552005

Analysis Report for Preparation of 2008 Culebra Potentiometric Surface Contour Map

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1 Introduction

This report documents the preparation of a contour map and particle track for the Culebra Member of the Rustler Formation in the vicinity of the Waste Isolation Pilot Plant (WIPP), for inclusion in the 2009 Annual Site Environmental Report (ASER). The driver for this analysis is the draft of the Stipulated Final Order sent to the New Mexico Environment Department (NMED) on May 28, 2009 (Moody, 2009).

Beginning with the ensemble of 100 calibrated MODFLOW transmissivity (T), horizontal anisotropy (A), and areal recharge (R) fields (Hart et al., 2009) used in WIPP performance assessment (PA), 3 average parameter fields are used as input to MODFLOW to simulate freshwater heads within and around the WIPP land withdrawal boundary (LWB). PEST is used to adjust a subset of the boundary conditions in the ensemble-average model to obtain the best-fit match between the observed freshwater heads from September 2008 and the model-predicted heads. The output of the averaged, PEST-calibrated MODFLOW model is both contoured and used to compute an advective particle trace forward from the WIPP waste handling shaft.

2 Scientific Approach

2.1 Overview

Steady-state groundwater flow simulations are carried out using the much the same software and approach used in the analysis report for AP-114 Task 7 (Hart et al., 2009) to create the calibrated fields used as inputs – see Table 1 for a summary of all software used in this analysis. The MODFLOW parameter fields (including transmissivity (T), anisotropy (A), and recharge (R)) used here are an ensemble average of the Culebra parameter fields used for WIPP PA in the CRA-2009 performance assessment baseline calculations (PABC). To clearly distinguish between the two MODFLOW models, the original MODFLOW model, which consists of 100 realizations of calibrated parameter fields (Hart et al., 2009), will be referred to as the "PA MODFLOW model". The model derived here from the PA MODFLOW model, used to construct the resulting contour map and particle track, is referred to as the "averaged MODFLOW model". The calibrated model T, A and R input fields, model boundary conditions, and other model input files are appropriately averaged across all 100 calibrated realizations to produce a single averaged steady-state MODFLOW flow model that can be used to predict regional Culebra groundwater flow across the WIPP site.

The calibration process that resulted in the 100 model realizations of the PA MODFLOW model used PEST to adjust spatially variable model parameters, while assuming fixed MODFLOW boundary conditions. The calibration targets for the PA MODFLOW model were both snapshots of undisturbed heads across the site and transient head responses to large-scale pumping tests. Hart et al. (2009) describe the forward model setup and PEST calibration effort for the CRA-2009 PABC. An analogous but much simpler process is used in the averaged MODFLOW model; here PEST is used to modify a subset of the MODFLOW boundary conditions (see boundaries marked in red on Figure 1). The calibration targets for PEST associated with the average MODFLOW model are the observed September 2008 freshwater heads at Culebra monitoring wells. Boundary conditions are modified while holding spatially variable model parameters (T, A, and R) constant; in the calibration of the PA MODFLOW model the boundary conditions were fixed, while adjusting the spatially variable parameters.

Table 1. Software used

Software	Version	Description	Status
MODFLOW-2000	1.6	Flow model	Acquired; qualified under NP 19-1 (Harbaugh et al., 2000)
PEST	9.11	Inverse model	Developed; qualified under NP 19-1 (Doherty, 2002)
DTRKMF	1.00	Particle tracker	Developed; qualified under NP 19-1
Golden Software Surfer	9	Contouring	Commercial off the shelf
Gnuplot	4.2	Plotting	Commercial off the shelf
Microsoft Excel	2007	Plotting	Commercial off the shelf
Python	2.3.4	Scripting Language	Commercial off the shelf

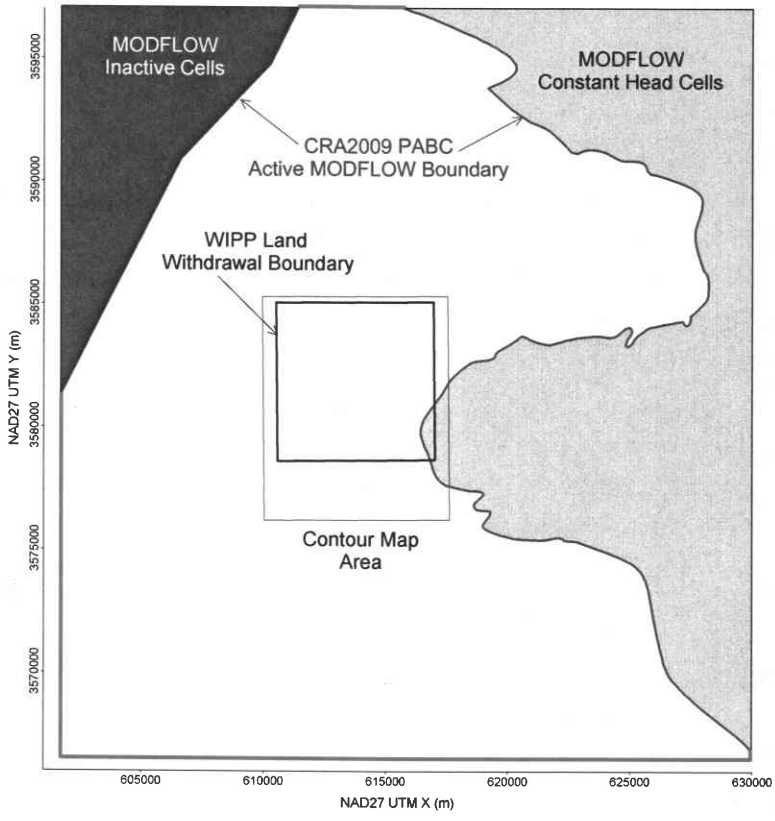

Figure 1. MODFLOW-2000 model domain, adjusted boundary conditions shown in red, contour area outlined in green.

The resulting heads from the PEST-calibrated ensemble-average flow model are contoured over an area surrounding the WIPP site using Surfer (a subset of the complete MODFLOW model domain – see the green rectangle surrounding the WIPP LWB in Figure 1). The trace made by a conservative (i.e., non-dispersive and non-reactive) particle released from the waste handling shaft to the WIPP land withdrawal boundary is computed from the resulting flow field in MODFLOW using DTRKMF, and also plotted with Surfer. Scatter plots statistics summarizing the fit of the PEST-calibrated model to the observed freshwater head at Culebra monitoring wells are created in Gnuplot and Excel. MODFLOW, PEST, DTRKMF, and the Bash and Python scripts written for this work were executed on the PA Linux cluster (`alice.sandia.gov`), while the commercial-off-the-shelf software Surfer, Gnuplot and Excel were executed on a Windows XP desktop computer with an Intel Xenon CPU.

2.2 Creating Ensemble Average MODFLOW Simulation

An ensemble-average MODFLOW model is used to compute both the freshwater head and flow vectors across the model domain; the heads are then contoured and the cell-by-cell flow vectors are used to compute particle tracks. The ensemble-averaged inputs are used to create a single average simulation that produces a single output, rather than averaging the 100 individual outputs of the Culebra flow model used for WIPP PA.

The MODFLOW model grid is a single layer, comprised of 307 rows and 284 columns, each model cell being 100 meters square. The modeling area spans 601,700 to 630,000 meters in the east-west direction, and 3,566,500 to 3,597,100 meters in the north-south direction, both in UTM NAD27 coordinates (zone 13).

The calibrated T, A, and R parameter fields from the PA MODFLOW model were checked out of the CVS repository using the `checkout_average_run_modflow.sh` script (all scripts are listed completely in the Appendix; input files are available on the attached media). The model inputs can be divided into two groups. The first group is the model inputs that are the same across all 100 calibrated realizations; these include the model grid definition, the boundary conditions, and the model solver parameters. The second group is the model inputs that are different for each realization; these include transmissivity (T), horizontal anisotropy (A), and vertical recharge (R). The constant model inputs in the first group are used directly in the averaged MODFLOW model (checked out from the CVS repository), while the inputs in the second group were averaged across all 100 calibrated model realizations using the Python script `average_realizations.py`. All three averaged parameters were log transformed before being averaged, since they vary over multiple orders of magnitude.

2.3 Boundary Conditions

The boundary conditions taken from the PA MODFLOW model are used as the baseline condition, from which PEST calibration proceeds. There are two types of boundary conditions in both MODFLOW models. The first type of condition includes geologic or hydrologic boundaries, which correspond to known physical features in the flow domain. The no-flow boundary along the axis of Nash Draw is a hydrologic boundary (i.e., the boundary along the dark gray region in Figure 1). Also, the constant head boundary along the halite margin corresponds to a geologic boundary (i.e., the eastern irregular boundary adjoining the light gray region in Figure 1).

Physical boundaries are believed to be well known, and are not adjusted in the PEST calibration. The second type of boundary condition includes the constant-head cells along the rest of the model domain; the linear southern, southwestern, and northern boundaries that coincide with the rectangular frame surrounding the model domain are all of this type (shown as a heavy red line in Figure 1). The value of specified head used along this second boundary type is adjusted in the PEST calibration process.

The Python script `boundary_types.py` is used to distinguish between the two different types of specified head boundary conditions based on the specified head value used in the PA MODFLOW model. All constant-head cells (specified by a value of -1 in the MODFLOW IBOUND array from the PA MODFLOW model) that have a starting head value greater than 1000 m (corresponding to the land surface) are left fixed and not adjusted in the PEST optimization. The remaining constant-head cells are distinguished by setting their IBOUND array value to -2 (which is still interpreted as a constant-head value by MODFLOW, but allows simpler discrimination between boundary conditions in scripts elsewhere).

Using the output from `boundary_types.py`, the Python script `surface_02_extrapolate.py` computes the heads at active (IBOUND=1) and adjustable constant-head boundary condition cells (IBOUND=-2), given parameter values for the surface to extrapolate.

2.4 PEST Calibration of Averaged MODFLOW Model to Observations

There are three major types of inputs to PEST. The first type of input includes the set of observed September 2008 freshwater head values used as targets for the PEST calibration. The second class of inputs includes the entire MODFLOW model setup derived from the PA MODFLOW model and described in the previous section, along with any pre- or post-processing scripts or programs needed; this comprises the forward model that PEST runs repeatedly to estimate sensitivities of model outputs to model inputs. The third type of input includes the PEST configuration files, which include parameter and observation groups, indicating which parameters in the MODFLOW model will be adjusted in the inverse simulation.

Freshwater head values used as targets for the PEST calibration were collected in September 2008 (Siegel, 2009) and are summarized in Table 2.

Table 2. Calibration targets used in PEST, from Siegel (2009).

Well I.D.	Date	Adjusted Freshwater Head (ft amsl)	Adjusted Freshwater Head (m amsl)	Density Used (g/cm^3)
AEC-7	09/22/08	3064.06	933.93	1.078
C-2737 (PIP)	09/24/08	3023.61	921.60	1.029
ERDA-9	09/24/08	3033.97	924.75	1.067
H-02b2	09/24/08	3050.51	929.79	1.014
H-03b2	09/24/08	3015.20	919.03	1.042
H-04b	09/24/08	3006.82	916.48	1.015
H-05b	09/22/08	3081.33	939.19	1.093
H-06bR	09/23/08	3074.22	937.02	1.036
H-07b1	09/23/08	2999.24	914.17	1.002
H-09c (PIP)	09/23/08	2997.25	913.56	1.001
H-10c	09/23/08	3024.16	921.72	1.001
H-11b4	09/22/08	3009.92	917.42	1.070
H-12	09/23/08	3007.71	916.75	1.097
H-15R	08/15/08	3045.09	928.14	1.199
H-16	09/25/08	3050.45	929.78	1.039
H-17	09/22/08	3007.52	916.69	1.133
H-19b0	09/24/08	3015.69	919.18	1.068
IMC-461	09/23/08	3046.33	928.52	1.005
SNL-01	09/23/08	3085.69	940.52	1.033
SNL-02	09/23/08	3074.57	937.13	1.012
SNL-03	09/23/08	3081.17	939.14	1.023
SNL-05	09/23/08	3077.77	938.11	1.010
SNL-08	09/22/08	3055.32	931.26	1.103
SNL-09	09/22/08	3057.49	931.92	1.024
SNL-10	09/22/08	3056.14	931.51	1.011
SNL-12	09/23/08	3003.45	915.45	1.005
SNL-13	09/22/08	3012.72	918.28	1.027
SNL-14	09/22/08	3006.17	916.28	1.048
SNL-16	09/22/08	3010.72	917.67	1.010
SNL-17	09/23/08	3007.36	916.64	1.006
SNL-18	09/23/08	3082.59	939.57	1.028
SNL-19	09/23/08	3073.61	936.84	1.003
WIPP-11	09/22/08	3084.85	940.26	1.038
WIPP-13	09/22/08	3081.86	939.35	1.053
WIPP-19	09/24/08	3063.27	933.68	1.044
WIPP-25 (PIP)	09/23/08	3069.43	935.56	1.011
WQSP-1	09/24/08	3078.05	938.19	1.048
WQSP-2	09/24/08	3086.54	940.78	1.048
WQSP-3	09/24/08	3076.08	937.59	1.146
WQSP-4	09/24/08	3016.16	919.33	1.075
WQSP-5	09/24/08	3013.67	918.57	1.025
WQSP-6	09/24/08	3022.96	921.40	1.014

To minimize the number of estimable parameters, and to ensure a degree of smoothness in the constant-head boundary condition values, a parametric surface is used to extrapolate the heads to the estimable boundary conditions. The surface is of the same form described in the analysis report for AP-114 Task 7. The parametric surface is given by the following equation:

$$h_{x,y} = A + B * (y + D * \text{sign}(y) * \text{abs}(y)^{exponent}) + C(E * x^3 + F * x^2 - x)$$

where sign(y) is the function returning 1 for $y>0$, -1 for $y<0$ and 0 for $y=0$ and x and y are coordinates scaled to the range $-1 \leq \{x,y\} \leq 1$. In Hart et al. (2009), the values A=928.0, B=8.0, C=1.2, D=1.0, *exponent*=0.5, E=1.0, and F=-1.0 are used with the above equation.

PEST was then used to estimate the values of parameters *A,B,C,D,E,F*, and *exponent* given the observed heads in Table 2. The Python script `surface_02_extrapolate.py` was used to compute the MODFLOW starting head input file (which is also used to specify the constant-head values) from the parameters *A-F* and *exponent*. Each forward run of the forward model corresponded to a call to the Bash script `run_02_model`. This script called the `surface_02_extrapolate.py` script, the MODFLOW-2000 v1.6 executable, and the qualified PEST utility `mod2obs.exe`, which is used to extract and interpolate model-predicted heads from the MODFLOW output files at observation well locations.

The PEST-specific input files (the third type of input) were generated from the observed heads using the Python script `create_pest_02_input.py`. The PEST input files include the instruction file (how to read the model output), the template files (how to write the model input files), and the PEST control file (listing the ranges and initial values for the estimable parameters and the weights associated with observations).

2.5 Figures Generated from Calibrated MODFLOW Model

The MODFLOW model is run predicatively using the ensemble-averaged model parameters, along with the PEST-calibrated boundary conditions. The resulting cell-by-cell flow budget is then used by DTRKMF to compute a particle track from the waste handling shaft to at least the edge of the WIPP land withdrawal boundary. The Python script `convert_dtrkmf_output_for_surfer.py` converts the IJK cell-based results of DTRKMF into a UTM x and y coordinate system, saving the results in the Surfer blanking file format to facilitate plotting with Surfer. The heads in the binary MODFLOW output file are converted to an ASCII Surfer grid format using the Python script `head_bin2ascii.py`.

The resulting particle trace and contours of the model-predicted head are plotted using Surfer 9 for an area including the WIPP land withdrawal boundary, similar to the region shown in previous versions of the ASER (e.g., see Figure 6.11 in (DOE, 2008)), see green outline in Figure 1. The modeled heads extracted from the MODFLOW output by `mod2obs.exe` are then merged into a common file for plotting using the Python script `merge_observed_modeled_heads.py`.

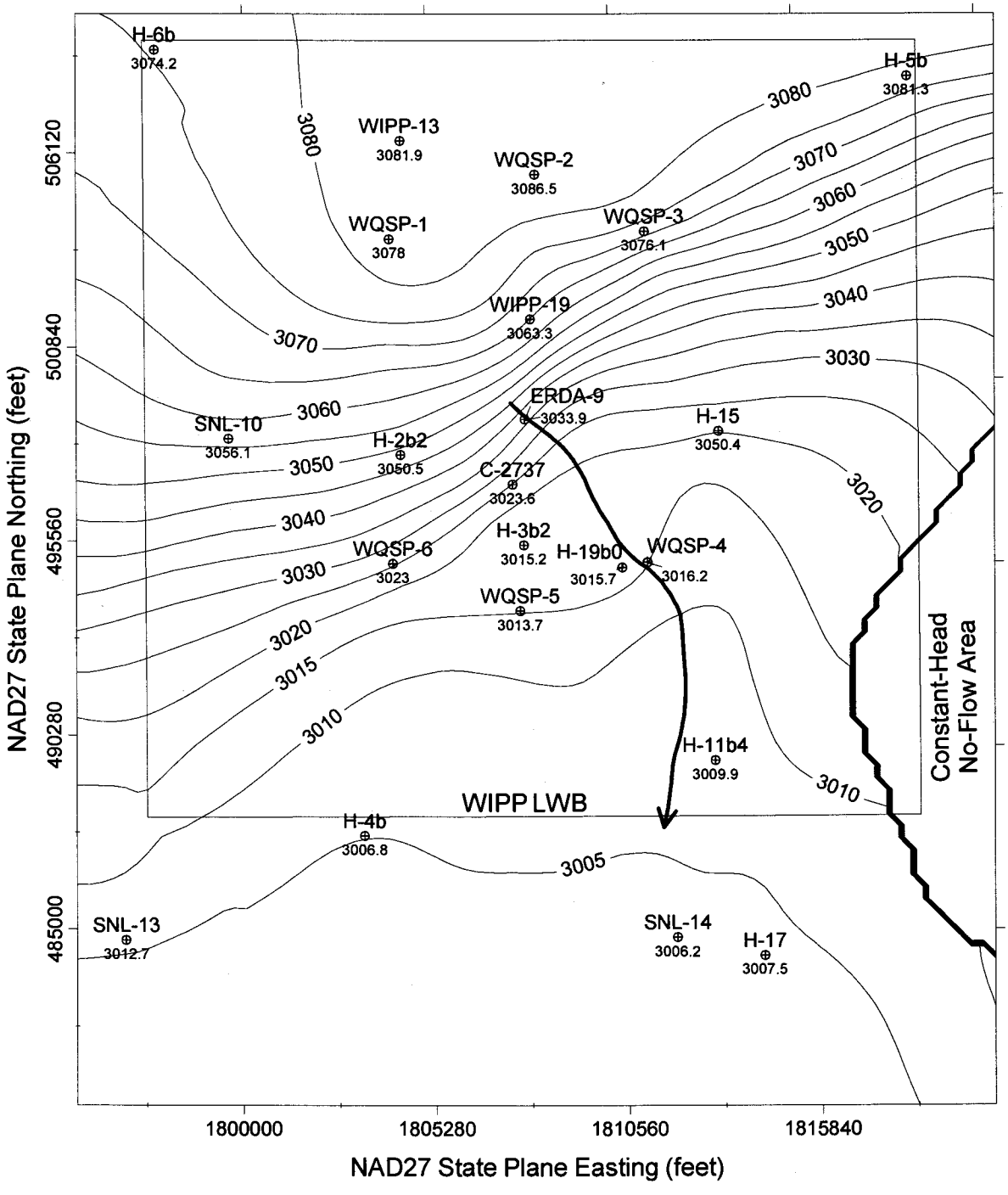

Figure 2. Model-generated September 2008 freshwater head contours with observed head listed at each well (5-foot contour interval) with blue water particle track from waste handling shaft to WIPP LWB

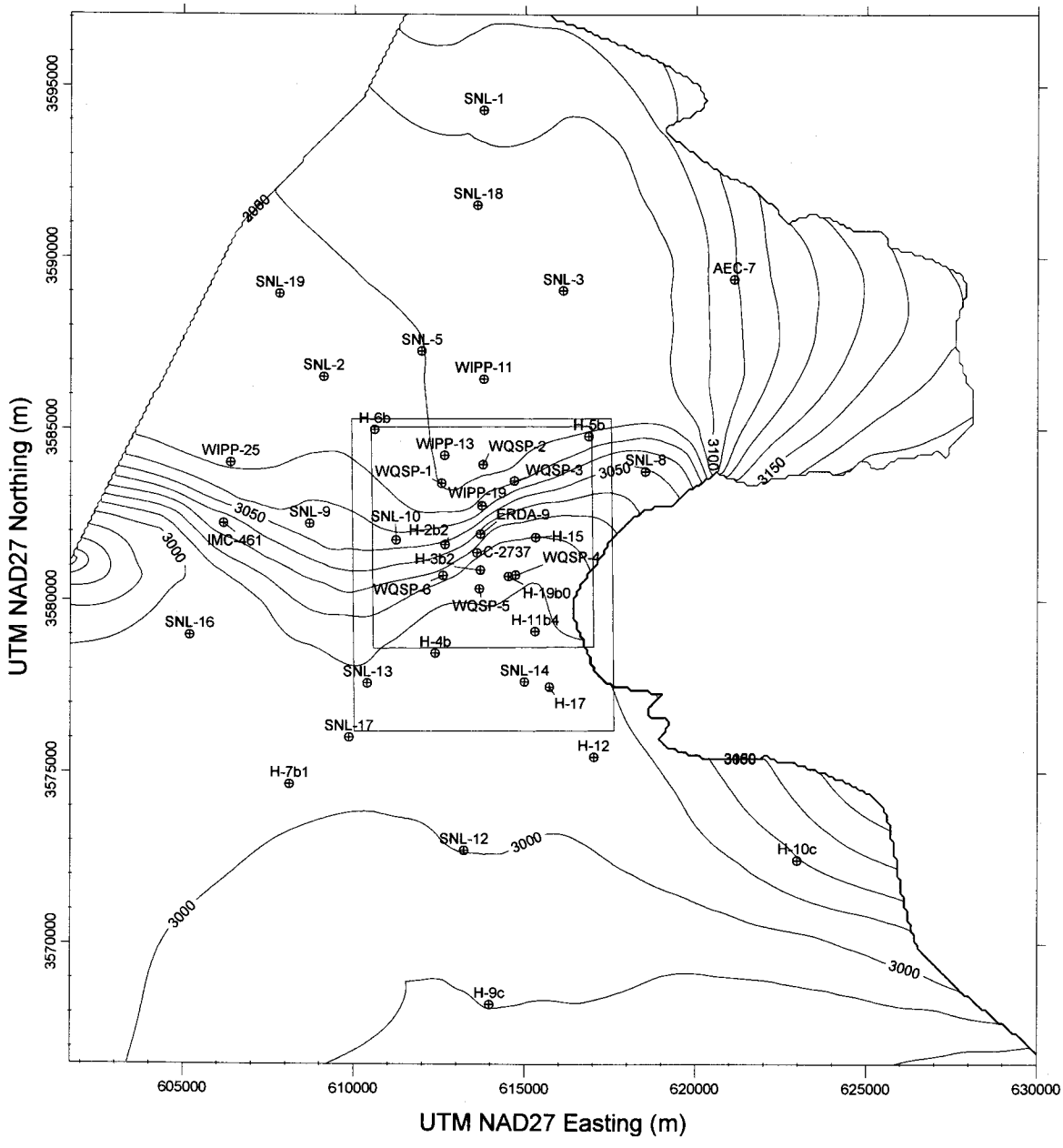

Figure 3. MODFLOW-modeled heads for entire model domain (10-foot contour interval). Green rectangle indicates region contoured in Figure 2, black square is WIPP LWB.

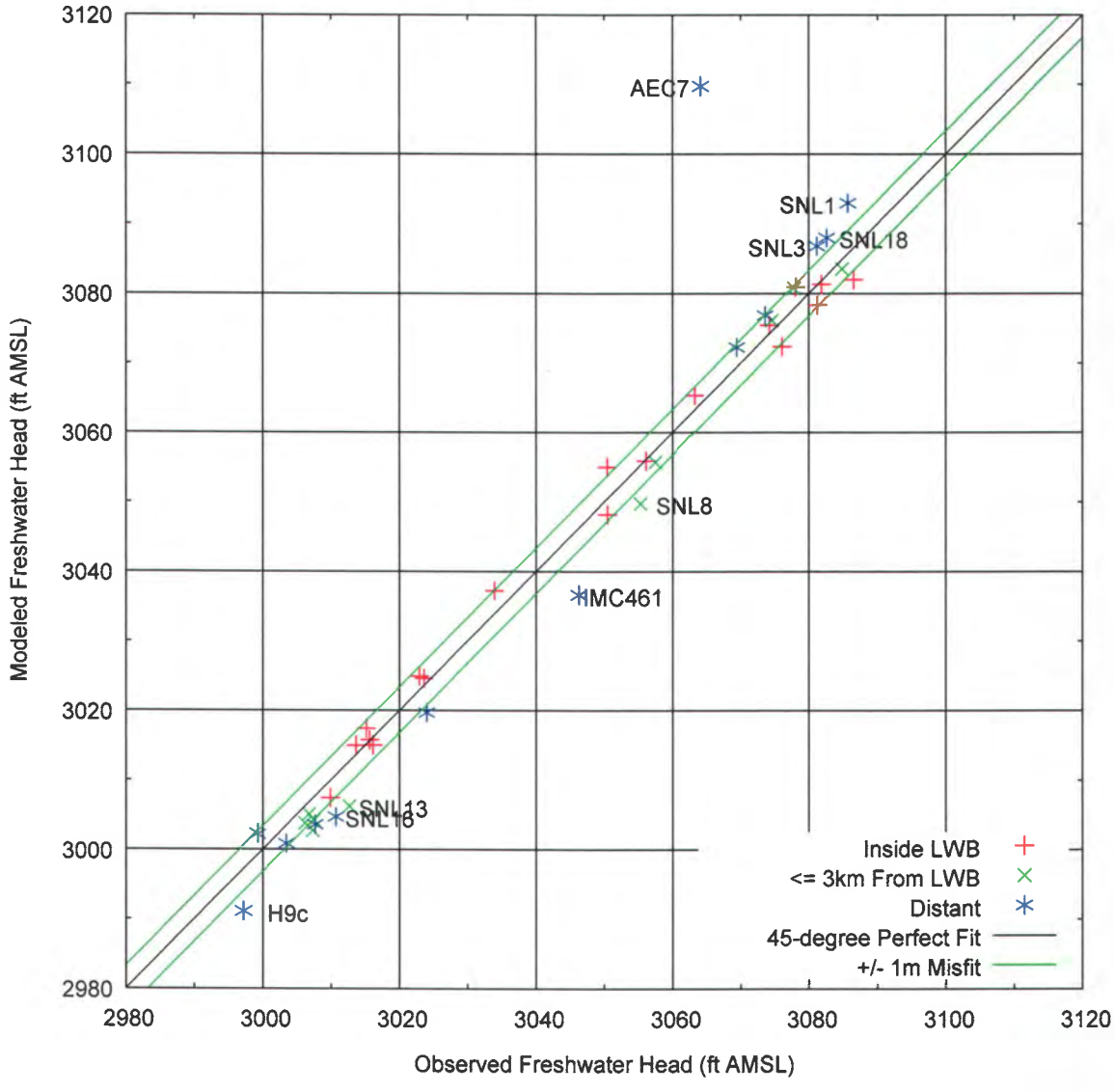

Figure 4. Measured vs. modeled scatter plot for PEST-calibrated MODFLOW-2000 generated heads and September 2008 observed freshwater heads

3 Results

3.1 Freshwater Head Contours

The model-generated freshwater head contours in Figure 2 show the known characteristics of groundwater flow in the Culebra at the WIPP site. There is a roughly east-west trending band of steeper gradients, corresponding to known lower transmissivities. The uncontoured region in the eastern part of the figure corresponds to the portion of the Culebra that is located stratigraphically between halite in other members of the Rustler Formation (Tamarisk Member above and Los Medaños Member below). This region east of the "halite margin" is represented as having high head but extremely low permeability, essentially serving as a no-flow boundary in this area.

3.2 Particle Track

The heavy blue line in Figure 2 shows the DTRKMF-predicted path a water particle would take through the Culebra from the coordinates corresponding to the WIPP waste handling shaft to the land withdrawal boundary (a computed path length of 4.079 km). Assuming a thickness of 4 m for the Culebra and a constant porosity of 16%, the travel time to the WIPP LWB is 5,715 years (output from DTRKMF is adjusted from a 7.75-m Culebra thickness), for an average velocity of 0.71 m/yr.

3.3 Measured vs. Modeled Fit

The scatter plot in Figure 4 shows measured and modeled freshwater heads at the observation locations used in the PEST calibration. The observations are divided into three groups, based on proximity to the WIPP site. Wells within the LWB are represented by red crosses, wells outside but within 3 km of the LWB are represented with green "x"s, and other wells within the MODFLOW model domain but distant from the WIPP site are given by a blue asterisk. These groupings were utilized in the PEST calibration; higher weights (2.5) were given to wells inside the LWB, and lower weights (0.4) were given to wells distant to the WIPP site, while wells in the middle received an intermediate weight (1.0). This allowed PEST to improve the fit of the model to observed heads inside the area contoured in Figure 2, at the expense of fitting wells closer to the boundary conditions (i.e., wells not shown in Figure 2).

The central diagonal line in Figure 4 represents a perfect model fit (1:1 or 45-degree slope); the two lines on either side of this represent a 1-m misfit above or below the perfect fit. Wells more than 1.5 m from the 1:1 line are labeled. AEC-7 has a large misfit (14 m), for two reasons. First, this well has historically had an anomalously low freshwater head elevation, lower than wells around it in all directions. Secondly, it did not have a May 2007 observation (due to ongoing well reconfiguration activities) and therefore was not included as a calibration target in the PA MODFLOW model calibration. The ensemble-average T, A, and R fields used here were not calibrated to accommodate this observation. This well is situated in a low-transmissivity region, and near the constant-head boundary associated with the halite margin, therefore PEST will not be able to improve this fit solely through adjustment of the boundary conditions indicated with red in Figure 1.

The R^2 value for the best-fit line through the measured vs. modeled data inside the WIPP LWB only is 0.991 (computed in Excel) and the slope of this best-fit line is 1.000. The R^2 value for a best-fit line through only the data from the intermediate zone is 0.993, with a slope of 0.999. The R^2 value for the best-fit line through the distant wells only is 0.8975, with a slope of 1.001. The R^2 for the best-fit line through all wells together is 0.944, with a slope of 1.000.

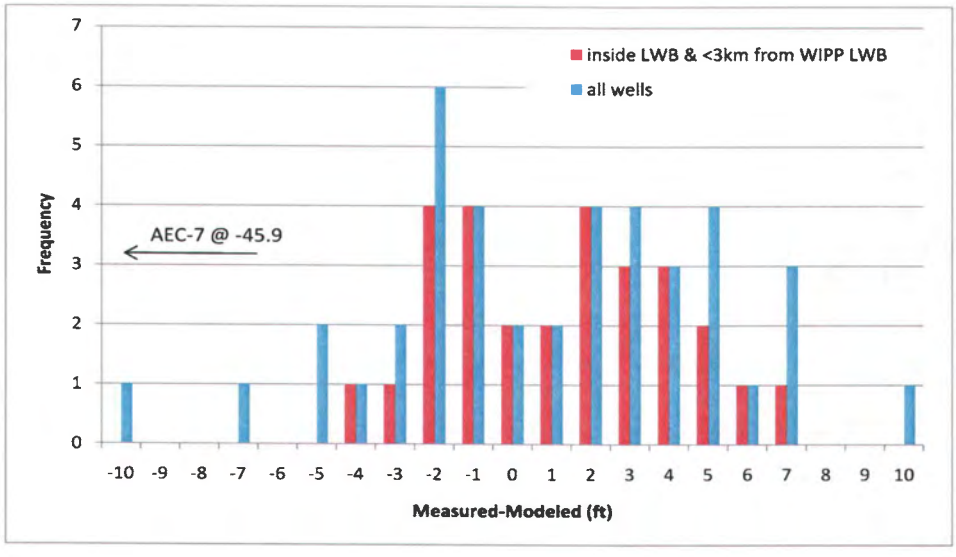

Figure 5. Histogram of Measured-Modeled errors

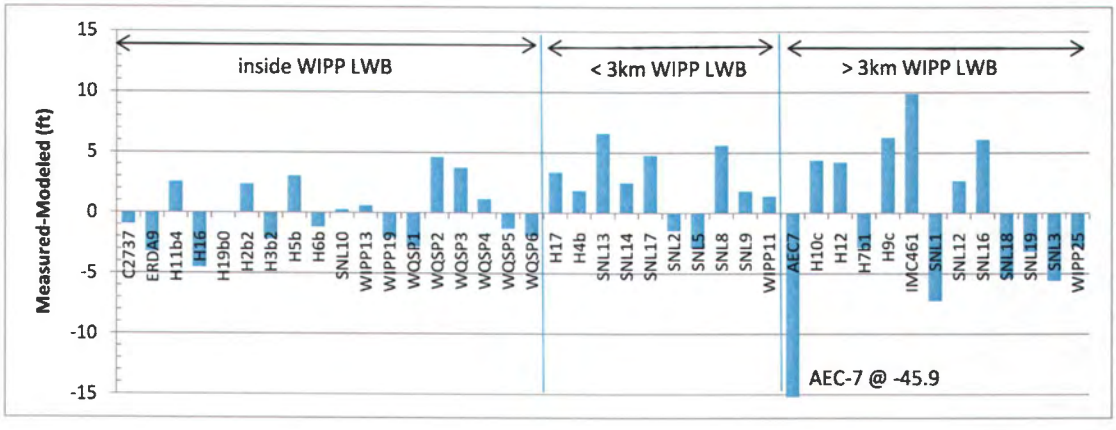

Figure 6. Measured-Modeled errors for each well with an observed September 2008 freshwater head

Figure 5 and Figure 6 show the distribution of errors resulting from the PEST-adjusted fit to observed data. The distribution in Figure 5 is roughly symmetric about 0, indicating there is not a strong bias, although the distribution is approximately bimodal.

Aside from AEC-7, and to a lesser degree some other distant wells whose modeled values do not greatly impact the contours shown in Figure 2, the model fit to the September 2008 observations is very good. The ensemble-average model captures the average Culebra behavior, while the PEST calibration improved the model fit to the specific September 2008 observations.

4 References

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5 Run Control Narrative

This appendix is a narrative describing the calculation process mentioned in the text, which produced the figures given there.

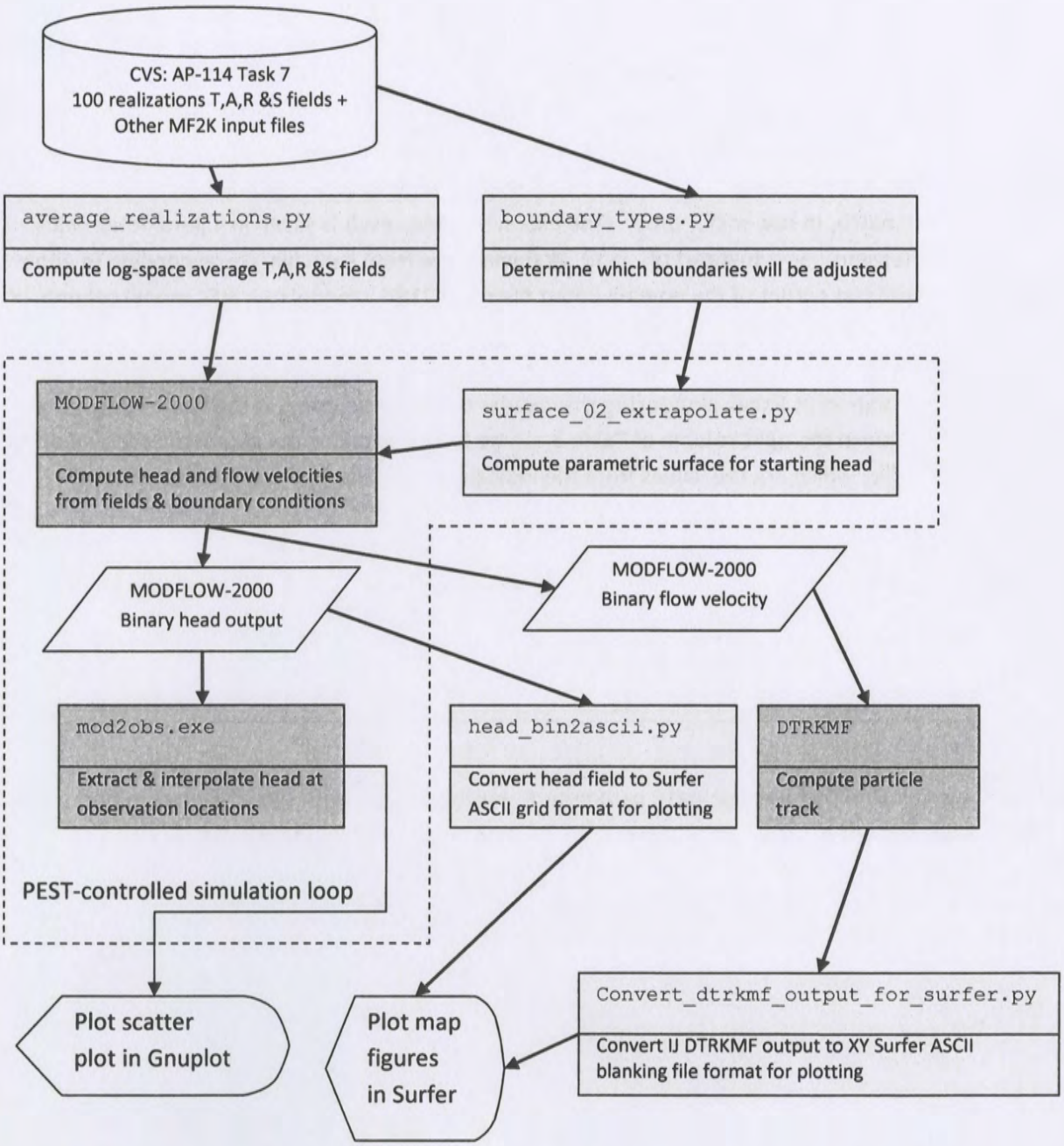

Figure 7. Process flowchart; dark gray indicates qualified programs, light gray are scripts written for this analysis

Figure 7 gives an overview of the driver script `checkout_average_run_modflow.sh` (§A-4.1);

this script first exports the 4 parameter fields (transmissivity (T), anisotropy (A), recharge (R), and storativity (S)) from CVS for each of the 100 realizations of MODFLOW, listed in the file `keepers` (see lines 17-26 of script). Some of the realizations are inside the `Update` or `Update2` subdirectories in CVS, which complicates the directory structure. An equivalent list `keepers_short` is made from `keepers`, and the directories are moved to match the flat directory structure (lines 31-53). At this point the directory structure has been modified but the MODFLOW input files checked out from CVS are unchanged.

The Python script `average_realizations.py` (§A-4.2) is called, which first reads in the `keepers_short` list, then reads in each of the 400 input files and computes the arithmetic average of the base-10 logarithm of the value at each cell across the 100 realizations. The 400 input files are saved as a flattened 2D matrix, in row-major order. The exponentiated result is saved in 4 parameter fields, each with the extension `.avg` instead of `.mod`. A single value from each file, corresponding to either the cell in the southeast corner of the domain (input file row 87188 = model row 307, model column 284 for K and A) or on the west edge of the domain (input file row 45157 = model row 161, model column 1 for R and S) is saved in the text file `parameter_representative_values.txt` to allow checking the calculation in Excel, comparing the results to the value given at the same row of the `.avg` file. The value in the right column of Table 3 can be found by taking the geometric average of the values in the text file, which are the values from the indicated line of each of the 100 realizations.

Table 3. Averaged values for representative model cells

Field	Input file row	Model row	Model column	Geometric average
K	87188	307	284	9.2583577E-09
A	87188	307	284	9.6317478E-01
R	45157	161	1	1.4970689E-19
S	45157	161	1	4.0388352E-03

Figure 8 shows plots of the average log10 parameters, which compare with similar figures in Hart (2009); inactive regions <1.0E-15 were reset to 1.0 to improve the plotted color scale. The rest of the calculations are done with these averaged fields.

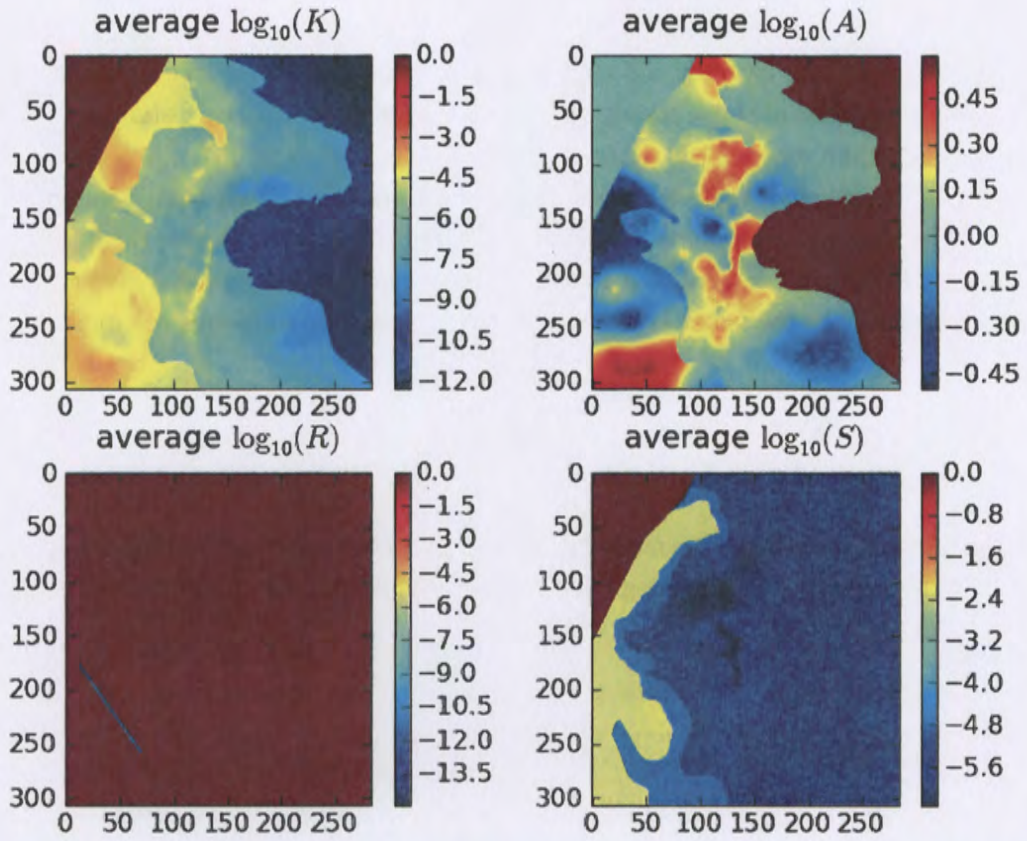

Figure 8. Plots of base-10 logarithms of average parameter fields; rows and columns are labeled on edges of figures.

Next a subdirectory is created, and the averaged MODFLOW model is run without any modifications by PEST. Subsequently, another directory will be created where PEST will be run to improve the fit of the model to observed heads at well locations.

The next portion of the driving script `checkout_average_run_modflow.sh` links copies of the input files needed to run MODFLOW-2000 and DTRKMF into the `original_average` run directory. Then MODFLOW-2000 is run with the name file `mf2k_head.nam`, producing binary head (`modeled_head.bin`) and binary cell-by-cell flow budget (`modeled_flow.bud`) files, as well as a text listing file (`modeled_head.lst`). DTRKMF is then run with the input files `dtrkmf.in` and `wippctrl.inp`, which utilizes the cell-by-cell budget file written by MODFLOW to generate a particle track output file, `dtrk.out`. The input file `wippctrl.inp` specifies the starting location of the particle in DTRKMF face-centered cell coordinates, the porosity of the aquifer (here 16%), and the coordinates of the corners of the WIPP LWB, since the calculation stops when the particle reaches the LWB.

The Python script `head_bin2ascii.py` (§A-4.7) converts the MODFLOW binary head file, which includes the steady-state head at every element in the flow model domain (307 rows × 284 columns) into a Surfer ASCII grid file format. This file is simply contoured in Surfer, no interpolation or gridding is needed. The Python script `convert_dtrkmf_output_for_surfer.py` (§A-4.9) reads the

DTRKMF output file `dtrk.out` and does two things. First it converts the row, column format of this output file to an X,Y format suitable for plotting, and second it converts the effective thickness of the Culebra from 7.75m to 4m. The following table shows the first 10 lines of the `dtrk.out` and the corresponding output of the Python script `dtrk_output_original_average.bln`. The first three columns of `dtrk.out` (top half of Table 4) after the header are cumulative time (red), column (blue), and row (green). The three columns in the blanking file (second half of Table 4) after the header are UTM NDA27 X (blue), UTM NAD27 Y (green), and adjusted cumulative time (red, which is faster than the original cumulative travel time by the factor 4/7.75=0.516). The conversion from row, column to X, Y is

$$X = 601700.0 + 100.0 * column$$
$$Y = 3597100.0 - 100.0 * row$$

since the I,J origin is the northwest corner of the model domain (601700,3597100), while the X,Y origin is the southwest corner of the domain. The blanking file is plotted directly in Surfer, since it now has the same coordinates as the ASCII head file.

Table 4. Comparison of first 10 lines of DTRKMF output and converted Surfer blanking file for `original_average`

```
         1         159
 0.00000000E+00  118.79  150.21 1.18790000E+04 1.50210000E+04 0.00000000E+00 1.85168267E-01 1.59999996E-01 1.00000000E+00
 5.53946616E+01  118.86  150.29 1.18859872E+04 1.50285080E+04 1.02562574E+01 1.85130032E-01 1.59999996E-01 1.00000000E+00
 1.10789323E+02  118.93  150.36 1.18929942E+04 1.50359947E+04 2.05104788E+01 1.85094756E-01 1.59999996E-01 1.00000000E+00
 1.66017959E+02  119.00  150.43 1.19000000E+04 1.50434379E+04 3.07321029E+01 1.85062532E-01 1.59999996E-01 1.00000000E+00
 3.27990509E+02  119.21  150.62 1.19206651E+04 1.50624751E+04 5.88294962E+01 1.73534671E-01 1.59999996E-01 1.00000000E+00
 4.89963060E+02  119.42  150.81 1.19415109E+04 1.50813473E+04 8.69490492E+01 1.73684593E-01 1.59999996E-01 1.00000000E+00
 6.51450155E+02  119.62  151.00 1.19624759E+04 1.51000000E+04 1.15010608E+02 1.73860152E-01 1.59999996E-01 1.00000000E+00
 7.40581455E+02  119.75  151.10 1.19749757E+04 1.51102419E+04 1.31170520E+02 1.81333000E-01 1.59999996E-01 1.00000000E+00
 8.29712755E+02  119.87  151.20 1.19874963E+04 1.51204665E+04 1.47335525E+02 1.81390626E-01 1.59999996E-01 1.00000000E+00
```

```
159,1
613579.0,3582079.0,0.00000000e+00
613586.0,3582071.0,2.85907931e+01
613593.0,3582064.0,5.71815861e+01
613600.0,3582057.0,8.56866885e+01
613621.0,3582038.0,1.69285424e+02
613642.0,3582019.0,2.52884160e+02
613662.0,3582000.0,3.36232338e+02
613675.0,3581990.0,3.82235590e+02
613687.0,3581980.0,4.28238841e+02
```

The PEST utility script `mod2obs.exe` is run to extract and interpolate the model-predicted heads at observation locations. The input files for mod2obs.exe were taken from AP-114 Task 7 in CVS. The observed head file has the wells and observed heads from September 2008, but is otherwise the same as that used in the model calibration in AP-114. The Python script `merge_observed_modeled_heads.py` (§A-4.8) simply puts the results from `mod2obs.exe` and the original observed heads in a single file together for plotting in Gnuplot and Excel.

A similar process to that described so far in this narrative is carried out in a new directory called `pest_02` (beginning line 146 of the driver script). The PEST calibration is carried out there, to keep it separate from the `original_average` simulation. Now the Python script `boundary_types.py` (§A-4.3) is also run, to create a new MODFLOW IBOUND array, where the two different types of boundary conditions are differentiated. This Python script uses the MODFLOW IBOUND array (`init_bnds_orig.inf` first ⅓ of Table 5) and the initial head array (`init_head_orig.mod`

middle ⅓ of Table 5) as inputs, and writes a new MODFLOW IBOUND array (`init_bnds.inf` bottom ⅓ of Table 5) with constant head nodes indicated in red in Figure 1 marked as -2 and other constant head nodes remaining as -1 as output. The script differentiates between these two types of boundary conditions by checking if the starting head is <1000m. Starting heads >1000m are associated with the constant head areas to the east of the halite margins (lighter gray areas in Figure 1).

Table 5. Input IBOUND, starting head, and output IBOUND array data corresponding to first row of MODFLOW model

0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 -1 -1
943 943 943 943 943 943 943 943 943 943 943 943 943 943 943 943 943 943 943 943 943 943 943 943 943 944 1087 1085 1085 1082 1081 1080 1079 1078 1078 1077 1077 1077 1077 1077 1077 1078 1078 1077 1078 1079 1081 1083 1084 1086 1086 1088 1090 1092 1095 1096 1098 1099 1099 1100 1100 1100 1101 1101 1101 1102 1104 1104 1104 1104 1104 1104 1105 1105 1105 1105 1106 1106 1106 1107 1107 1109 1110 1110 1112 1113 1113 1114 1114 1115 1115 1115 1116 1116 1116 1117 1118 1119 1120 1121 1122 1123 1124 1125 1126 1127 1127 1129 1129 1130 1131 1132 1132 1133 1133 1134 1135 1136 1137 1137 1138 1139 1139 1140 1140 1140 1141 1142 1142 1142 1142 1142 1143 1143 1144 1144 1144 1145 1145 1146 1147 1147 1147 1147 1148 1148 1147 1146 1144 1143 1143 1145 1147 1148 1149 1150 1149 1149 1148 1149 1149 1151 1151 1151 1150 1152 1153 1154 1155
0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -2 -1 -1

Table 5 shows the data corresponding to the northernmost row of the MODFLOW model domain (284 entries long) for the two input files and one output file. In the top IBOUND array, the values are either 0 or -1, indicating either inactive (the region northwest of the no-flow area shown in dark gray in Figure 1) or constant head (both red and light gray cells in Figure 1). The first 284 values from the initial head file (reformatted from scientific notation to integers to facilitate printing) show a jump from approximately 944 (in blue) to >1000 (in red). These same cells are colored in the output, showing how the initial head value is used to distinguish between the two types of constant head boundaries. MODFLOW treats any cells as constant head which have an IBOUND entry < 0, so both -2 and -1 are the same to MODFLOW, but allow distinguishing between them in the Python script which extrapolates the heads to the boundaries.

The required PEST input files are created by the Python script `create_pest_02_input.py` (§A-4.4). This script writes **1)** the PEST instruction file (`modeled_head.ins`), which shows PEST how to extracting the model-predicted heads from the `mod2obs.exe` output; **2)** the PEST template file

(surface_par_params.ptf), which shows PEST how to write the input file for the surface extrapolation script; **3)** the PEST parameter file (surface_par_params.par), which lists the starting parameter values to use when checking the PEST input; **4)** the PEST control file (bc_adjust_2008ASER.pst), which has PEST-related parameters, definitions of extrapolation surface parameters, and the observations and weights that PEST is adjusting the model inputs to fit. The observed heads are read as an input file in the PEST borehole sample file format (meas_head_2008ASER.smp), and the weights are read in from the input file (obs_loc_2008ASER.dat).

PEST runs the "forward model" many times, adjusting inputs and reading the resulting outputs using the instruction and template files created above. The forward model actually consists of a Bash shell script (run_02_model) that simply calls a pre-processing Python script surface_02_extrapolate.py (§A-4.5), the MODFLOW-2000 executable, and the PEST utility mod2obs.exe as a post-processing step. The script redirects the output of each step to /dev/null to minimize screen output while running PEST, since PEST will run the forward model many dozens of times.

The pre-processing Python script surface_02_extrapolate.py reads the new IBOUND array created in a previous step (with -2 now indicating which constant head boundaries should be modified), the initial head file used in AP-114 Task 7 (init_head_orig.mod), two files listing the relative X and Y coordinates of the model cells (rel_{x,y}_coord.dat), and an input file listing the coefficients of the parametric equation used to define the initial head surface. This script then cycles over the elements in the domain, writing the original starting head value if the IBOUND value is -1 or 0, and writing the value corresponding to the parametric equation if the IBOUND value is -2 or 1. Using the parameters corresponding to those used in AP-114 Task 7, the output starting head file should be identical to that used in AP-114 Task 7.

After PEST has converged to the optimum solution for the given observed heads and weights, it runs the forward model one more time with the optimum parameters. The post-processing Python scripts for creating the Surfer ASCII grid file and Surfer blanking file from the MODFLOW and DTRKMF output are run and the results are plotted in Surfer.

A-1 Input Files

input file name	file type	description
average_realizations.py	Python script	average 100 realizations
boundary_types.py	Python script	distinguish different BC types
checkout_average_run_modflow.sh	Bash script	main routine: checkout files, run MODFLOW and PEST, call Python scripts
convert_dtrkmf_output_for_surfer.py	Python script	convert DTRKMF IJ output to Surfer X,Y blanking format
create_pest_02_input.py	Python script	create PEST input files from observed data
dtrkmf.in	input listing	responses to DTRKMF prompts
head_bin2ascii.py	Python script	convert MODFLOW binary output to Surfer ASCII grid format
keepers	input	listing of 100 realizations from CVS
meas_head_2008ASER.smp	input	observed September 2008 heads in `mod2obs.exe` bore sample file format
merge_observed_modeled_heads.py	Python script	paste observed head and model-generated heads into one file
mod2obs_files.dat	file listing	files needed to run `mod2obs.exe`
mod2obs_head.in	input listing	responses to `mod2obs.exe` prompts
modflow_files.dat	file listing	files needed to run MODFLOW
obs_loc_2008ASER.dat	input	listing of wells and geographic groupings
pest_02_files.dat	file listing	files needed to run PEST
rel_x_coord.dat	input	relative coordinate $1 \leq x \leq 1$
rel_y_coord.dat	input	relative coordinate $1 \leq y \leq 1$
run_02_model	Bash script	PEST model: execute MODFLOW and do pre- and post-processing
settings.fig	input	`mod2obs.exe` input file
spec_domain.spc	input	`mod2obs.exe` input file
spec_wells.crd	input	`mod2obs.exe` input file
surface_02_extrapolate.py	Python script	compute starting head from parameter and coordinate inputs
wippctrl.inp	input	DTRKMF input file

Table A-1: Input Files

A-2 Output Files

output file name	file type	description
ASER_boundary{,_state_plane}.bln	Surfer blanking file	coordinates defining contouring area in UTM and state plane coordinates
contour_map_with_particle_feet.srf	Surfer Plot file	plot of local contours
contour_map_with_particle.emf	enhanced metafile	plot of local contours
contours_whole_domain.srf	Surfer Plot file	plot of regional contours
contours_whole_domain.emf	enhanced metafile	plot of regional contours
dtrk_output_pest_02{,_state_plane}.bln	Surfer blanking file	coords defining particle track: output from convert_dtrkmf_output_for_surfer.py in UTM & state plane (converted via Surfer)
modeled_head_pest_02.grd	Surfer grid file	model-generated heads: output from head_bin2ascii.py
modeled_head_pest_02_ ... contours_feet_UTM.grd	Surfer grid file	modeled_head_pest_02.grd contour values in feet
modeled_head_pest_02_... contours_feet_state_plane.grd	Surfer grid file	modeled_head_pest_02.grd projected to state plane coords using Surfer
modeled_vs_observed_head.xlsx	Excel spreadsheet	modeled_vs_observed_head_pest_02.txt plotted for histograms & regression R^2 values
modeled_vs_observed_head_pest_02.txt	text file	output of merge_observed_modeled_heads.py
parameter_representative_values.txt	run control	point extracted from each realization
plot_scatter_plots.gnu	gnuplot input	input for plotting scatter_pest_02_feet.emf
scatter_pest_02_feet.emf	enhanced metafile	output from gnuplot
wipp_boundary{,_state_plane}.bln	Surfer blanking file	coordinates defining WIPP LWB in UTM and state plane coordinates
well_data_with_names_and_observed.dat	coords/names	well coordinates and names for plotting

Table A-2: Output Files

A-3 Directory Listings

```
C:\>dir input
 Volume in drive C is DriveC
 Volume Serial Number is 542A-10F7
 Directory of C:\input

08/28/2009  03:26 PM    <DIR>          .
08/28/2009  03:26 PM    <DIR>          ..
08/28/2009  03:06 PM             2,057 average_realizations.py
08/25/2009  01:00 PM             2,088 boundary_types.py
08/25/2009  12:58 PM             6,650 checkout_average_run_modflow.sh
08/25/2009  01:04 PM               630 convert_dtrkmf_output_for_surfer.py
08/25/2009  01:01 PM             2,514 create_pest_02_input.py
08/07/2009  11:17 AM                48 dtrkmf.in
08/25/2009  01:03 PM             3,862 head_bin2ascii.py
08/07/2009  11:17 AM             1,091 keepers
08/26/2009  11:04 AM             1,497 meas_head_2008ASER.smp
08/07/2009  12:16 PM               437 merge_observed_modeled_heads.py
08/07/2009  11:18 AM                76 mod2obs_files.dat
08/07/2009  11:25 AM               138 mod2obs_head.in
08/07/2009  11:18 AM               372 modflow_files.dat
08/07/2009  11:26 AM               390 obs_loc_2008ASER.dat
08/07/2009  11:18 AM               182 pest_02_files.dat
08/07/2009  11:18 AM         2,397,670 rel_x_coord.dat
08/07/2009  11:18 AM         2,397,528 rel_y_coord.dat
08/25/2009  01:02 PM               295 run_02_model
08/07/2009  11:19 AM                26 settings.fig
08/07/2009  11:19 AM                47 spec_domain.spc
08/07/2009  11:19 AM             1,659 spec_wells.crd
08/25/2009  01:02 PM             2,463 surface_02_extrapolate.py
08/07/2009  11:17 AM               506 wippctrl.inp
              23 File(s)      4,822,226 bytes
               2 Dir(s)  153,963,347,968 bytes free
```

```
C:\output>dir
 Volume in drive C is DriveC
 Volume Serial Number is 542A-10F7
 Directory of C:\output

08/28/2009  03:47 PM    <DIR>          .
08/28/2009  03:47 PM    <DIR>          ..
08/28/2009  01:30 PM           548,148 contours_whole_domain.emf
08/28/2009  01:30 PM           947,380 contours_whole_domain.srf
08/28/2009  01:43 PM           684,880 contour_map_with_particle.emf
08/28/2009  01:43 PM         1,183,187 contour_map_with_particle_feet.srf
08/26/2009  11:44 AM         1,133,833 modeled_head_pest_02.grd
08/28/2009  01:33 PM           697,604 modeled_head_pest_02_contours_feet_state_plane.grd
08/28/2009  01:29 PM         1,577,862 modeled_head_pest_02_contours_feet_UTM.grd
08/26/2009  11:44 AM             1,077 modeled_vs_observed_head_pest_02.txt
08/28/2009  01:11 PM            30,102 modeled_vs_observed_head_pest_02.xlsx
08/28/2009  03:08 PM             9,081 parameter_representative_values.txt
08/28/2009  12:46 PM             2,451 plot_scatter_plots.gnu
08/28/2009  12:48 PM            33,900 scatter_pest_02_feet.emf
08/28/2009  03:18 PM    <DIR>          state_plane_data
08/28/2009  03:18 PM    <DIR>          utm_data
08/28/2009  01:25 PM             2,701 well_data_with_names_and_observed.dat
              13 File(s)      6,852,206 bytes
               4 Dir(s)  153,963,302,912 bytes free
C:\>dir output\state_plane_data
 Volume in drive C is DriveC
 Volume Serial Number is 542A-10F7
 Directory of C:\output\state_plane_data

08/28/2009  03:18 PM    <DIR>          .
08/28/2009  03:18 PM    <DIR>          ..
08/24/2009  12:59 PM               168 ASER_boundary_state_plane.bln
08/28/2009  01:36 PM             6,934 dtrk_output_pest_02_state_plane.bln
08/24/2009  12:59 PM               166 wipp_boundary_state_plane.bln
               3 File(s)          7,268 bytes
               2 Dir(s)  153,963,360,256 bytes free

C:\>dir output\utm_data
 Volume in drive C is DriveC
 Volume Serial Number is 542A-10F7
 Directory of C:\output\utm_data

08/28/2009  03:18 PM    <DIR>          .
08/28/2009  03:18 PM    <DIR>          ..
07/31/2009  03:57 PM                85 ASER_boundary.bln
08/26/2009  11:44 AM             5,412 dtrk_output_pest_02.bln
04/13/2009  10:08 AM               105 wipp_boundary.bln
               3 File(s)          5,602 bytes
               2 Dir(s)  153,963,360,256 bytes free
```

A-4 Script Source Listing

A-4.1 Bash shell script `checkout_average_run_modflow.sh`

```
#!/bin/bash

# this script makes the following directory substructure
#
#  current_dir \------ Outputs  (calibrated parameter fields - INPUTS)
#               \----- Inputs   (other modflow files - INPUTS)
#                \---- original_average  (foward sim using average fields)
#                 \--- bin      (MODFLOW and DTRKMF binaries)
#                  \-- pest_0?  (pest-adjusted results)

echo " ^^^^^^^^^^^^^^^^^^^^ "
echo " checking out T fields"
echo " ^^^^^^^^^^^^^^^^^^^^ "

# these will checkout the calibrated parameter-field data into subdirectories
# checkout things that are different for each of the 100 realiztaions
for d in `cat keepers`
do
  cvs -d /nfs/data/CVSLIB/Tfields checkout Outputs/${d}/modeled_{K,A,R,S}_field.mod
done

# checkout MODFLOW input files that are constant for across all realizations
cvs -d /nfs/data/CVSLIB/Tfields checkout Inputs/data/elev_{top,bot}.mod
cvs -d /nfs/data/CVSLIB/Tfields checkout Inputs/data/init_{bnds.inf,head.mod}
cvs -d /nfs/data/CVSLIB/Tfields checkout Inputs/modflow/mf2k_culebra.{lmg,lpf}
cvs -d /nfs/data/CVSLIB/Tfields checkout Inputs/modflow/mf2k_head.{ba6,nam,oc,dis,rch}

# modify the path of "updated" T-fields, so they are all at the
# same level in the directory structure (simplifying scripts elsewhere)

if [ -a keepers_short ]
then
    rm keepers_short
fi
touch keepers_short

for d in `cat keepers`
do
  bn=`basename ${d}`
  # test whether it is a compount path
  if [ ${d} != ${bn} ]
      then
      dn=`dirname ${d}`
      mv ./Outputs/${d} ./Outputs/

      # put an empty file in the directory to indicate
      # what the directory was previously named
      touch ./Outputs/${bn}/${dn}
```

```
  fi

  # create a keepers list without directories
  echo ${bn} >> keepers_short
done

echo " ^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^ "
echo " perform averaging across all realizations "
echo " ^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^ "

python average_realizations.py

# checkout MODFLOW / DTRKMF executables
cvs -d /nfs/data/CVSLIB/MODFLOW2K checkout bin/mf2k/mf2k_1.6.release
cvs -d /nfs/data/CVSLIB/MODFLOW2K checkout bin/dtrkmf/dtrkmf_v0100

# check out pest and obs2mod binaries
cd bin
cvs -d /nfs/data/CVSLIB/PEST checkout Builds/Linux/pest.exe
cvs -d /nfs/data/CVSLIB/PEST checkout Builds/Linux/mod2obs.exe
cd ..

echo " ^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^ "
echo " setup copies of files constant between all realizations "
echo " ^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^ "

# directory for putting original base-case results in
od=original_average

if [ -d ${od} ]
then
    echo ${od}" directory exists: removing and re-creating"
    rm -rf ${od}
fi

mkdir ${od}
cd ${od}
echo `pwd`

# link to unchanged input files
for file in `cat ../modflow_files.dat`
do
  ln -sf ${file} .
done

# link to averaged files computed in previous step
for f in {A,R,K,S}
do
  ln -sf ../modeled_${f}_field.avg ./modeled_${f}_field.mod
done

```

```
ln -sf elev_top.mod fort.33
ln -sf elev_bot.mod fort.34

echo "^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^"
echo " run original MODFLOW and DTRKMF and export results for plotting"
echo "^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^"

# run MODFLOW, producing average head and CCF
../bin/mf2k/mf2k_1.6.release mf2k_head.nam

# run DTRKMF, producing particle track (from ccf)
../bin/dtrkmf/dtrkmf_v0100 <dtrkmf.in

# convert binary MODFLOW head output to Surfer ascii grid file format
ln -sf ../head_bin2ascii.py .
python head_bin2ascii.py
mv modeled_head_asciihed.grd modeled_head_${od}.grd

# convert DTRKMF output from cells to X,Y and
# save in Surfer blanking file format
ln -sf ../convert_dtrkmf_output_for_surfer.py .
python convert_dtrkmf_output_for_surfer.py
mv dtrk_output.bln dtrk_output_${od}.bln

# extract head results at well locations and merge with observed
# head file for easy scatter plotting in Excel (tab delimited)
for file in `cat ../mod2obs_files.dat`
do
  ln -sf ${file} .
done

ln -sf ../meas_head_2008ASER.smp .
ln -sf ../obs_loc_2008ASER.dat .
../bin/Builds/Linux/mod2obs.exe <mod2obs_head.in
ln -sf ../merge_observed_modeled_heads.py
python merge_observed_modeled_heads.py
mv both_heads.smp modeled_vs_observed_head_${od}.txt

# go back down into root directory
cd ..
echo `pwd`

echo "^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^"
echo " setup and run PEST to optimize parametric surface to set BC "
echo "^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^^"

for p in pest_02
do

  if [ -d ${p} ]
      then
```

```
151         echo ${p}" directory exists: removing and re-creating"
152         rm -rf ${p}
153     fi
154
155     mkdir ${p}
156     cd ${p}
157     echo `pwd`
158
159     # link to unchanged input files
160     for file in `cat ../modflow_files.dat`
161       do
162       ln -sf ${file} .
163     done
164
165     # link to averaged files computed in previous step
166     for f in {A,R,K,S}
167       do
168       ln -sf ../modeled_${f}_field.avg ./modeled_${f}_field.mod
169     done
170
171     # link to mod2obs files (needed for pest)
172     for file in `cat ../mod2obs_files.dat`
173       do
174       ln -sf ${file} .
175     done
176
177     # link to pest files
178     for file in `cat ../${p}_files.dat`
179       do
180       ln -s ${file} .
181     done
182
183     # rename 'original' versions of files to be modified by pest
184     rm init_head.mod
185     ln -sf ../Inputs/data/init_head.mod ./init_head_orig.mod
186     rm init_bnds.inf
187     ln -sf ../Inputs/data/init_bnds.inf ./init_bnds_orig.inf
188
189     # create new ibound array for easier modification during PEST
190     # optimization iterations
191     python boundary_types.py
192
193     # create the necessary input files from observations
194     python create_${p}_input.py
195
196     # run pest
197     ../bin/Builds/Linux/pest.exe bc_adjust_2008ASER
198
199     # last output files should be best run
200     # extract all the stuff from that output
201     ########################################
```

```
202
203     ln -sf elev_top.mod fort.33
204     ln -sf elev_bot.mod fort.34
205
206     ../bin/dtrkmf/dtrkmf_v0100 <dtrkmf.in
207
208     ln -sf ../head_bin2ascii.py .
209     python head_bin2ascii.py
210     mv modeled_head_asciihed.grd modeled_head_${p}.grd
211
212     ln -sf ../convert_dtrkmf_output_for_surfer.py .
213     python convert_dtrkmf_output_for_surfer.py
214     mv dtrk_output.bln dtrk_output_${p}.bln
215
216     for file in `cat ../mod2obs_files.dat`
217     do
218       ln -sf ${file} .
219     done
220
221     ../bin/Builds/Linux/mod2obs.exe <mod2obs_head.in
222     ln -sf ../merge_observed_modeled_heads.py
223     python merge_observed_modeled_heads.py
224     mv both_heads.smp modeled_vs_observed_head_${p}.txt
225
226     cd ..
227   done
```

A-4.2 Python script `average_realizations.py`

```
from math import log10,pow

nrow = 307
ncol = 284
nel = nrow*ncol
nfr = 100  # number of fields (realizations)
nft = 4    # number of field types

debug = True  # set to True to get output described in RunControl narrative

def floatload(filename):
    """Reads file (a list of strings, one per row) into a list of strings."""
    f = open(filename,'r')
    m = [float(line.rstrip()) for line in f]
    f.close()
    return m

types = ['K','A','R','S']

# get list of 100 best calibrated fields
flist = open('keepers_short','r')
runs = flist.read().strip().split('\n')
flist.close()

# initialize to help speed lists up a bit
# nfr (100) realizations of each
fields = []
for i in xrange(nft):
    fields.append([None]*nfr)
    for i in xrange(nfr):
        # each realization being nel (87188) elements
        fields[-1][i] = [None]*nel

# read in all realizations
print 'reading ...'
for i,run in enumerate(runs):
    print i,run
    for j,t in enumerate(types):
        fields[j][i][0:nel] = floatload('Outputs/'+run+'/modeled_'+t+'_field.mod')

# save file with one cell from each realization for checking in Excel
if debug:
    print 'writing debugging output for checking'
    fd = open('parameter_representative_values.txt','w')
    fd.write('%s %18s %18s %18s %18s\n'%
             ('rzn',types[0],types[1],types[2],types[3]))
    for i,run in enumerate(runs):
        fd.write('%s %.14e %.14e %.14e %.14e\n' %
                 (run,fields[0][i][-1],fields[1][i][-1],
```

```
                    fields[2][i][159*284],fields[3][i][159*284]))
    fd.close()

# open up files for writing
fh = []
for t in types:
    fh.append(open('modeled_'+ t +'_field.avg','w'))

# transpose fields to allow slicing across realizations, rather than across cells
for j in range(len(types)):
    fields[j] = zip(*(fields[j]))

print 'writing ...'
# do averaging across 100 realizations
for i in xrange(nel):
    if i%10000 == 0:
        print i
    for h,d in zip(fh,fields):
        h.write('%18.11e\n' % pow(10.0,sum(map(log10,d[i]))/nfr) )

for h in fh:
    h.close()
```

A-4.3 Python script boundary_types.py

```
from itertools import chain

nx = 284        # number columns in model grid
ny = 307        # number rows
nel = nx*ny

def intload(filename):
    """Reads file (a 2D integer array) as a list of lists.
    Outer list is rows, inner lists are columns."""
    f = open(filename,'r')
    m = [[int(v) for v in line.rstrip().split()] for line in f]
    f.close()
    return m

def intsave(filename,m):
    """Writes file as a list of lists as a 2D integer array, format '%3i'.
    Outer list is rows, inner lists are columns."""
    f = open(filename,'w')
    for row in m:
        f.write(' '.join(['%2i' % col for col in row]) + '\n')
    f.close()

def floatload(filename):
    """Reads file (a list of real numbers, one number each row)
    into a list of floats."""
    f = open(filename,'r')
    m = [float(line.rstrip()) for line in f]
    f.close()
    return m

def reshapev2m(v):
    """Reshape a vector that was previously reshaped in C-major order
    from a matrix, back into a matrix (here a list of lists)."""
    m = [None]*ny
    for i,(lo,hi) in enumerate(zip(xrange(0,nel-nx+1,nx), xrange(nx,nel+1,nx))):
        m[i] = v[lo:hi]
    return m

##########################################

# read in original MODFLOW IBOUND array (only 0,1, and -1)
ibound = intload('init_bnds_orig.inf')

# read in initial heads
h = reshapev2m(floatload('init_head_orig.mod'))

# discriminate between two types of constant head boundaries
# -1) CH, where value > 1000 (area east of halite margin)
# -2) CH, where value < 1000 (single row/column of cells along domain edge
```

```
50
51 for i,row in enumerate(ibound):
52     for j,val in enumerate(row):
53         # is this constant head and is starting head less than 1000m ?
54         if ibound[i][j] == -1 and h[i][j] < 1000.0:
55             ibound[i][j] = -2
56
57 # save new IBOUND array that allows easy discrimination between types
58 # in python script during PEST optimization runs, and is still handled
59 # the same by MODFLOW since all ibound values < 0 are constant head.
60 intsave('init_bnds.inf',ibound)
```

A-4.4 Python script create_pest_02_input.py

```
prefix = '2008ASER'

#################################################
## pest instruction file reads output from mod2obs
fin = open('meas_head_' + prefix  + '.smp','r')

# each well is a [name,head] pair
wells = [[line.split()[0],line.split()[3]] for line in fin]
fin.close()

fout = open('modeled_head.ins','w')
fout.write('pif @\n')
for i,well in enumerate(wells):
        fout.write("l1 [%s]39:46\n" % well[0])
fout.close()

# exponential surface used to set initial head everywhere
# except east of the halite margins, where the land surface is used.
# initial guesses come from AP-114 Task report
params = [928.0, 8.0, 1.2, 1.0, 1.0, -1.0, 0.5]
pnames = ['a',   'b', 'c', 'd', 'e', 'f', 'exp']

###############################
## pest template file
ftmp = open('surface_par_params.ptf','w')
ftmp.write('ptf @\n')
for n in pnames:
        ftmp.write('@      %s       @\n' % n)
ftmp.close()

###################
## pest parameter file

fpar = open('surface_par_params.par','w')
fpar.write('double  point\n')
for n,p in zip(pnames,params):
    fpar.write('%s  %.2f  1.0  0.0\n' % (n,p))
fpar.close()

####################
## pest control file
f = open('bc_adjust_'+ prefix +'.pst','w')

f.write("""pcf
* control data
restart estimation
%i %i 1 0 1
1 1 double point 1 0 0
5.0 2.0 0.4 0.001 10
```

```
3.0 3.0 1.0E-3
0.1
30 0.001 6 6 0.0001 4
1 1 1
* parameter groups
bc relative 0.005 0.0001 switch 2.0 parabolic
""" % (len(params),len(wells)))

f.write('* parameter data\n')
for n,p in zip(pnames,params):
        if p > 0:
                f.write('%s none relative %.3f %.3f %.3f bc 1.0 0.0 1\n' %
                        (n, p, -2.0*p, 3.0*p))
        else:
                f.write('%s none relative %.3f %.3f %.3f bc 1.0 0.0 1\n' %
                        (n, p, 3.0*p, -2.0*p))

f.write("""* observation groups
ss_head
* observation data
""")

## read in observation weighting group definitions
fin = open('obs_loc_'+prefix+'.dat','r')
location = [line.rstrip().split()[1] for line in fin]
fin.close()

weights = []

for l in location:
    # inside LWB
    if l == '0':
        weights.append(2.5)
    # near LWB
    if l == '1':
        weights.append(1.0)
    # distant to LWB
    if l == '2':
        weights.append(0.4)

for name,head,w in zip(zip(*wells)[0],zip(*wells)[1],weights):
    f.write('%s  %s  %.3f   ss_head\n' % (name,head,w))
f.write("""* model command line
./run_02_model
* model input/output
surface_par_params.ptf surface_par_params.in
modeled_head.ins modeled_head.smp
""")
f.close()
```

A-4.5 Python script surface_02_extrapolate.py

```
from itertools import chain
from math import sqrt

def matload(filename):
    """Reads file (a 2D string array) as a list of lists.
    Outer list is rows, inner lists are columns."""
    f = open(filename,'r')
    m = [line.rstrip().split() for line in f]
    f.close()
    return m

def floatload(filename):
    """Reads file (a list of real numbers, one number each row)
    into a list of floats."""
    f = open(filename,'r')
    m = [float(line.rstrip()) for line in f]
    f.close()
    return m

def reshapem2v(m):
    """Reshapes a rectangular matrix into a vector in same fashion
    as numpy.reshape(), which is C-major order"""
    return list(chain(*m))

def sign(x):
    """ sign function"""
    if x<0:
        return -1
    elif x>0:
        return +1
    else:
        return 0

########################################

# read in modified IBOUND array, with the cells to modify set to -2
ibound = reshapem2v(matload('init_bnds.inf'))

h = floatload('init_head_orig.mod')

# these are relative coordinates, -1 <= x,y < +1
x = floatload('rel_x_coord.dat')
y = floatload('rel_y_coord.dat')

# unpack surface parameters (one per line)
# z = A + B*(y + D*sign(y)*sqrt(abs(y)))+C*(E*x**3 - F*x**2 - x)

finput = open('surface_par_params.in','r')
try:
```

```
    a,b,c,d,e,f,exp = [float(line.rstrip()) for line in finput]
except ValueError:
    # python doesn't like 'D' in 1.2D-4 notation used by PEST sometimes.
    finput.seek(0)
    lines = [line.rstrip() for line in finput]
    for i in range(len(lines)):
        lines[i] = lines[i].replace('D','E')
    a,b,c,d,e,f,exp = [float(line) for line in lines]

finput.close()

# file to output initial/boundary head for MODFLOW model
fout = open('init_head.mod','w')
for i in xrange(len(ibound)):
    if ibound[i] == '-2' or ibound[i] == '1':
        # apply surface to active cells (ibound=1) -> starting guess
        # and non-geologic boundary conditions (ibound=-2) -> constant head
        if y[i] == 0:
            fout.write('%.7e \n' % (a + c*(e*x[i]**3 + f*x[i]**2 - x[i])))
        else:
            fout.write('%.7e \n' % (a + b*(y[i] + d*sign(y[i])*abs(y[i])**exp) +
                                    c*(e*x[i]**3 + f*x[i]**2 - x[i])))
    else:
        # use land surface at constant head east of halite boundary
        # ibound=0 doesn't matter (inactive)
        fout.write('%.7e\n' % h[i])

fout.close()
```

A-4.6 Bash shell script `run_02_model`

```
#!/bin/bash

#echo 'step 1: surface extrapolate'
python surface_02_extrapolate.py

# run modflow
#echo 'step 2: run modflow'
../bin/mf2k/mf2k_1.6.release mf2k_head.nam  >/dev/null

# run mod2obs
#echo 'step 3: extract observations'
../bin/Builds/Linux/mod2obs.exe < mod2obs_head.in  >/dev/null
```

A-4.7 Python script `head_bin2ascii.py`

```
import struct

class FortranFile(file):
    """ modified from May 2007 Enthought-dev mailing list
    post by Neil Martinsen-Burrell"""

    def __init__(self,fname, mode='r', buf=0):
        file.__init__(self, fname, mode, buf)
        self.ENDIAN = '<'  # little endian
        self.di = 4  # default integer (could be 8 on 64-bit)

    def readReals(self, prec='f'):
        """Read in an array of reals (default single precision)
        with error checking"""
        # read header (length of record)
        l = struct.unpack(self.ENDIAN+'i',self.read(self.di))[0]
        data_str = self.read(l)
        len_real = struct.calcsize(prec)
        if l % len_real != 0:
            raise IOError('Error reading array of reals from data file')
        num = l/len_real
        reals = struct.unpack(self.ENDIAN+str(num)+prec,data_str)
        # check footer
        if struct.unpack(self.ENDIAN+'i',self.read(self.di))[0] != l:
            raise IOError('Error reading array of reals from data file')
        return list(reals)

    def readInts(self):
        """Read in an array of integers with error checking"""
        l = struct.unpack('i',self.read(self.di))[0]
        data_str = self.read(l)
        len_int = struct.calcsize('i')
        if l % len_int != 0:
            raise IOError('Error reading array of integers from data file')
        num = l/len_int
        ints = struct.unpack(str(num)+'i',data_str)
        if struct.unpack(self.ENDIAN+'i',self.read(self.di))[0] != l:
            raise IOError('Error reading array of integers from data file')
        return list(ints)

    def readRecord(self):
        """Read a single fortran record (potentially mixed reals and ints)"""
        dat = self.read(self.di)
        if len(dat) == 0:
            raise IOError('Empy record header')
        l = struct.unpack(self.ENDIAN+'i',dat)[0]
        data_str = self.read(l)
        if len(data_str) != l:
            raise IOError('Didn''t read enough data')
```

```
50          check = self.read(self.di)
51          if len(check) != 4:
52              raise IOError('Didn''t read enough data')
53          if struct.unpack(self.ENDIAN+'i',check)[0] != 1:
54              raise IOError('Error reading record from data file')
55          return data_str
56
57  def reshapev2m(v,nx,ny):
58      """Reshape a vector that was previously reshaped in C-major order
59      from a matrix, back into a C-major order matrix (here a list of lists)."""
60      m = [None]*ny
61      n = nx*ny
62      for i,(lo,hi) in enumerate(zip(xrange(0,n-nx+1,nx), xrange(nx,n+1,nx))):
63          m[i] = v[lo:hi]
64      return m
65
66  def floatmatsave(filehandle,m):
67      """Writes array to open filehandle, format '568%e12.5'.
68      Outer list is rows, inner lists are columns."""
69
70      for row in m:
71          f.write(''.join([' %12.5e' % col for col in row]) + '\n')
72
73  # open file and set endian-ness
74  ff = FortranFile('modeled_head.bin')
75
76  # currently this assumes a single-layer MODFLOW model
77  # (or at least only one layer of output)
78
79  # format of MODFLOW header in binary layer array
80  fmt = '<2i2f16s3i'
81  # little endian, 2 integers, 2 floats,
82  #    16-character string (4 element array of 4-byte strings), 3 integers
83
84  while True:
85      try:
86          # read in header
87          h = ff.readRecord()
88
89      except IOError:
90          # exit while loop
91          break
92
93      else:
94          # unpack header
95          kstp,kper,pertim,totim,text,ncol,nrow,ilay = struct.unpack(fmt,h)
96
97          # print status/confirmation to terminal
98          print kstp,kper,pertim,totim,text,ncol,nrow,ilay
99
100         h = ff.readReals()
```

```
101
102 ff.close()
103
104 xmin, xmax = (601700.0,630000.0)
105 ymin, ymax = (3566500.0,3597100.0)
106 hmin = min(h)
107 hmax = max(h)
108
109 # write output in Surfer ASCII grid format
110 f = open('modeled_head_asciihed.grd','w')
111 f.write("""DSAA
112 %i %i
113 %.1f %.1f
114 %.1f %.1f
115 %.8e %.8e
116 """ %(ncol,nrow,xmin,xmax,ymin,ymax,hmin,hmax) )
117 hmat = reshapev2m(h,ncol,nrow)
118
119 # MODFLOW starts data in upper-left corner
120 # Surfer expects data starting in lower-left corner
121 # flip array in row direction
122
123 floatmatsave(f,hmat[::-1])
124 f.close()
```

A-4.8 Python script `merge_observed_modeled_heads.py`

```
fobs = open('meas_head_2008ASER.smp','r')
fmod = open('modeled_head.smp','r')
fwgt = open('obs_loc_2008ASER.dat')
fout = open('both_heads.smp','w')

fout.write('\t'.join(['#NAME','OBSERVED','MODELED','WEIGHT'])+'\n')

for obs,mod,w in zip(fobs,fmod,fwgt):
    fout.write('\t'.join([obs.split()[0],obs.split()[3],mod.split()[3],
                          w.rstrip().split()[1]])+'\n')

fobs.close()
fmod.close()
fwgt.close()
fout.close()
```

A-4.9 Python script `convert_dtrkmf_output_for_surfer.py`

```
1
2  # grid origin for dtrkmf cell -> x,y conversion
3  x0 = 601700.0
4  y0 = 3597100.0
5
6  dx = 100.0
7  dy = 100.0
8
9  fout = open('dtrk_output.bln','w')
10
11 # read in all results for saving particle tracks
12 fin = open('dtrk.out','r')
13 results = [l.split() for l in fin.readlines()[1:]]
14 fin.close()
15
16 npts = len(results)
17
18 # write Surfer blanking file header
19 fout.write('%i,1\n' % npts)
20
21 # write x,y location and time
22 for pt in results:
23     x = float(pt[1])*dx + x0
24     y = y0 - float(pt[2])*dy
25     t = float(pt[0])/7.75*4.0  # convert 7.75m to 4m Cuelbra thickness
26     fout.write('%.1f,%.1f,%.8e\n' % (x,y,t))
27
28 fout.close()
```