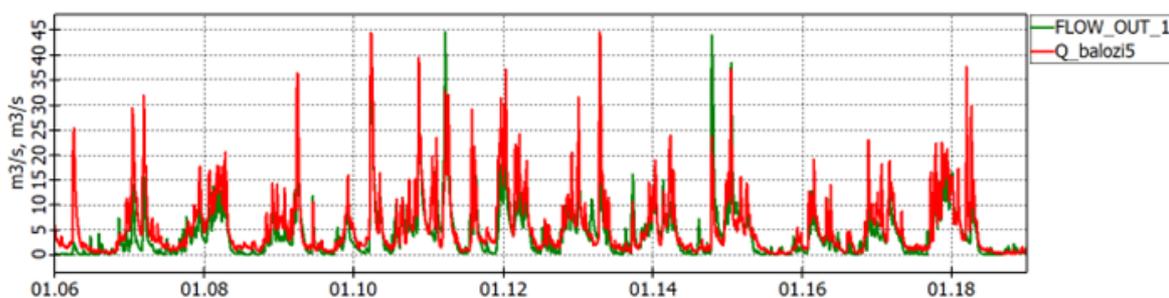


MODELLING RESULTS FOR REFERENCE AND BASELINE SCENARIOS

DELIVERABLE R3



Prepared within the LIFE GoodWater IP Action C1: “Development of the water quality and quantity system for the territory of Latvia”

Rīga, 2022

Modelling results for Reference and Baseline scenarios

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Summary

This document is a description of the electronic deliverable: Modelling results for Reference and Baseline scenarios. This document describes the organisation of the electronic deliverables. It also summarises the debugging of SWAT+ software elements, adding PAICSWAT software and dedicated original scripts for the postprocessing of the modelling results (Chapter 4).

Document is written in English, it contains 30 pages, 16 figures, 4 tables and 4 references.



Table of contents

Summary	3
Introduction	5
1. SWAT+ debugging	6
1.1. Source code changes	6
1.2. Documentation	15
1.3. SWAT+ Editor	15
2. Structure of modelling results	16
3. PAICSWAT	20
3.1. Description of PAICSWAT	20
3.2. Illustrations of modelling results	21
4. Scripts for postprocessing of SWAT+ modelling results	26
4.1. Water and nutrient balance	26
4.2. Calibration overview	27
4.3. Overview of concentration in reaches	29
References	30



Introduction

This document is a description of the electronic deliverable: modelling results for reference and baseline scenarios. The modelling results are obtained by the calibrated and validated modelling system for the water quantity and quality in the territory of Latvia delivered as Deliverable R2 in PAIC (2022). The modelling system is developed on the basis of the data base prepared within Deliverable R1, PAIC (2020).

The documentation of the modelling and its results for reference and baseline scenarios will be presented in the Deliverable R4 – "Documentation of development, calibration, validation and results of SWAT+ modelling system".

This document describes the organization of the elements of the modelling system as complemented for the modelling itself with debugging of SWAT+ software elements (Chapter 1), adding PAICSWAT software (Chapter 3) and dedicated original scripts (Chapter 4) for the postprocessing of the modelling results.

The organisation of the electronic deliverables is described in the Chapter 2.



1.SWAT+ debugging

The errors in SWAT+ elements and respective corrections are described in this Section.

1.1. Source code changes

The list of the SWAT+ errors and discrepancy with SWAT+ documentation¹ and SWAT+ executable (rev60.5.4) and corresponding SWAT+ source code was communicated to TAMUS team and by part corrected in rev60.5.4 of 13-Apr-2022:

- Documentation about recall input file format (we use it for pointsources) is not corresponding to what is in the SWAT+ source code and SWAT+ executable. Documentation states that first two columns of data for daily, monthly and yearly timesteps are IYR and ISTEP, while in source code there are 6 columns (jday, mo, day_mo, iyr, ob_typ, ob_name) (see recall read.f90, line 150).
- Point source (recall) data format written by SWAT+ editor does not correspond to point source format used by SWAT+. It seems that it corresponds to the SWAT+ documentation, but not to source code.
- If pointsource is monthly, then only first calendar month of each year (January) is applied (read from file). There should be line `istep = istep + 1` in monthly case (line 180 of recall_read.f90), otherwise all 12 monthly data will be assigned for January only (istep=1).
- It seems that if weather files start with the later date than simulation start, weather generator is used for the whole simulation period, instead of only for the missing dates.
- Seems that weather generator dewpoint values cannot anymore be entered as relative humidity, despite that documentation states "If all twelve months are less than one, the model assumes relative humidity is input". Comparing SWAT and SWAT+ source codes, it seems that relevant code part is missing in SWAT+.
- For some of the input files documentation states that "The title line is not processed by the model and may be left blank". It seems that it is not anymore the case (comparing to SWAT2012). For instance, if header or title lines in atmo.cli is blank then atmo.cli is read incorrectly. This is related to change to how the header lines are read from files. SWAT2012 used formatted input (e.g. `read (127,5101) titldum`, where line 5101 contained format) – that allowed to read the blank line also. Now in SWAT+ wildcard format is used (e.g. `read (127,*,iostat=eof) titldum`) – in that case if the line is blank

¹ inputs_swatplus_rev60_5.docx from <https://swatplus.gitbook.io/docs/user/io>

reading continues to the next non-blank line and real data could be misinterpreted as the comment.

- Names of files in pcp.cli, tmp.cli etc. and station names in atmospheric deposition atmodep.cli must be in sorted order (not stated anywhere in documentation!), otherwise potentially incorrect data will be assigned (searching procedure in search.f90 works only on sorted arrays (!), but on reading no sorting is performed!).
- If atmo.cli file is provided then even for the weather stations without assigned atmospheric deposition (set to null), some undocumented default values for wet deposition is used (c_NH4=1 mg/l, c_NO3=0.2 mg/l) (see type atmospheric_deposition of climate_module.f90).

The further changes listed below are made by PAIC in the reference SWAT+ source code and presented in github repository²:

1. Wrong output parameter unit. File *ru_day.txt*
<https://github.com/andrejstmh/SWATplus/commit/4d8678d81f01cd79a1a7c74fb90c691ae9aa8b8d#diff-8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372>
2. Wrong output parameter unit. File *channel_sd_day.txt*
<https://github.com/andrejstmh/SWATplus/commit/cdf7c476bc205e16098fde4f558aa3983192a5f3#diff-8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372>
3. Output variables increase all the time (time step). File *soil_nutcarb_out.txt*
<https://github.com/andrejstmh/SWATplus/commit/c6c259351c2577d5bdbfee3fa5bc98b0f491ef57#diff-8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372>
4. Incorrectly calculated channel depth. That leads to very small channel outflow. File *ch_watqual4.f90* (rchdep = 0.)
<https://github.com/andrejstmh/SWATplus/commit/764564e06fb0d8786a42789cdbc5b5d9fe7fddcc#diff-8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372>
5. Variables are misnamed (exchanged surq_gen <=> surq_cont)
<https://github.com/andrejstmh/SWATplus/commit/7a9c54eed89cbbfd5f8667ac8916ac4877320060#diff-8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372>

² https://github.com/andrejstmh/SWATplus/commits/main/source_codes

6. Surfaces runoff processes are not running in regular case. Mismatched (exchanged) wetland and regular cases.
<https://github.com/andrejstmh/SWATplus/commit/d8a575c61e33339f1dd82c1e379880ac08c80ec9#diff-8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372>
7. Coefficients cn3_swf and perco are initialised to constant values if the tile drain is on
<https://github.com/andrejstmh/SWATplus/commit/d588df775f37fa55a06da175d5b30d32ee0b07eb#diff-8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372>
8. Evapotranspiration model differs from SWAT (coefficient esco) comparing SWAT theoretical documentation TWRI(2011)³ Eq 2:2.3.17 with the source code rev.627/etact.f
<https://github.com/andrejstmh/SWATplus/commit/e652f28d1665f32b06b63d664c681d4bf8af4fb3#diff-8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372>
9. Evapotranspiration model differs from SWAT. According to TWRI(2011) Eq 2:2.3.18 & source code rev.627/etact.f
<https://github.com/andrejstmh/SWATplus/commit/f5401037f379defa733ee107d249287d37e6d19c#diff-8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372>
10. Aquifer nitrogen flow units and initial content are incorrect
<https://github.com/andrejstmh/SWATplus/commit/21af44f544e4af0d2aca112228dff29d91dfd7#diff-8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372>
11. The computation of the travel time of the lateral flow is switched on
<https://github.com/andrejstmh/SWATplus/commit/cf5e71eb7a838335e35b01e74ff4e57771a5e18d#diff-8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372>
12. The computation of the lateral subsurface flow from first soil layer is switched on
<https://github.com/andrejstmh/SWATplus/commit/d2563c25ab187bbf6e77b9c3db435807d6e47a#diff-8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372>
13. Redundant line with wrong code is removed
<https://github.com/andrejstmh/SWATplus/commit/cd931978ea8544d064d399069a888>

³ <https://swat.tamu.edu/media/99192/swat2009-theory.pdf>

[6cc7e407440#diff-](#)

[8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372](#)

14. *dep_imp* variable is introduced (like in SWAT v.627).

<https://github.com/andrejstmh/SWATplus/commit/a494ec84e131c3c85d16a3213bf1aef893904e34#diff->

[8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372](#)

15. A possibility to control output of files *wetland_*.txt* separately from the *reservoir_*.txt* is introduced:

<https://github.com/andrejstmh/SWATplus/commit/df27eed54c9bdf82bac88884c7f287df7c8c054#diff->

[8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372](#)

16. Module *wetland_control.f90* containing uninitialized variables is bypassed

<https://github.com/andrejstmh/SWATplus/commit/2a60358aceb00005a0a046c0a4961c7538539485#diff->

[8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372](#)

17. Delay time for aquifers is introduced (like in SWAT v.627)

<https://github.com/andrejstmh/SWATplus/commit/de55a5e7d4e1e92aac5196ec3fceed92eeee6ee0#diff->

[8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372](#)

18. Evaporation module is changed (like in SWAT v.627)

<https://github.com/andrejstmh/SWATplus/commit/3b1e2bff30186326fbad9f98c7c156317f5b10ec#diff->

[8fed27fc4815cb7df6aa86a3e660491b04b6323635c44972575ab52fc6a6d372](#)

19. Error in search procedure. Uninitialized variable "aqu_found"

<https://github.com/andrejstmh/SWATplus/commit/a64dd6aeca7c620caeac491b6499d1655ddab248>

20. Corrected snow pack temperature

<https://github.com/andrejstmh/SWATplus/commit/6e248856a34351c3450bcee6bfc38b63f67beadf>

21. Corrected output_losses_header1

<https://github.com/andrejstmh/SWATplus/commit/062b171a219a900fc538abefe07933097e0c15fa>

22. Half-life of NO₃ implemented
<https://github.com/andrejstmh/SWATplus/commit/8bea0a7c778f50df9824567e3126cbad79f77893>
23. Update fresh residue plants and soil layers procedure
<https://github.com/andrejstmh/SWATplus/commit/88c0b3b45718724fd75a92f45758715cd8b12ff9>
24. Fresh residue tillage mixing added
<https://github.com/andrejstmh/SWATplus/commit/028e92666381df3e130dc4454b95aed7c95e4c04>
25. Mgt_kill operation: add above ground mass to soil fresh residue pool
<https://github.com/andrejstmh/SWATplus/commit/8b8756b00ef574b38d25491643a81cb14fdcce1f>
<https://github.com/andrejstmh/SWATplus/commit/b177c68d0b0f7f88dc84cbf9f08bb12cb4a9193e>
26. Fresh organic nutrients remineralization added
<https://github.com/andrejstmh/SWATplus/commit/7cdf7b30a0a1dfcae45b9881db9768e11db91b7b>
<https://github.com/andrejstmh/SWATplus/commit/831cfd17466b27b95510ac2f60aaca42ed8049a3>
27. Bugfix for tillage of fresh organic residue mixing
<https://github.com/andrejstmh/SWATplus/commit/6d89f0b887f043c8638a765009feb363c6e6e8d5>
28. Corrected soil nitrate denitrification formula
<https://github.com/andrejstmh/SWATplus/commit/d367c61606555e4cfcf4093410caef757b6c3514>
29. Organic fertilizer added to fresh organic pool (as in SWAT2012)
<https://github.com/andrejstmh/SWATplus/commit/6db7844c520e7c8036d7ead14fef4ead57fc0823>
30. Harvest index routines modified to correspond to SWAT2012
<https://github.com/andrejstmh/SWATplus/commit/f873c980f69592fd8b0e44ca74384c15d2fad922>
31. Corrected sediment yield calculations
<https://github.com/andrejstmh/SWATplus/commit/64060bd9ea1ae88d04ca839b77918300514261e5>



32. Distribution of plant nitrogen uptake like in SWAT2012
<https://github.com/andrejstmh/SWATplus/commit/324111ac9ca82dac768ae193bc644b85495148fd>
33. Leaf area index growth dependent on total stress like in SWAT2012
<https://github.com/andrejstmh/SWATplus/commit/d1d4d26ec4b1ef4746589d5334cad599088f4>
34. Possibility to input initial organic phosphorus and nitrogen in soil layer implemented (as in SWAT2012)
<https://github.com/andrejstmh/SWATplus/commit/85bfa1e220cff7a72f7013849f862a08801b2632>
35. Changed routine for the biomass growth (bioday)
<https://github.com/andrejstmh/SWATplus/commit/b46ea5b770bdd8a764222a3ccbc3756048d365da>
36. Implemented soil carbon initialization from file
<https://github.com/andrejstmh/SWATplus/commit/9697215d7508345b67086a18f9e3fef4db87885>
37. Restored possibility to select soil phosphorus model as in SWAT2012
<https://github.com/andrejstmh/SWATplus/commit/62d51dcec2b3fc2ab7a2d6e72bfba466a324e7ab>
38. Fixed units of conversion factor in soil initialization routine
<https://github.com/andrejstmh/SWATplus/commit/c737f7f73be45517ead1e0d28b5e1a28e403fa9b>
39. Implemented soil organic phosphorus initialization
<https://github.com/andrejstmh/SWATplus/commit/dfe657d3117125b6d448b6d1175994514a1e1ddc>
40. Implemented phosphorus in surface runoff from residue (as in SWAT2012)
<https://github.com/andrejstmh/SWATplus/commit/566d7567fbff3a2ec9153c9af9ef756833a3c9c>
41. Implemented initialization of soil mineral phosphorus un nitrogen from input file (as in SWAT2012)
<https://github.com/andrejstmh/SWATplus/commit/9aa67e5e64418c19406620b306cfb72da8b260e7>



- <https://github.com/andrejstmh/SWATplus/commit/7f26f6e9dda43bc5a71f00a497f2ec2ddce99cac>
42. Formula for initial stable soil phosphorus pool should use carbon concentration not mass per area
<https://github.com/andrejstmh/SWATplus/commit/f78905f02d36c9a33b999268b5ea5d53441dfca6>
43. Code merged with swat official version 13apr2022-Version-60_5_4
<https://github.com/andrejstmh/SWATplus/commit/f9cfe948775a3871dd9fa983f726e0fc2155f193>
44. Simulation.out header fixed
<https://github.com/andrejstmh/SWATplus/commit/2b9f3b90b40b1071f1f958e9a7e9384edfa6941b>
45. PSP calculation should contain carbon concentration not mass per area
<https://github.com/andrejstmh/SWATplus/commit/733b0a2d1ee1b3f896eec74a57cfb6bc47c3fcb7>
46. Fix for concentration based mixing in tillage
<https://github.com/andrejstmh/SWATplus/commit/e44dba159f2ddf06ac4cc010f1223815a51dbeb5>
47. Coefficient nperco_lchtile enabled
<https://github.com/andrejstmh/SWATplus/commit/9764d1842217c09a356772fb682cee966ece3a3d>
48. Revert soil soluble phosphorus routines to previous SWAT+ version to correspond to SWAT2012
<https://github.com/andrejstmh/SWATplus/commit/8b8ff55f980a5d4a400140485af3976aa1362d06>
49. Bugfix in phosphorus sediment calculations
<https://github.com/andrejstmh/SWATplus/commit/dbb3cae2fa72168cadd0484c8950545adbbf5a6a>
50. Change LAI, add leaf mass to residue at dormancy for perennials
<https://github.com/andrejstmh/SWATplus/commit/4ca991f5dfc015956c6430034b014eb7d48439fa>

51. Implemented leaf senescence dependent on plant accumulated phu
<https://github.com/andrejstmh/SWATplus/commit/5f6a9fd085b46f18be99ad096384b34678636501>
52. Leaf drop should affect total plant mass – fixed
<https://github.com/andrejstmh/SWATplus/commit/765d01eefb290e78a0b94446688517ae3a5e660e>
53. Dormancy for forests implemented
<https://github.com/andrejstmh/SWATplus/commit/96e5dfc9801168ac95bb636b54959da55cbd178c>
54. Plant partitioning for forests, additional parameter included
<https://github.com/andrejstmh/SWATplus/commit/dc4874937d95365ea11d9d606feeb97f4662135e>
55. End of year management operations added (as in SWAT2012)
<https://github.com/andrejstmh/SWATplus/commit/27ec883b0edbdd338688dd54822734b7d907fce9>
56. Bugfix - no need to reset phuacc after harvest only operation
<https://github.com/andrejstmh/SWATplus/commit/c73dfce5eeb51ac9ecd64f711c3bb7e7eee4eb7b>
57. Nutrients outflowing from reservoir was not subtracted from reservoir - fixed
<https://github.com/andrejstmh/SWATplus/commit/7ea920867cd7ea185434dc97ec22aa9de074896a>
58. Implemented reservoir settling rate as in SWAT2012
<https://github.com/andrejstmh/SWATplus/commit/448473165355da58056918f2ac84ddb01aac2cbd>
59. Reservoir initialization bug fix
<https://github.com/andrejstmh/SWATplus/commit/5f0d0c9f02cb3f8c73f094072618395a2bed32f3>
60. Minimal flow from reservoir implemented
<https://github.com/andrejstmh/SWATplus/commit/730267719a4c2d464ea5c98e269dda3530b352f9>
<https://github.com/andrejstmh/SWATplus/commit/25d93a1eba9754bbd09c10b1b820945dc4f99514>

61. Repaired aquifer flow numerical scheme that was unstable (implemented order of operations as was in SWAT2012)
<https://github.com/andrejstmh/SWATplus/commit/dfde88519d5ede6951e3e01e90df43509d48a759>
62. Corrected uninitialized output variable in action "hru_fr_update"
<https://github.com/andrejstmh/SWATplus/commit/9b75dd914539b7d061946399e25cf8acce87c02>
63. Aquifer nutrient output dimension changed to kg/ha
<https://github.com/andrejstmh/SWATplus/commit/bec9744e2e3639b3a17323f5711443e0c30fcee5>
64. Corrected routing unit output
<https://github.com/andrejstmh/SWATplus/commit/1a248048fd12b4ec0c298cd99509406ea0413ea8>
65. Routing unit number added to routing unit output
<https://github.com/andrejstmh/SWATplus/commit/339081f285aa0f7148cbd6019b2467ce2a482a3f>
66. Enabled and bugfix source code for grassed waterways
<https://github.com/andrejstmh/SWATplus/commit/3fc00b39edc86951e374f07a24fa02ab50cdbaa1>
67. Bugfix_in plant determination at scheduled management operation
<https://github.com/andrejstmh/SWATplus/commit/db46d5a3ec55775db9221dbab37fac9c6dd8f0a0>



1.2. Documentation

The subsection documents errors and missing parts of SWAT input/output documentation⁴. This may include differences between the documentation and the source code. The following problems are found in *outputs_swatplus_rev60_5.pdf*:

- Document does not contain information about *soil_nutcarb_out*.
- Wrong documented unit. File *aquifer_day.txt* (**seepno3** kg(N)/ha)
- Wrong documented unit. File *channel_sd_day.txt* (**flo_in, flo_out** m3/s)
- Document does not contain information about input parameter **latq_co**
- Inconsistent parameter name **cn3_swf**

1.3. SWAT+ Editor

Errors found in SWAT+ Editor

1. There is no way to setup daily and monthly output for *soil_nutcarb_out.txt* in the SWATPlus editor.
2. Wrong name of the parameter. HRU/ Hydrology Properties /Plant ET curve number coefficient.
3. There is no possibility to create setup without routing units.
4. Error in the save procedure (file *ari/fileio/hru.py* class *Hru_data_hru* variable *field*).

⁴ <https://swatplus.gitbook.io/docs/user/io>

2. Structure of modelling results

The water quality modelling system is electronically delivered containing all the data and scripts needed for the creation of the system, see PAIC (2022). The water quality modelling system is generated automatically by executing the script *main.py* from folder *projectDir/Scripts*.

Main.py also runs the calculation of the modelling system and creates the results in *setupsDir* folders. The folder structure where the results can be found would be as follows:

LVSWAT2022(projectDir)/Watersheds(setupsDir)/RiverBasinName/WatershedName/ScenarioName/

For, example, the result folder for baseline scenario in Venta river basin district and the 1st Abava watershed will be

LVSWAT2022/Watersheds/Venta/Abava_1/Baseline/

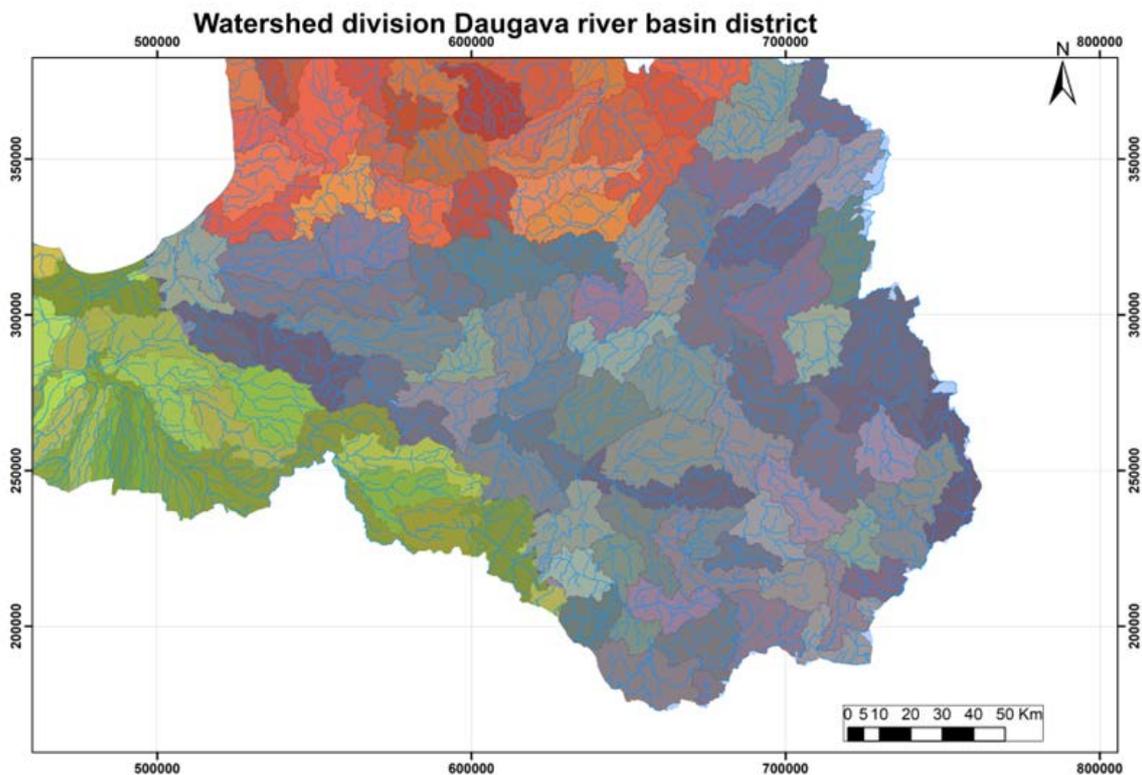


Figure 1: Watersheds of Daugava RBD.

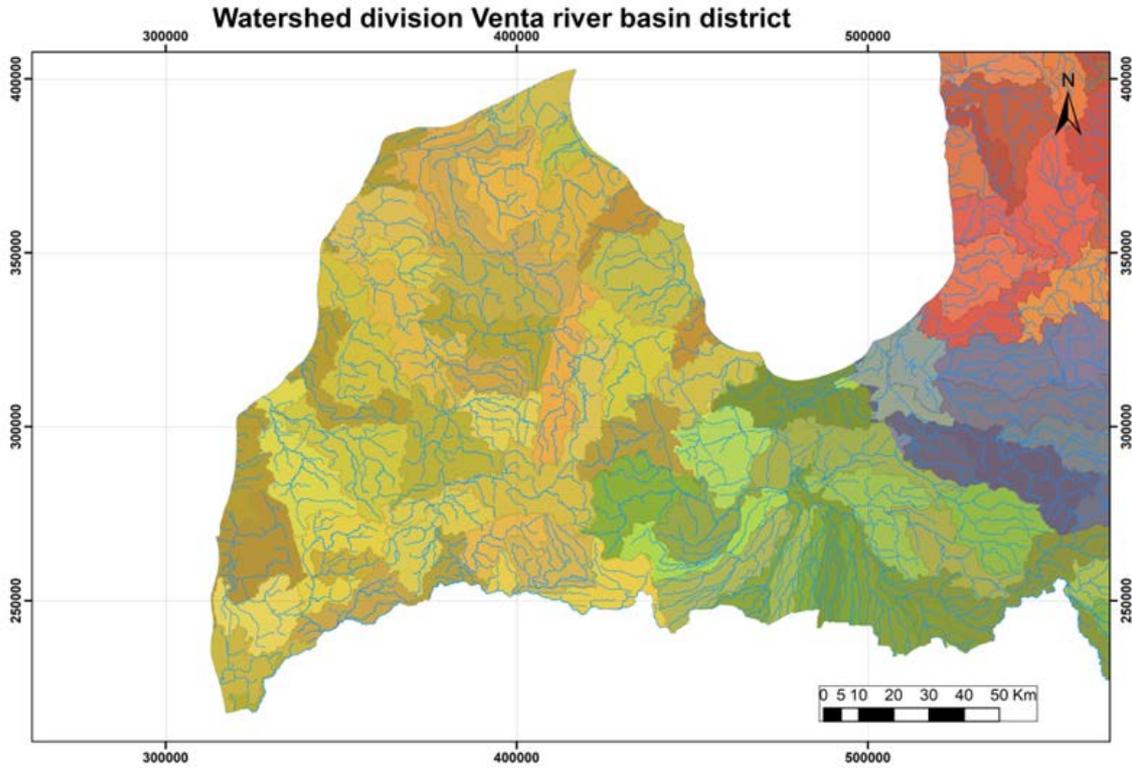


Figure 2: Watersheds of Venta RBD.

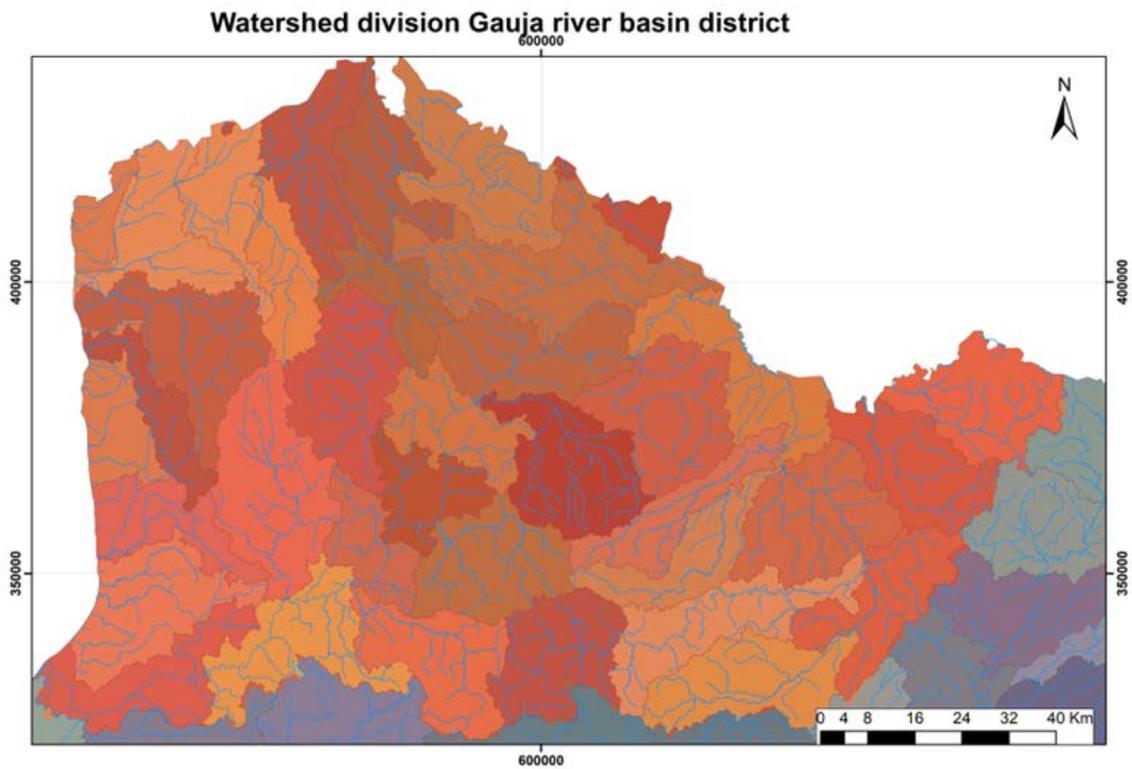


Figure 3: Watersheds of Gauja RBD.

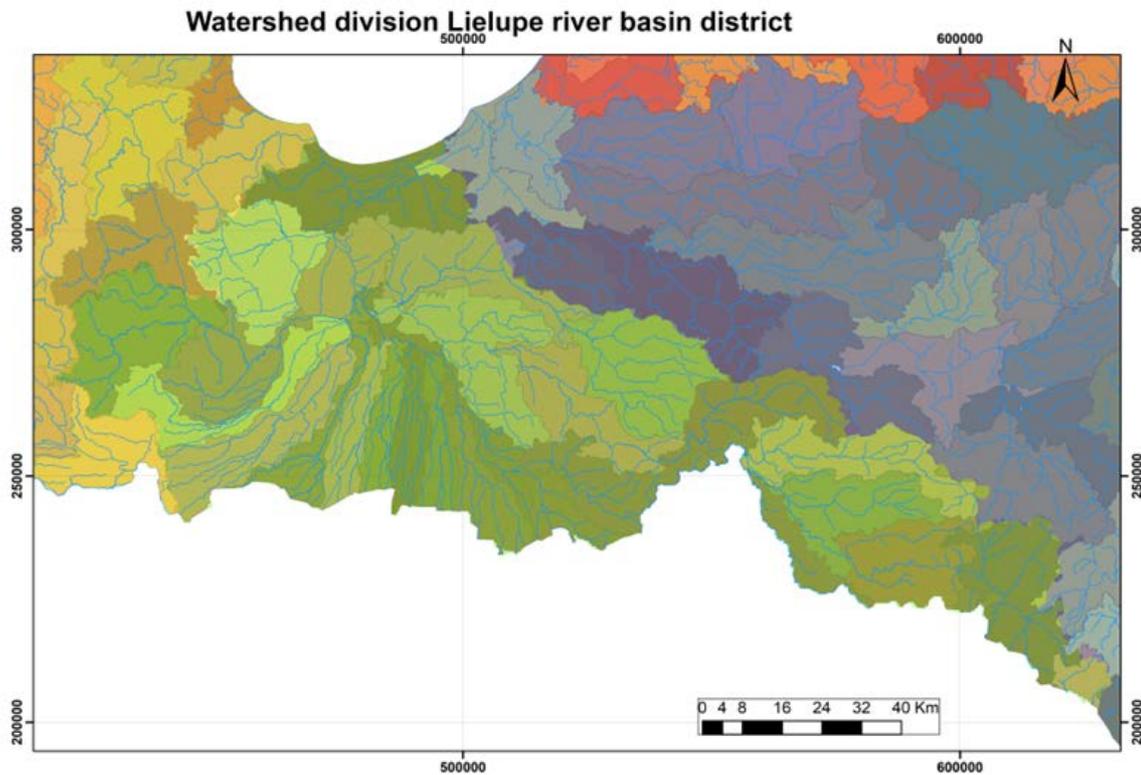


Figure 4: Watersheds of Lielupe RBD.

The SWAT+ model for the specific watershed can be found in this folder, together with its input and output files for the specific scenario. The output file extension .out indicates these files while most of the other files are text input files for the model to run.

Time period of calculation is 2006-2018 (with 2006 and 2007 used for warm-up of the model).

The model system for Latvia consists of 4 river basin districts which are divided into 169 watersheds, which are further divided into 3469 catchments. The division of 4 RBDs into watersheds is shown in Figures 1-4, while illustrations of further division of two watersheds in the catchments – in Figures 5-6.

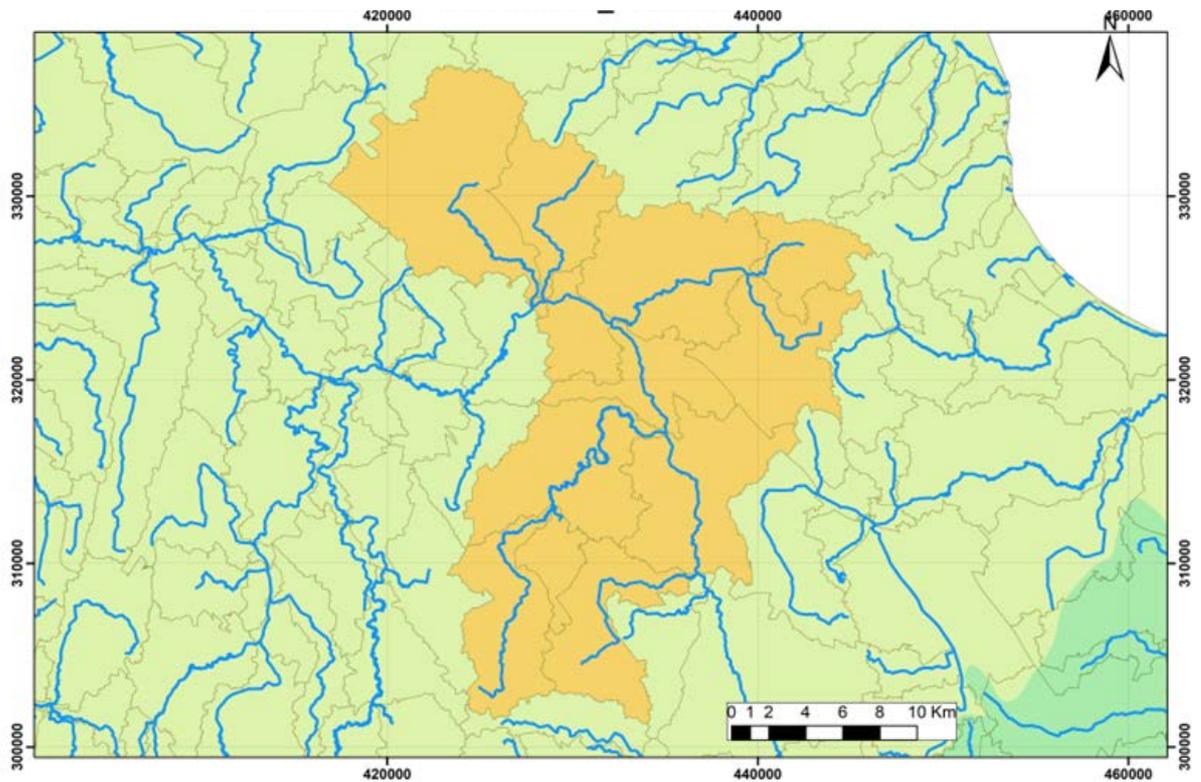


Figure 5: Catchments of Abava watershed.

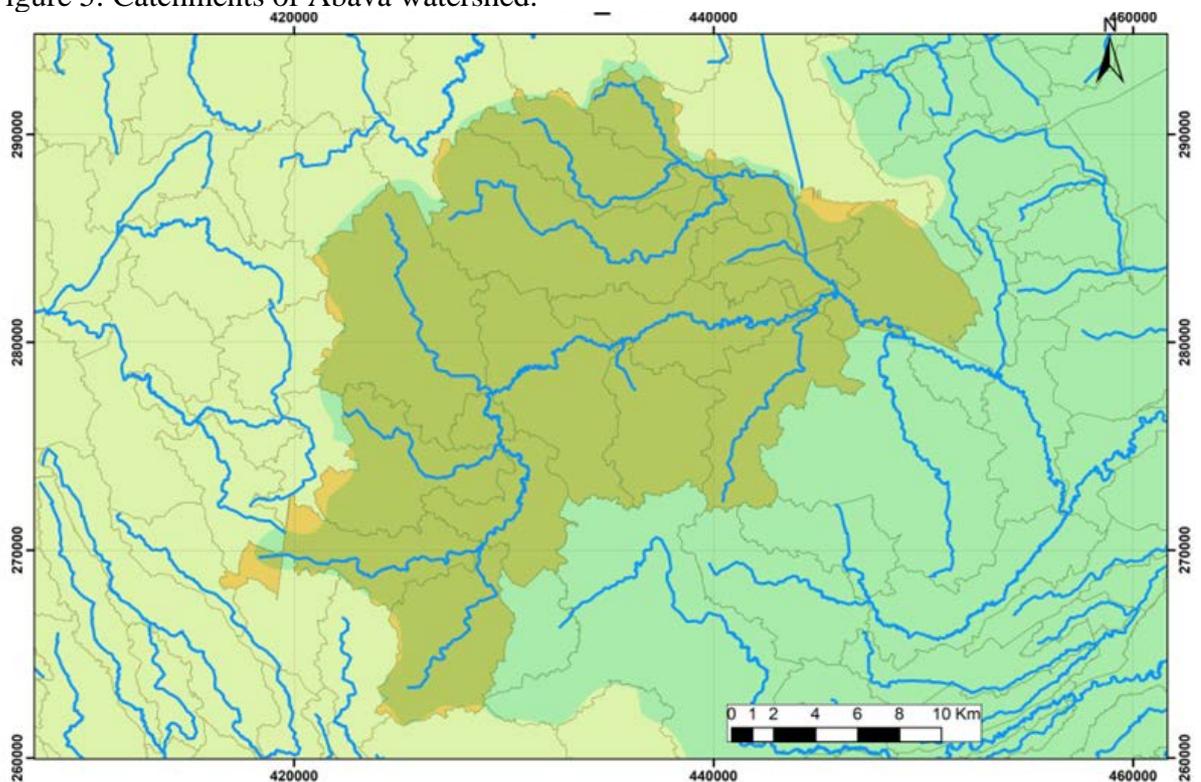


Figure 6: Catchments of Bērze watershed.

3. PAICSWAT

3.1. Description of PAICSWAT

The main tool for visualization and analysis of SWAT+ modeling results is PAICSWAT software by PAIC. Output results of modelling system could be visualized by PAICSWAT software (PAIC, 2012). The visualization works by converting output results calculated by SWAT+ to the SWAT2012 format which could be loaded into PAICSWAT. The preparation step includes running of *prepare_PAICSWAT.py* script located in the main scripts directory. It includes two subroutines:

1. *initPAICSWAT.prepare_PAICSWAT* – this subroutine regenerates *PAICSWAT.ini* configuration file. The location of the configuration file is provided in *settings.PAICSWAT_ROOT* configuration variable (by default it is located in *PAICSWAT* subdirectory of the project directory). In case if the *overwriteData* parameter is *True* then the catchment and river geometrical data in format required by PAICSWAT will be regenerated as well. Default location of PAICSWAT geometrical data is under subdirectory *PAICSWAT/Data* of the project directory.

```
initPAICSWAT.prepare_PAICSWAT(project,  
PAICSWATDataDir=settings.PAICSWAT_ROOT + "/Data",  
PAICSWAT_INI =settings.PAICSWAT_INI,  
overwriteData = True)
```

2. *initPAICSWAT.FillPaicSWAT_forsetups* – this subroutine generates two SWAT2012 files that are required for the PAICSWAT to function – *file_paicswat.cio* and *fig.fig*. These files are copied to each of the watershed setups.

```
initPAICSWAT.FillPaicSWAT_forsetups(project, watershednamelist =  
watershednamelist)
```

Each change of the calculation directory (provided in *settings.defaultCaseSubDir*) requires re-running of the preparation procedure.

Conversion of SWAT+ results to SWAT2012 format are achieved by the subroutine *convertOutputToPAICSWAT.convertAll*. The calculation loop in the main script (*main.py*) automatically call this conversion subroutine after calculation of each of the watersheds (for that particular watershed). It also can be called explicitly by script *convert_results.py* where selection of watersheds for conversion are implemented.

The executable of PAICSWAT software (*PAICSWAT.exe*) is provided under *PAICSWAT* subdirectory of the project directory. It could be run from there after preparation steps are accomplished. The *PAICSWAT* functionality compatible between SWAT+ and SWAT2012 includes:

- Selection of catchment and watershed on the PAICSWAT map.
- Loading of calculation setups (Menu *File->Load case*).
- Loading of compatible setup results (Menu *File->Load from directory*).
- Loading and visualisation of time-graphs in PAICSWAT data format (*.sta+.txt*) (e.g. observation and weather data).
- Visualisation of time-graphs of outputs at catchment level (reaches, subbasins, reservoirs (PAICSWAT tabs *Model->RCH*, *Model->SUB*, *Model->RES*)).
- Comparison of model outputs to observations (PAICSWAT tab *Calibration*).
- Statistics and correlations of abovementioned graphs (Buttons *Statistics*, *Correlations*, *Concentrations*, *Monthly value*, *Normalize by area*).

Incompatible functionality includes, but are not limited to:

- Visualisation of results on a HRU level.
- Displaying and modification of SWAT2012 input files.
- Running of calculation by SWAT2012.

3.2. Illustrations of modelling results

The modelling results for Bērze_2 watershed are illustrated with the use of PAICSWAT.

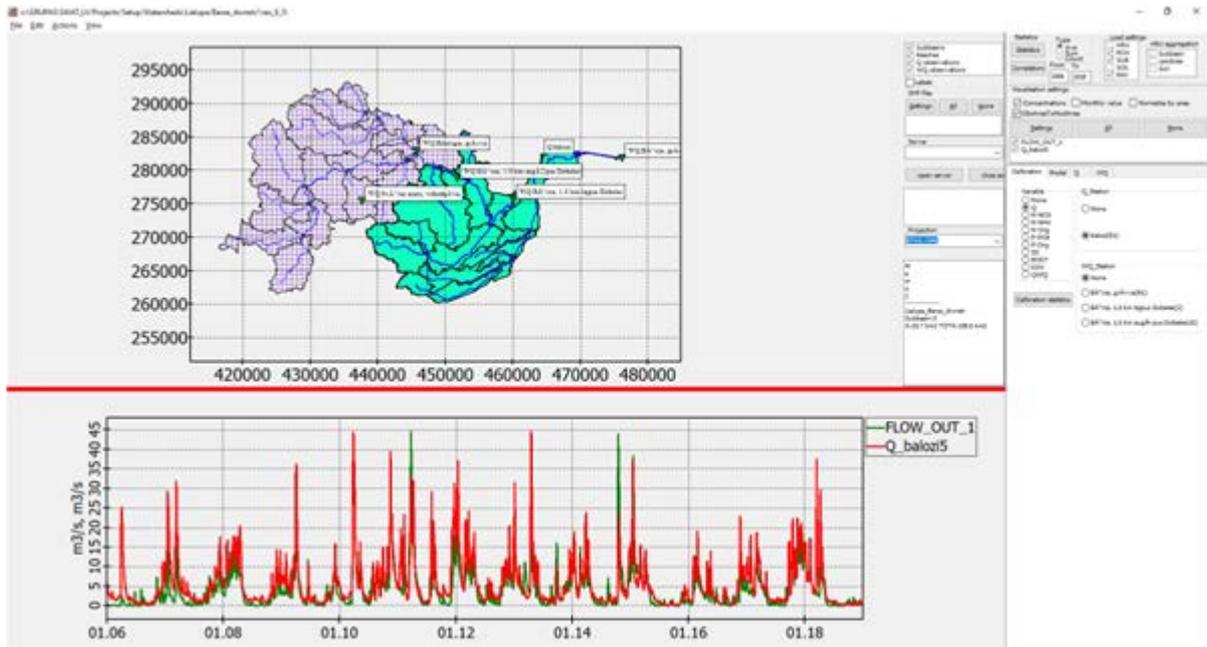


Figure 7: Screenshot of PAICSWAT.

The screenshot of PAICSWAT is shown in Figure 7. It consists of map of the watershed under consideration (Figure 8), time graph of observed and modelled parameters (Figure 9) and several control panels. Parameter selection and calibration panel is shown in Figure 10.

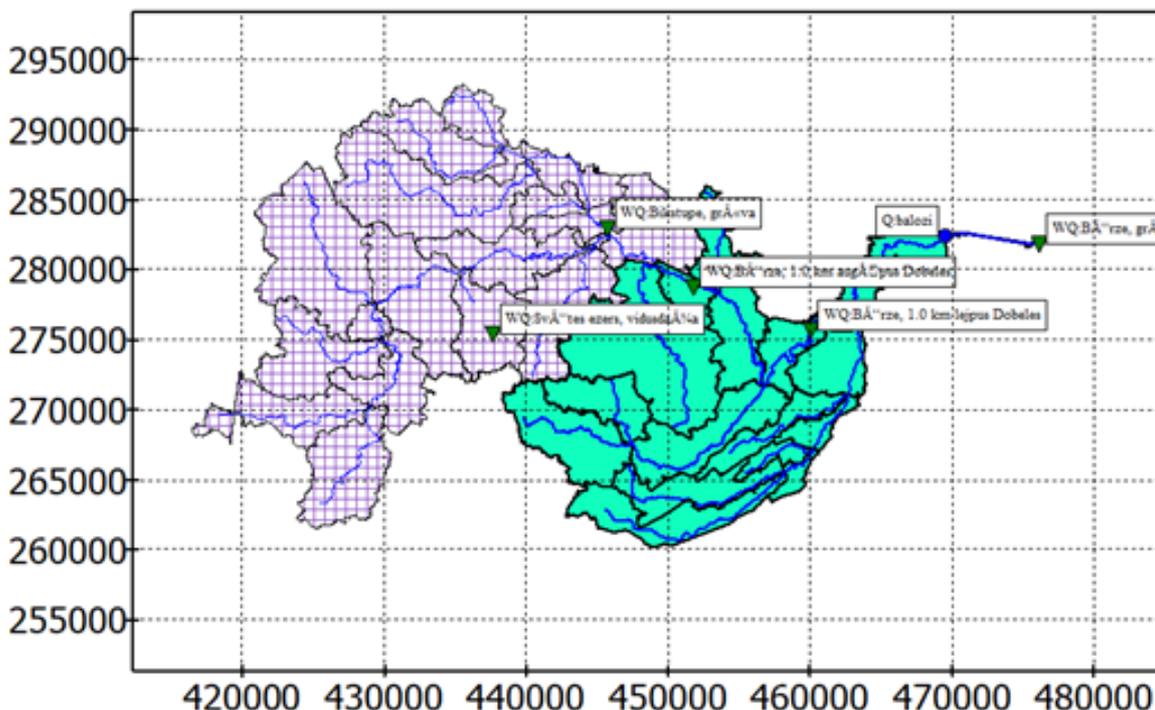


Figure 8: PAICSWAT: map of Bērze_2 watershed.

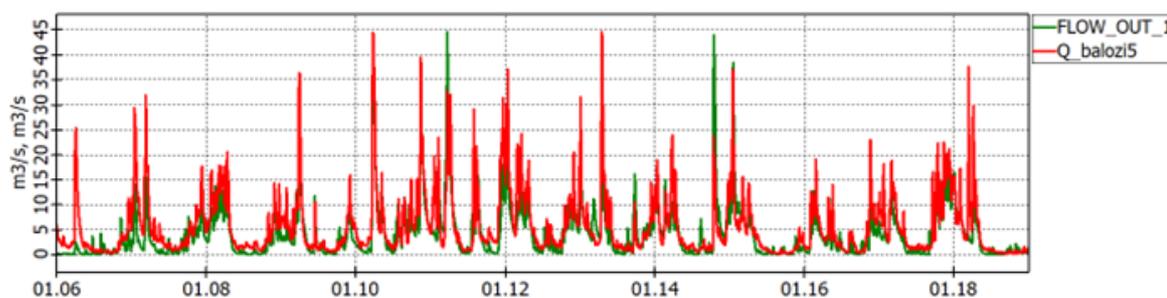


Figure 9: Observed vs modelled Discharge at Bērze-Baloži station.

PAICSWAT allows statistical analysis (Figure 12) and calculation of seasonal parameter values (Figure 11) from observed and modeled data series in selected (figure 13) stations.

Visualisation of water quality parameters and water balance components by PAICSWAT are given in Figures 14-16.

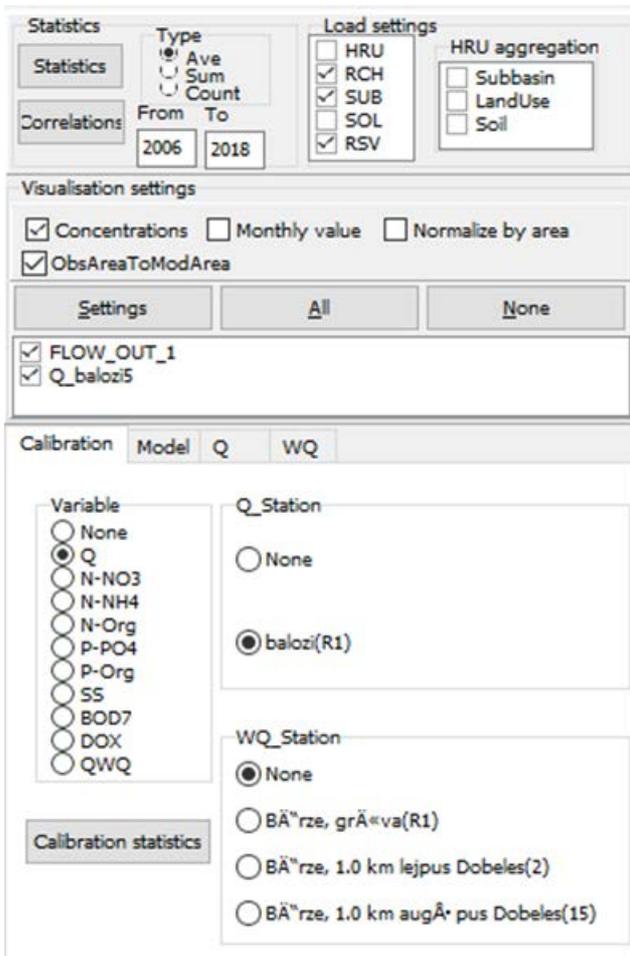


Figure 10: PAICSWAT selection and calibration panel.

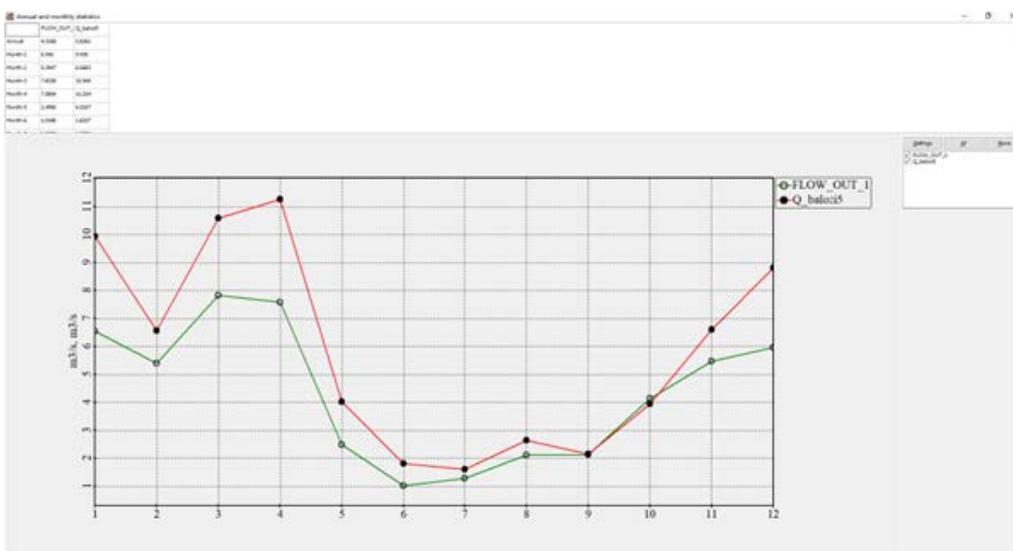


Figure 11: Yearly overview of discharge Bērze(Baloži) station Observed vs modeled, PAICSWAT statistics visualization screen.

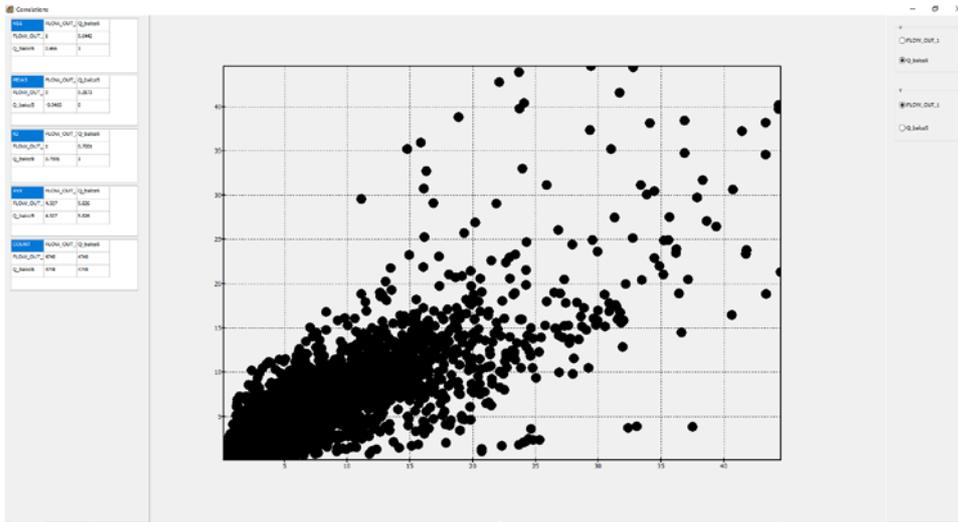


Figure 12: Statistical overview of discharge at Bērze (Baloži) station Observed vs modeled. PAICSWAT Correlations visualization screen.

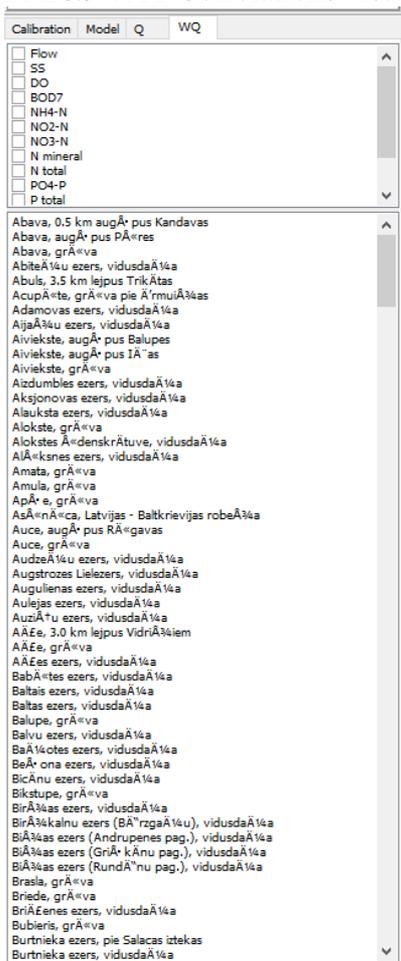


Figure 13: Water quality station selector in PAICSWAT for visualizing specific stations and parameter.

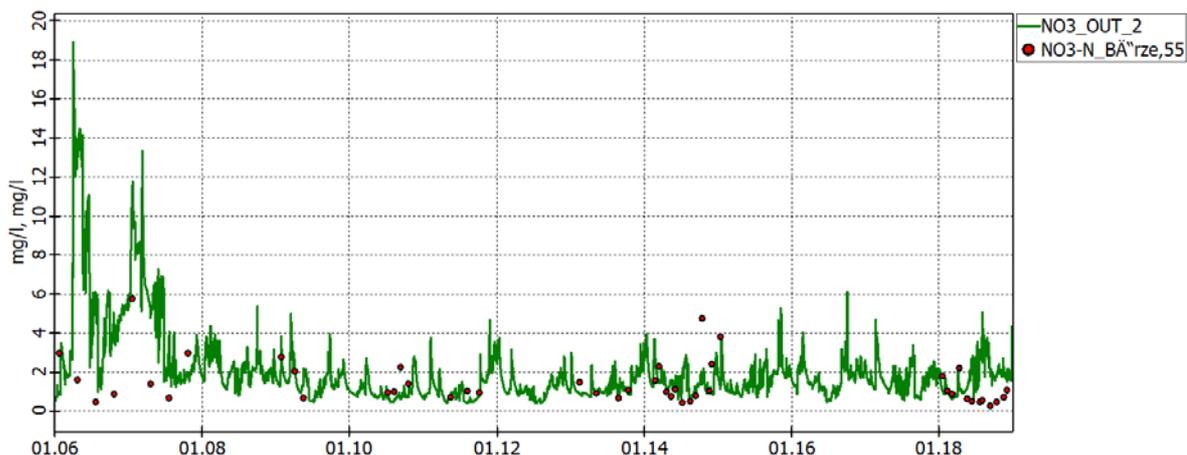


Figure 14: Mineral nitrogen (N-NO3) concentrations at measurement station Bērze- Bērze, 1km downstream Dobe.

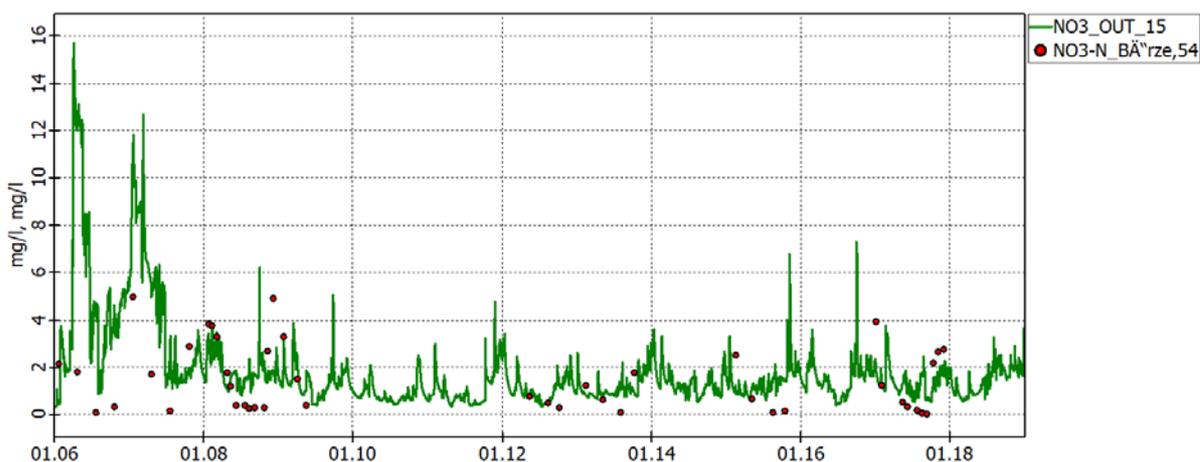


Figure 15: Mineral nitrogen (N-NO3) concentrations at measurement station Bērze- Bērze, 1km upstream Dobe.

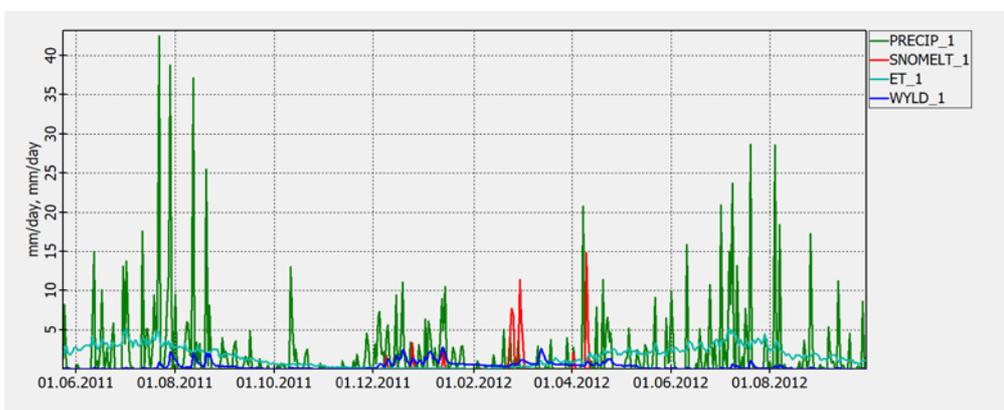


Figure 16: Time-series of water balance components in a Bērze_2 watershed.

4. Scripts for postprocessing of SWAT+ modelling results

4.1. Water and nutrient balance

Script for the calculation of a catchment level water balance from the calculation results is *waterbalance.py*. It is located in the main script directory of the project. It allows to calculate water-balance results in all LV territory or for selected watersheds (configurable by variable *watershednamelist*). The results are placed in the text file, the name of this file is user configurable (variable *wb_TXTFILE*). Internally, script uses subroutine *result_processing.prepare_sub_waterbalance* that returns the water balance table in *pandas Dataframe* format. The fields of the table are summarised in the Table 1.

Table 1. Fields of the water balance table.

Field	Description
catchmentid	Global Id of the catchment (corresponds to field <i>catchmentid</i> of <i>catchments</i> table under schema <i>catchments</i> of PostgreSQL database of project)
CNMIN_a	Average concentration of mineral nitrogen in a yield from the catchment, mg/l
CNTOT_a	Average concentration of total nitrogen in a yield from the catchment, mg/l
CPMIN_a	Average concentration of mineral phosphorus in a yield from the catchment, mg/l
CPTOT_a	Average concentration of total phosphorus in yield from the catchment, mg/l
ET	Average evapotranspiration in catchment, mm/day
GWNO3	Groundwater mineral nitrogen yield from the catchment, kg/ha/day
GW_Q	Groundwater water yield from the catchment, mm/day
LATNO3	Lateral mineral nitrogen yield from the catchment, kg/ha/day
LATQ	Lateral water yield from the catchment, mm/day
NMIN_YLD	Average mineral nitrogen yield from the catchment, kg/ha/day
NSURQ	Surface flow mineral nitrogen yield from the catchment, kg/ha/day
ORGN	Organic nitrogen yield from the catchment, kg/ha/day
ORGP	Organic phosphorus yield from the catchment, kg/ha/day
PET	Average potential evapotranspiration in catchment, mm/day
PMIN_YLD	Average mineral phosphorus yield from the catchment, kg/ha/day
PREC	Average precipitation in catchment, mm/day
SEDP	Average phosphorus yield in sediment from the catchment, kg/ha/day
SNOMELT	Average snowmelt in catchment, mm/day
SOLP	Average soluble phosphorus yield from the catchment, kg/ha/day
SURQ	Average surface water flow in catchment, mm/day
TILEQ	Average tile drain water flow in catchment, mm/day

Field	Description
TNO3	Average mineral nitrogen yield in tile flow from the catchment, kg/ha/day
TOTN_YLD	Toal nitrogen yield from the catchment, kg/ha/day
TOTP_YLD	Total phosphorus yield from the catchment, kg/ha/day
WYLD	Average water yield from the catchment, mm/day

Resulting text file can be loaded into QGIS as a text layer, joined on the field *catchmentid* to the *catchments* table under schema *catchments* of PostgreSQL database of the project to allow for visualisation of catchment water and nutrient balance components.

4.2. Calibration overview

Script for the calibration overview is *calibration_overview.py* located in the main *scripts* directory of the project. By running it the Nash-Sutcliffe efficiency coefficients and bias of model results in comparison to observation results can be calculated. The resulting text file is determined by variable *statistics_TXTFILE* while list of watersheds could be provided in *watershednamelist* (in case of empty list – for the whole territory). Script allows for automatic distinction between calibration and validation periods based on the variable *calibration_fraction* (0.0 – all period is validation, 1.0 – all period is calibration).

Before running the *calibration_overview* script, the script *prepare_calibration* should be run at least once. This script prepares connections between observation stations and model catchments in tables *qobs_catchments* and *wqobs_catchments* of scheme *calibration* of PostgreSQL database of the project (see configuration in *settings.Q_STATIONS_CATCHMENTS* and *settings.WQ_STATIONS_CATCHMENTS*).

The resulting text file contains columns listed in Table 2; possible “substances” are listed in Table 3.

Table 2: Columns of the calibration assessment text file.

Column	Description
STAID	Observation station identifier, corresponds to identifier in station files of observed data <i>Qobs.sta</i> and <i>Wqobs.sta</i>
STANAME	Observation station name
SUBSTANCE	Abbreviation of substance (see Table below)
CATCHMENTID	Global Id of the catchment (corresponds to field <i>catchmentid</i> of <i>catchments</i> table under schema <i>catchments</i> of PostgreSQL database of project)
IS_RESERVOIR	<i>True</i> if the output of catchment is from reservoir, <i>False</i> otherwise
DATACOUNT	Count of data points for comparison
CALIB_PER	Time period of calibration

Column	Description
CALIB_COUNT	Count of data points for comparison in calibration period
CALIB_NSE	NSE of daily time series for a substance in calibration period
CALIB_PBIAS	PBIAS of daily time series for a substance in calibration period
VALID_PER	Time period of validation
VALID_COUNT	Count of data points for comparison in validation period
VALID_NSE	NSE of daily time series for a substance in validation period
VALID_PBIAS	PBIAS of daily time series for a substance in validation period
CALIB_MON_PBIAS	PBIAS of monthly averages for a substance in calibration period
CALIB_MONR2	R2 of monthly averages for a substance in calibration period
VALID_MON_PBIAS	PBIAS of monthly averages for a substance in validation period
VALID_MONR2	R2 of monthly averages for a substance in validation period

Table 3: List of available parameters (“substances”) in the calibration assessment text file.

Parameter	Description
Q	Discharge, m ³ /s
NO3	Concentration of nitrate nitrogen, mg/l
NORG	Concentration of organic nitrogen, mg/l
NTOT	Concentration of total nitrogen, mg/l
PO4	Concentration of phosphate phosphorus, mg/l
PORG	Concentration of organic phosphorus, mg/l
PTOT	Concentration of total phosphorus, mg/l
QNO3	Mass flow of nitrate nitrogen, g/s
QNTOT	Mass flow of total nitrogen, g/s
QPO4	Mass flow of phosphate phosphorus, g/s
QPTOT	Mass flow of total phosphorus, g/s

The output text file can be loaded into QGIS as a text layer, joined on the field *STAIID* to either the *qobs_catchments* table (for substance *Q*) or the *wqobs_catchments* table (for any other substance) under schema *calibration* of PostgreSQL database of the project for the visualisation of calibration metrics in stations.

4.3. Overview of concentration in reaches

Script for the reach concentration overview is *reachConcentrations.py* located in the main *scripts* directory of the project. By running it the mean concentrations and loads of substances in rivers and reservoirs per catchment can be calculated. The resulting text file is determined by variable *output_TXTFILE* while the list of watersheds could be provided in *watershednamelist* (in case of an empty list – for the whole territory of LV). The resulting text file contains columns listed in Table 4; possible “substances” are listed in Table 3.

Table 4: Columns of the reach concentration text file.

Column	Description
CATCHMENTID	Global Id of the catchment (corresponds to field <i>catchmentid</i> of <i>catchments</i> table under schema <i>catchments</i> of PostgreSQL database of project)
SUBSTANCE	Abbreviation of substance (see Table above)
rch	Substance’s concentration in the catchment’s river
rsv	Substance’s concentration in the catchment’s reservoir
tot	Substance’s concentration in the catchment

The output text file can be used for for the visualisation of the concentration in the reaches of watersheds. It should be loaded into QGIS as a text layer, joined on the field CATCHMENTID to either the *catchments* table of QGIS project or the *catchments* table under schema *catchments* of PostgreSQL database of the project.

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