ahead projections denoted by color-coded dashed lines with terminal points are shown for each model.

The statistics in from the retrospective analysis with forecasting, including mohn's rho and forecast bias, can also be called without plotting using the function SShcbias()

```
SShcbias(retroI.phk,quant="SSB",verbose=F)
     type
              peel
                             Rho ForcastRho
   1
      SSB
              2018 -0.189511890 -0.21290187
   2
      SSB
              2017 -0.138534518 -0.16468977
              2016
   3
      SSB
                     0.048946154 0.17079324
      SSB
              2015 -0.017403026 -0.06301299
   4
      SSB
   5
              2014 -0.032761073 -0.03298098
   6
      SSB
              2013
                     0.001623057 -0.01285596
      SSB
              2012
                    0.059944948 -0.03946034
   7
   8
      SSB Combined -0.038242335 -0.05072981
SShcbias(retroI.phk,quant="F",verbose=F)
     type
              peel
                             Rho
                                 ForcastRho
   1
        F
              2018
                     0.235928805
                                  0.22117119
   2
        F
              2017
                     0.154620273
                                  0.22678319
   3
        F
              2016 -0.096502898 -0.16382344
        F
   4
              2015 -0.008902238
                                  0.20182834
        F
   5
              2014
                     0.034769554
                                  0.02664839
   6
        F
              2013
                     0.003690925
                                  0.02342468
        F
   7
              2012
                     0.049728943
                                  0.04693738
   8
          Combined
                     0.053333338
                                  0.08328139
```

2.3 Hindcast Cross-Validation and prediction skill

Retrospective forecasting provides an indication of the expected bias of in modelled quantities when updated with new data, but is not suitable for model validation, which should based on actual observations that are unknown to the model. To address this, Kell et al. (2016) proposed the used of hindcasting cross-validation techniques (HCXval) for stock assessment model validation, where observations (e.g. CPUE, length comps) are compared to their predicted future values. The key concept behind the HCXval approach is 'prediction skill', which is defined as any measure of the accuracy of a forecasted value to the actual observed value that is not known by the model (Kell et al., 2021).

Implementing the Hindcast Cross-Validation (HCxval) diagnostic in Stock Synthesis requires the same model outputs that are already produced with generated by r4ss:SS_doRetro() as described in Section 3.1. Therefore, there is no additional computationally intensive step needed for HCxval if conducted in conjunction with retrospective analysis. As robust measure of prediction skill, we implemented the mean absolute scaled error (MASE). In brief, the MASE score scales the mean absolute error (MAE) of forecasts (i.e., prediction residuals) to MAE of a naïve in-sample prediction, which is realized in the form of a simple 'persistence algorithm', i.e. tomorrow's weather will be the same as today's (see Eq. 3, p.5 in Carvalho and Winker et al. 2021). A MASE score > 1 indicates that the average model forecasts are worse than a random walk. Conversely, a MASE score of 0.5 indicates that the model forecasts twice as accurately as a naïve baseline prediction; thus, the model has prediction skill.

HCxval is implemented for Stock Synthesis using function SSplotHCxval() to produce the novel HCxval diagnostic plot and computes the MASE scores for CPUE indices, mean lengths or mean ages that have observations falling within the hindcast evaluation period.

Plotting HCxval for abudance indices requires the same step of summarizing the list of retrospective runs as for the retrospective analysis, which, of course, only needs be done once. Here, we summarize the retro runs for shortfin make.

```
retroI.sma <- r4ss::SSsummarize(retro.sma, verbose=F)
```

```
sspar(mfrow=c(3,2),plot.cex=0.8)
hci = SSplotHCxval(retroI.sma,add=T,verbose=F,ylimAdj = 1.3,legendcex = 0.7)
```

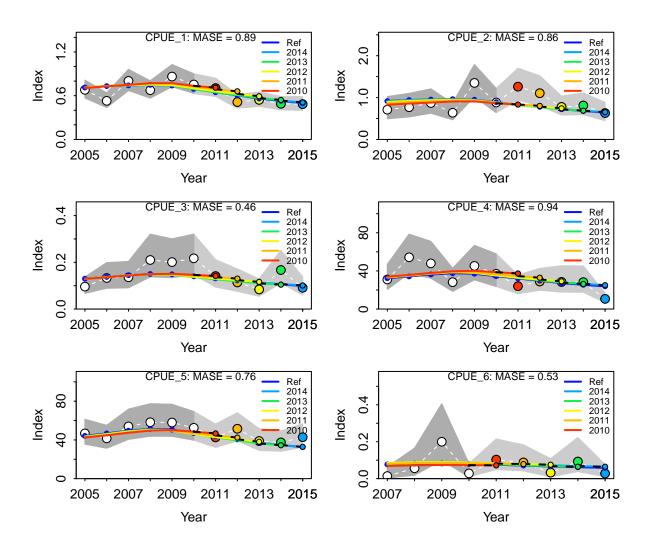


Figure 7: Hindcasting cross-validation (HCxval) results CPUE fits, showing observed (large points connected with dashed line), fitted (solid lines) and one-yearahead forecast values (small terminal points). HCxval was performed using one reference model (Ref) and five hindcast model runs (solid lines) relative to the expected CPUE. The observations used for crossvalidation are highlighted as color-coded solid circles with associated 95 % confidence intervals. The model reference year refers to the endpoints of each one-year-ahead forecast and the corresponding observation (i.e., year of peel \pm 1). The mean absolute scaled error (MASE) score associated with each CPUE

By comparison, the forecast length- and age-composition are somewhat hidden as "ghost files" of Stock Synthesis report.sso. To extract and summarize the composition in the form of oberved and expected mean lengths and mean ages, respectively, ss3diags provides the function SSretroComps().

```
retroC.sma = SSretroComps(retro.sma)

sspar(mfrow=c(1,2),plot.cex=0.8)
hcl = SSplotHCxval(retroC.sma,subplots="len",add=T,verbose=F,
ylimAdj = 1.3,legendcex = 0.7,indexselect = c(1,2))
```

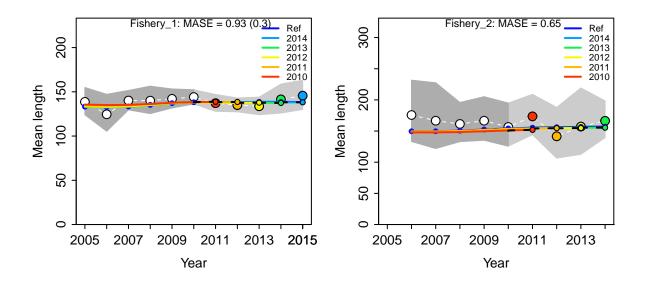


Figure 8: Hindcasting cross-validation (HCxval) results for mean lengths. Note that MASE values in breakets are adjusted MASE values for cases where naive predictions have a Mean-Absolute-Error below 0.1

Note that Figure 8, provides some additional, so called adjusted MASE values, in brackets. This gets invoked in cases where the inter-annual variation in the observed values is very small (default MAE < 0.1 for naive predictions $\log(y[t+1])-\log(y[t])$). The reasoning is that prediction residuals must be already very accurate to fall below this threshold. The adjusted MASE essential keep the naive prediction MAE denominator of the MASE to a maximum. Below we show the effect of changing adjustment threshold from the default MAE.base.adj = 0.1

to a larger value MAE.base.adj = 0.15

where MASE is the ratio of the mean absolute error of the prediction residuals MAE.PR to the residuals of the naive predictions MAE.base

```
mase1$MAE.PR/mase1$MAE.base
[1] 0.9265301 0.6504563

mase1$MASE
[1] 0.9265301 0.6504563

and MASE.adj

mase1$MAE.PR/pmax(mase1$MAE.base,0.1)
[1] 0.2981727 0.6504563

mase1$MASE.adj
[1] 0.2981727 0.6504563
```

Note that applying HCxval for composition data requires to correctly specify the composition data type that was fitted in the model. For example, age composition data need be specified as "age" in SSplotHCxval and SSmase, as shown below for the Pacific hake model.

```
retroC.phk = SSretroComps(retro.phk) # summarize comps
sspar(mfrow=c(1,2),plot.cex=0.8)
hcl = SSplotHCxval(retroC.phk, subplots="age", add=T,
verbose=F,ylimAdj = 1.3,legendcex = 0.7,indexselect = c(1,2))
SSmase(retroC.phk,quants="age")
   Computing MASE with all 7 of 7 prediction residuals for Index Fishery
   Computing MASE with only 4 of 7 prediction residuals for Index Acoustic_Survey
   Warning: Unequal spacing of naive predictions residuals may influence the interpretation of MASE
   MASE stats by Index:
               Index Season
                                MASE
                                         MAE.PR MAE.base MASE.adj n.eval
            Fishery 1 0.6319721 0.09063054 0.1434091 0.6319721
   1
   2 Acoustic_Survey
                         1 0.3556384 0.05181084 0.1456841 0.3556384
```

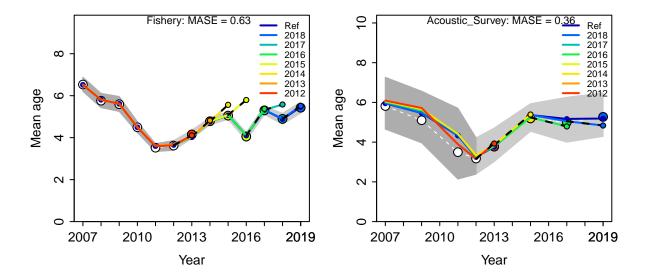


Figure 9: Hindcasting cross-validation (HCxval) results for mean lengths. Note that MASE values in breakets are adjusted MASE values for cases where naive predictions have a Mean-Absolute-Error below 0.1

2.4 Uncertainty with ss3diags

The management advice frameworks increasingly require translating the estimated uncertainty about the stock status into probabilistic statements (Kell et al. 2016). A classical example is the Kobe framework that is use in tuna Regional Fisheries Management Organisations around the world. The key quantities of interest are typically the ratios SSB/SSB_{MSY} and F/F_{MSY} . While is reasonably straight forward in Stock Synthesis to approximate uncertainty of individual quantities (e.g. SSB) from the asymptotic standard errors (SE) derived from the hessian matrix using the delta method, the joint distribution of SSB/SSB_{MSY} and F/F_{MSY}

requires to adequately account for the covariance structure between these two derived quantities. Joint distributions were therefore typical constructed using bootstrap or Markov Chain Monte-Carlo (MCMC) methods. However, these methods can be computationally intense and time consuming in integrated assessments.

As an alternative, ss3diags implements a rapid delta-Multivariate lognormal approximation with SSdeltaMVLN() to generate joint error distributions for SSB/SSB_{ref} and F/F_{ref} , where the ref may refer to MSY, but also, e.g., SSB_{40} and F_{40} . In Stock Synthesis, these ratios are determined by the derived quantaties Bratio and F, where either can take the form of ratios (e.g. F/F_{ref}) or absolute value (e.g. absF) depending on settings in the starter.ss file.

Let Bratio be $u = SSB/SSB_{ref}$, F be v = F, and $w = F_{ref}$ be the F reference point of interest (e.g. F_{MSY}), with $x = \log(u)$, $y = \log(v)$ and $z = \log(w)$, then the variance-covariance matrix VCM has the form:

$$VCM_{x,y,z} = \begin{pmatrix} \sigma_x^2 & cov_{x,y} & cov_{x,y} \\ cov_{x,y} & \sigma_y^2 & cov_{y,z} \\ cov_{x,z} & cov_{y,z} & \sigma_z^2 \end{pmatrix}$$

where, e.g., σ_x^2 is the variance of x and $cov_{x,y}$ is the covariance of x and y. Deriving those requires to conduct a few normal to lognormal transformations. First, we approximate the variances as:

$$\sigma_x^2 = \log\left(1 + \left(\frac{SE_u}{u}\right)^2\right)$$

where SE_u , SE_v and SE_z are the asymptotic standard error estimates for $u = SSB/SSB_{ref}$, v = F and $z = F_{ref}$.

The corresponding covariance for x and y, can then be approximated on the log-scale by:

$$COV_{x,y} = \log\left(1 + \rho_{u,v}\sqrt{\sigma_x^2\sigma_y^2}\right)$$

where $rho_{u,v}$ denotes the correlation of u and v.

To generate a joint distribution of $\tilde{u} = SSB/SSB_{ref}$, $\tilde{v} = F$ and $\tilde{z} = F_{ref}$, we use a multivariate random generator, available in the R package 'mvtnorm', to obtain a large number (e.g. nsim = 10,000) iterations, such that

$$JD(\tilde{u}, \tilde{v}, \tilde{w}) = \exp(MVN(\mu_{x,y,z}, VCM_{x,y,z}))$$

so that

$$S\tilde{S}B/S\tilde{S}B_{MSY} = \tilde{u}$$

and

$$\tilde{F}/\tilde{F}_{MSY} = \tilde{v}/\tilde{w}$$

The type of reference points dependents on settings in the starter.ss file that determine the derived quanties Bratio and Fvalue.

In the first example, we consider the ss3sma Stock Synthesis model, which was run with starter.ss settings that are commonly to produce target Kobe plot estimates of SSB/SSB_{MSY} and F/F_{MSY} in tRFMO assessments:

- 2 # Depletion basis: 2=rel SPBmsy; 3=rel X*SPB_styr; 4=rel X*SPB_endyr
- 2 # F_report_basis: 0=raw_F_report; 1=F/Fspr; 2=F/Fmsy; 3=F/Fbtgt

We provide the the function SSsettingsBratioF(ss3sma) to the starter.ss settings:

This function is also inbuilt in the SSdeltaMVLN() to prevent misleading results. To generate a joint MVLN of SSB/SSB_{MSY} and F/F_{MSY} , we can use the default options

```
mvln = SSdeltaMVLN(ss3sma,run="SMA")
starter.sso with Bratio: SSB/SSBMSY and F: (F)/(Fmsy)
```

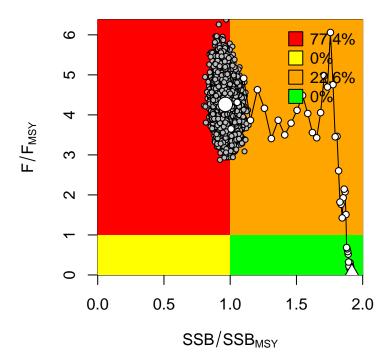


Figure 10: Kobe phase plots showing MVLN Kobe probability distributions of SSB/SSB_{MSY} and F/F_{MSY} for North Atlantic shortfin make Stock model.

The SSdeltaMVLN() output include the maximum the maximum likelihood estimates (mles) and the MVLN monte-Carlo distributions RE of SSB/SSB_{MSY} , F/F_{MSY} and F. Note the additional quanties SSB and RE are generated independely from lognormal distributions for practical reasons. These can be plotted by

```
sspar(mfrow=c(3,2),plot.cex = 0.7)
SSplotEnsemble(mvln$kb,ylabs=mvln$labels,add=T,verbose=F)
```

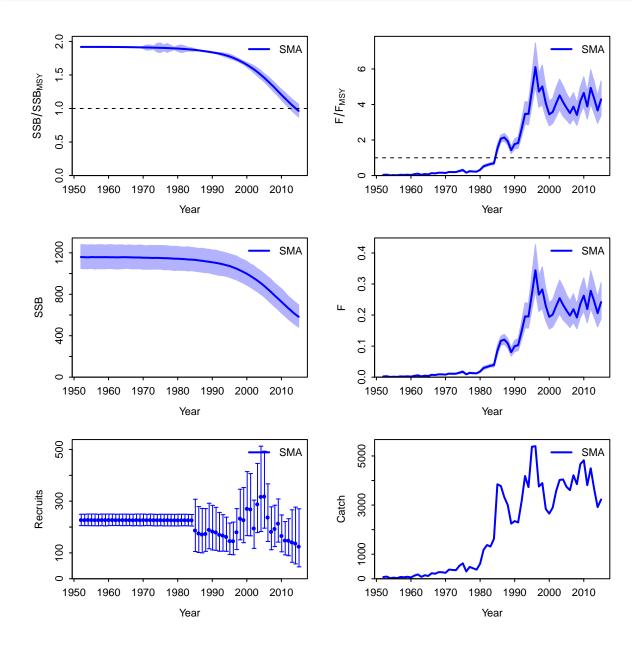


Figure 11: Distributions for SSB/SSB_{MSY} F/F_{MSY} , SSB, Recruitment and Catch trajectories for the North Atlantic shortfin make SS3 model

The SSdeltaMVLN() provides the option to set alternative Fref values, but this is only possible for the recommended starter.ss option 0 for F_report_basis. For option 2, SSdeltaMVLN() get an error if Fref where to be changed.

```
mvln = SSdeltaMVLN(ss3sma,run="SMA", Fref="Btgt")
    starter.sso with Bratio: SSB/SSBMSY and F: (F)/(Fmsy)
   Error in SSdeltaMVLN(ss3sma, run = "SMA", Fref = "Btgt"): Fref = Btgt option conflicts with (F)/(Fms
```

By comparison, the Pacific Hake base case model ss3phk is run with settings that are common in NOAA assessments, with Bratio set SSB/SSB_0 and F is typically kept at absolute quantity.

- 1 # Depletion basis: 2=rel SPBmsy; 3=rel X*SPB_styr; 4=rel X*SPB_endyr
- 0 # F_report_basis: 0=raw_F_report; 1=F/Fspr; 2=F/Fmsy; 3=F/Fbtgt

The management quantaties are in this case SSB/SSB_{40} and F/F_{spr40} , where the target of 40% is specified in the forecast.ss file.

```
0.4 # SPR target (e.g. 0.40) 0.4 # Biomass target (e.g. 0.40)
mvln = SSdeltaMVLN(ss3phk,run="Pac.Hake", Fref="SPR",plot=TRUE)
```



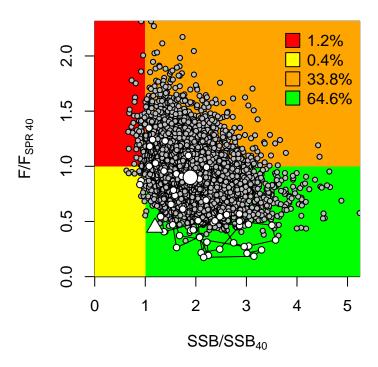


Figure 12: Kobe phase plots showing MVLN Kobe probability distributions of SSB/SSB_{40} and F/F_{SPR40} for North Atlantic shortfin make Stock model.

```
sspar(mfrow=c(3,2),plot.cex = 0.7)
SSplotEnsemble(mvln$kb,ylabs=mvln$labels,add=T,verbose=F)
```

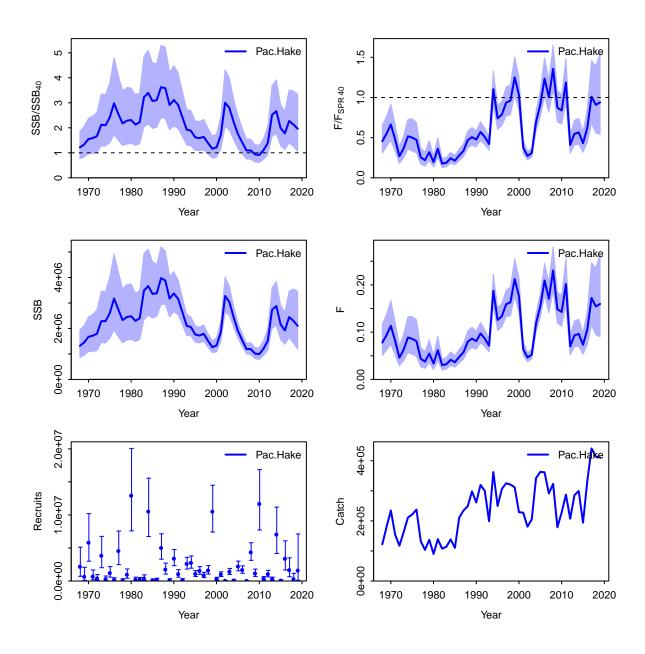


Figure 13: Distributions for SSB/SSB_{40} F/F_{SPR40} , SSB, Recruitment and Catch trajectories for the Pacific Hake SS3 model

It is important to note that MVLN approximation differs notably from MCMC posterior for this model as has been documented in Stewart et al. (2013) and Taylor et al. (2021). Such differences may be more likely to occur in cases where key parameters such as steepness h or natural M are estimated using informative priors, which can result in left skewed (non-lognormal) distributions of the benchmarks F_{ref} and B_{ref} .

In other instances, mismatches between the SSdeltaMVLN and MCMC may also be caused by poor performance of the latter due to poor regularization (Monnahan et al., 2019).

To facilitate comparison between the SSdeltaMLVN() and MCMC outputs, we provide the function SSdiagsMCMC(), which we illustrate on the example of the Stock Synthesis model for the ICES Gulf of Bothian Herring stock (ICES, 2021).

SSdiagsMCMC() requires loading both the report.sso and MCMC output in the posterior.sso file, where the MCMC was in this case run in the subfolder of the assessment file /mcmc.

```
# Alread loaded to ss3diags
ss3her = SS_output("gob_her")
mcmc.her = SSreadMCMC("gob_her/mcmc")
```

The options and output of SSdiagsMCMC() is largely identical to SSdeltaMVLN. In this case, the starter.ss is the same as for ss3phk, only the we use Fref = "Btgt" for illustration with SB_{50} and F_{SB50} for this illustration.

```
mvln = SSdeltaMVLN(ss3her,Fref="Btgt",plot=F,run="mvln")
    starter.sso with Bratio: SSB/SSBO and F: _abs_F

mcmc = SSdiagsMCMC(mcmc.her,ss3her,Fref="Btgt",plot=F,run="mcmc")
    starter.sso with Bratio: SSB/SSBO and F: _abs_F
```

Comparing detla-MVLN with MCMC simply requires combining the \$kb outputs by, e.g.,

```
sspar(mfrow=c(1,1),plot.cex = 0.8)
SSplotKobe(rbind(mvln$kb,mcmc$kb),joint=F,xlab=mvln$labels[1],
ylab=mvln$labels[2],fill=F)
```

```
1    Red  0.08002401
2    Orange  3.06091828
3    Yellow  0.00000000
4    Green 96.85905772

sspar(mfrow=c(2,2),plot.cex = 0.7)
SSplotEnsemble(rbind(mvln$kb,mcmc$kb),ylabs=mvln$labels,
add=T,subplots = c("stock", "harvest", "SSB", "F"),
verbose=F)
```

This works equally for joining a model ensemble.

Percent

Quadrant

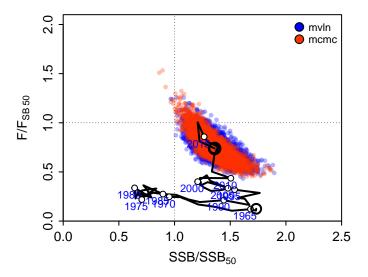


Figure 14: Kobe phase plot comparing MVLN and MCMC posterior distributions of SSB/SSB_{50} and F/F_{50} for the ICES Gulf of Bothia Herring SS3 model

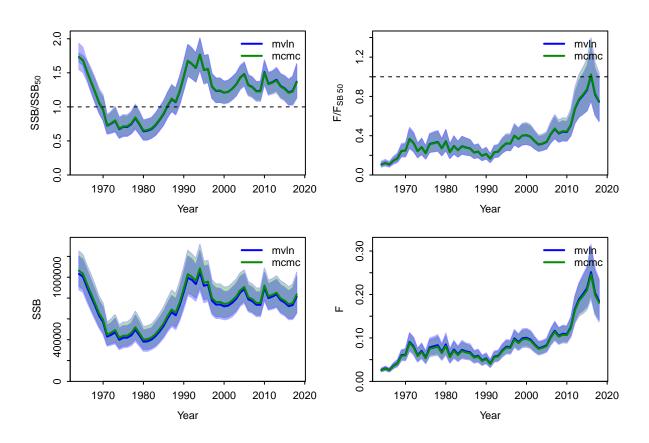


Figure 15: Comparison of MVLN and MCMC posterios for $SSB/SSB_{40}\ F/F_{SB40},\ SSB$ for the ICES Gulf of Bothia Herring SS3 model

3 Cookbook Recipies

3.1 Retrospetive with hindcasts

Retrospective analysis can be run for Stock Synthesis using the function r4ss::SS_doRetro() available in r4ss. This setup of the retrospective analysis has the advantage that forecasts are conducted automatically given the catch. This makes in possible to apply retrospective forecasting and hindcast cross-validations of observations based on the same output.

```
library(r4ss)
```

In the following we provide a step-by-step cookbook recipe for retrospective analysis in Stock Synthesis.

3.1.1 Identify restrospective period

The first step is specifying the range pf peels that will then determine the end.yr.vec of runs in r4ss::SS_doRetro()

```
start.retro <- 0  # end year of reference year
end.retro <- 7  # number of years for retrospective e.g.,</pre>
```

3.1.2 Identify the base directory

The second step is to specify the path directory that holds the folder with the base case run. In this case the Pacific Hake folder with a model folder 'Reference Run"

```
dirname.base = "C:/Users/henni/Dropbox/ss3diags_demo/PacificHake"
run = "Reference_Run"

model.run <- file.path(dirname.base,run)

model.run</pre>
```

DAT and CONTROL files

The third step is to specify the names of the data and control files. Note these files are named differently from the DATA.ss and CONTROL.ss. In this case

```
DAT = "phk.dat"
CTL = "phk.ctl"
```

The names of the DAT and CONTROL are declared on the top of the 'starter.ss', e.g.

#C Hake starter file phk.dat phk.ctl

3.1.3 Create a subdirectory for the Retrospectives

There are several ways to organise the retrospective output sturcture. Here we simply create a new subfolder for the retrospective runs output

```
dir.retro <- paste0(dirname.base,'/Retro_',run)
dir.create(path=dir.retro, showWarnings = F)</pre>
```

We also create a subdirectory for the retrospective model folders

```
dir.create(path=file.path(dirname.Retrospective, "retros"), showWarnings = F)
```

3.1.4 Copy model run files to new retrospective folder

```
file.copy(file.path(model.run, "starter.ss_new"), file.path(dir.retro, "starter.ss"))
file.copy(file.path(model.run, "control.ss_new"),
file.path(dir.retro, CTL))

file.copy(file.path(model.run, "data.ss_new"),
file.path(dir.retro, DAT))

file.copy(file.path(model.run, "forecast.ss"),
file.path(dir.retro, "forecast.ss"))

file.copy(file.path(model.run, "SS.exe"),
file.path(dir.retro, "SS.exe"))

# Automatically ignored for models without wtatage.ss
file.copy(file.path(model.run, "wtatage.ss"),
file.path(dir.retro, "wtatage.ss"))
```

3.1.5 Modify Starter.ss file

Modifying the Starter File helps to speed up model runs

```
starter <- readLines(paste0(dir.retro, "/starter.ss"))

#[8] "2 # run display detail (0,1,2)"
linen <- grep("# run display detail", starter)
starter[linen] <- paste0(1, " # run display detail (0,1,2)")
# write modified starter.ss
write(starter, file.path(dir.retro, "starter.ss"))</pre>
```

Execute retrospective runs

Now all is readt to run the retrospective analyses with r4SS function SS_doRetro Ideally, the runs should be done with Hessian to allow evaluating the retrospective trajectories with respect to confidence interval coverage of the reference model.

```
r4ss::SS_doRetro(masterdir=dir.retro, oldsubdir="",
newsubdir="", years=start.retro:-end.retro)
```

However, for larger models it might be desirable to shorten run times, by not inverting the hessian, using the option extras = "-nohess" (much faster)

```
r4ss::SS_doRetro(masterdir=dir.retro,oldsubdir="",
newsubdir="", years=start.retro:-end.retro,extras = "-nohess")
```

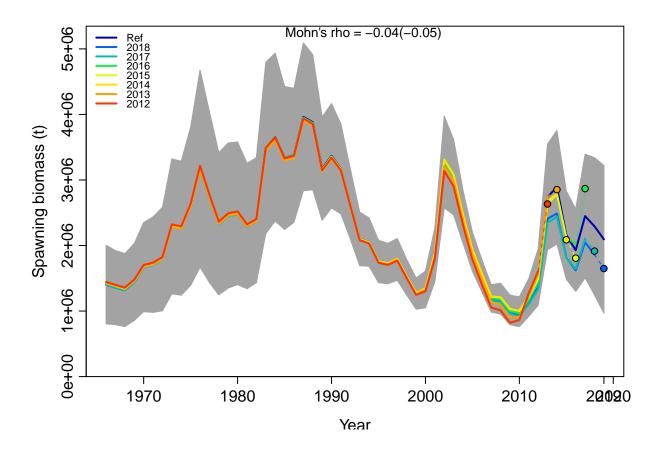
3.1.6 Read SS doRetro() output

It is often useful to save the retro model runs as .rdata file for further processing with ss3diags, considering that reading the models with r4ss::SSgetoutput() can be quite time-consuming for more complex models.

```
save(retro.phk, file=file.path(dir.retro,"retro.phk.rdata"))
```

3.1.7 Quick Check

```
library(ss3diags)
check.retro =r4ss::SSsummarize(retro.phk)
   Summarizing 8 models:
   imodel=1/8
    N active pars = 237
   imodel=2/8
    N active pars = 237
   imodel=3/8
    N active pars = 237
   imodel=4/8
    N active pars = 237
   imodel=5/8
    N active pars = 237
   imodel=6/8
    N active pars = 237
   imodel=7/8
    N active pars = 237
   imodel=8/8
     N active pars = 237
   Summary finished. To avoid printing details above, use 'verbose = FALSE'.
sspar(mfrow=c(1,1))
SSplotRetro(check.retro,forecast = T,add=T,legendcex = 0.7,legendloc = "topleft")
Plotting Retrospective pattern
```



 ${\tt Mohn's\ Rho\ stats,\ including\ one\ step\ ahead\ forecasts:}$

```
type
                         Rho ForecastRho
   SSB
           2018 -0.189511890 -0.21290187
1
2
  SSB
           2017 -0.138534518 -0.16468977
3
  SSB
           2016 0.048946154 0.17079324
4
  SSB
           2015 -0.017403026 -0.06301299
5
  SSB
           2014 -0.032761073 -0.03298098
           2013 0.001623057 -0.01285596
6
  SSB
7
   SSB
           2012 0.059944948 -0.03946034
   SSB Combined -0.038242335 -0.05072981
```