Repeating the previous Surrogate building sage (Chapter 7) but with addition of the anode consumption related data as an extra input variable to the surrogate model.

Note: The anode consumption related varible is not one of p-value or conductivity related parameter so will be dealt in a different way.

Also this experiment uses the experimental results from The Parameter\_CAxx\_CBxx.mlx files (0,5,10,15). So make sure to undertake those experiments in advance which are similar to 2 parameter estimation using surrogate (Chapter 7)

```
response_data_type = {'voltage', 'normal current density', 'Z_ELECTRIC_FIELD','anode related'
%response_data_type = {'voltage', 'Z_CURRENT_DENSITY', 'anode related'};
%calibration_data_type = {'voltage'};
%calibration_data_type = {'voltage', 'current density', 'Z_ELECTRIC_FIELD'};
metric = 'nmsq';
IPs_IDs = 32820:32868;
IPs IDs1 = IPs IDs(1:2:end);
IDs_current_density = [14390, 7400, 4060, 16000, 19860, 23802, 30002, 23822,21212, 8437];
%IDs_current_density =[27791
                                         27813
                                                        27839
                                                                      278591;
%IDs_current_density = IPs_IDs1;
anode_related_IDs = 1:25:266;
IDs = {py.list(IPs_IDs1), py.list(IDs_current_density),py.list(IPs_IDs1), py.list(anode_relat
IDs_mat_arr = {IPs_IDs1, IDs_current_density.', IPs_IDs1,anode_related_IDs.'};
%IDs_types = {'Internal Points','Mesh Points', 'Internal Points','CONSUMPTION_FACTOR','ANODE_
%IDs_types = {'Internal Points','Mesh Points', 'Internal Points','ANODE_CURRENT'};
IDs_types = {'Internal Points','Mesh Points', 'Internal Points', 'MASS_LOSS_RATE'};
%IDs_types = {'Internal Points', 'Internal Points', 'MASS_LOSS_RATE'};
%py.list(1:2)
parameters_base = {'CA', 'CB'};
```

The average anode consumption for different year time is something the CP model should predict, but this approach deals it as an Inverse problem.

The average anode consumption rate is added for each of the year during data sampling.

```
years = [0, 5,10,15];
all_anode_IDs = 1:266;
```

```
root_folder1 = 'D:\DOE_nd_data_generation\TIme_step\readyToTimeStepUsingV10_B';
average cons factor = average anode CONSUMPTION FACTOR(root folder1, years, all anode IDs);
average_cons_factor
average\_cons\_factor = 1 \times 4
           0.0754
                   0.2350
                             0.4450
anode IDs = 1:25:255;
res_folder = 'D:\DOE_nd_data_generation\TIme_step\readyToTimeStepUsingV10_B\year_10\Calibrati
anode_cons_rate = get_anode_data(res_folder, anode_IDs, 'CONSUMPTION_FACTOR');
%anode_cons_rate
anode_cons_rate = 11 \times 2
           0.2117
   1.0000
 226.0000
           0.2365
 101.0000
           0.2522
 201.0000
           0.2313
  76.0000
           0.2707
 176.0000
           0.2229
  51.0000
           0.2345
 151.0000
           0.2417
  26.0000
           0.2155
 251.0000
           0.2372
%DOE experiment for 2 varaibles using Central Composite Design
Central composite points = ccdesign(2, 'type', 'inscribed', 'center' , 1);
DOE_range1 = [0.005, 0.2; 0.005, 0.15];
DOE_range10 = [0.15, 0.35; 0.100, 0.25];
DOE_range15 = [0.15, 0.275; 0.1100, 0.19];
DOE sample points= cell(1, length(years)); snapshots years = DOE sample points;
DOE_sample_points{1} = reverse_normalization(Central_composite_points, DOE_range1);
DOE_sample_points{2} = reverse_normalization(Central_composite_points, DOE_range1);
DOE_sample_points{3} = reverse_normalization(Central_composite_points, DOE_range10);
DOE sample points{4} = reverse normalization(Central composite points, DOE range15);
for i = 1:length(years)
    source parameters = {'BARE', 'BARE'};
    if years(i) <10</pre>
        parameters = strcat(parameters_base, '0', cellstr(string(years(i))));
```

else

```
parameters = strcat(parameters_base, cellstr(string(years(i))));
end

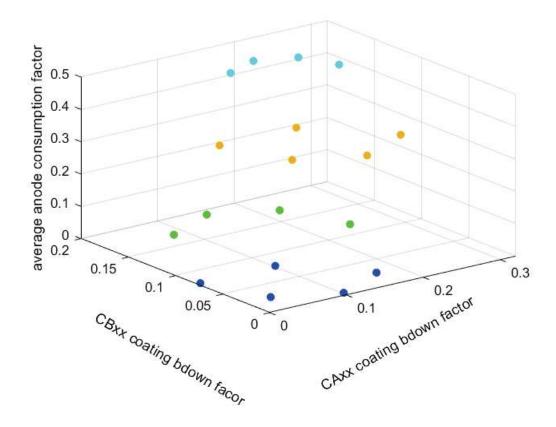
parameters_np_array1 = convert_arr_to_python_2d_list(DOE_sample_points{i});
simulation_seed_folder = fullfile(root_folder1,strcat('year_',string(years(i))), 'Initial_collection_dir = fullfile(root_folder1,strcat('year_',string(years(i))), 'Simulation_resusinapshots_py = py.BEASY_IN_OUT2.snapshots_for_given_parameters_and_IDs(py.list(source_parameters_i));
end

end
```

Data samples collected at different time-span, now will be used in surrogate-building for the three variables case (PhD Thesis section 8.5.3)

```
figure;
%scatter3( DOE sample points3(:,1), DOE sample points3(:,2),DOE sample points3(:,3),'filled
%scatter3( DOE sample points collected(:,1), DOE sample points collected(:,2),DOE sample poi
sampl idx = 1;
scatter3( DOE sample points collected cells{sampl idx}(:,1),
    DOE_sample_points_collected_cells{sampl_idx}(:,2),DOE_sample_points_collected_cells{sampl}
hold on;
sampl idx = 2;
scatter3( DOE_sample_points_collected_cells{sampl_idx}(:,1),
    DOE_sample_points_collected_cells{sampl_idx}(:,2),DOE_sample_points_collected_cells{sample}
sampl idx = 3:
scatter3( DOE sample points collected cells{sampl idx}(:,1),
    DOE_sample_points_collected_cells{sampl_idx}(:,2),DOE_sample_points_collected_cells{sampl
sampl idx = 4;
scatter3( DOE sample points collected cells{sampl idx}(:,1), ...
    DOE_sample_points_collected_cells{sampl_idx}(:,2),DOE_sample_points_collected_cells{sampl}
```

```
xlabel('CAxx coating bdown factor', 'Rotation', 30);
ylabel('CBxx coating bdown facor', 'Rotation', -30);
zlabel('average anode consumption factor');
```



```
%csvwrite('Snapshots_3d.csv', snapshots_collected )
%csvwrite('DOE_sample_points_collected_3d.csv', DOE_sample_points_collected )
```

```
%csvwrite('DOE_sample_points_collected_3d.csv', DOE_sample_points_collected )
```

To build the 3 dimensional polynomial fit surrogate, this tool relies upon the MATLAB toolbox 'polyfitn'. The toolbox polyfitn need to be downloaded into the system and linked if doesnot exist within the MATLAB.

```
surrogates_3d = response_surface(DOE_sample_points_collected, snapshots_collected,2);
```

## surrogates\_3d

surrogates\_3d = 1×71 cell

1	1×1 struct							
	1	2	3	4	5	6	7	8

```
surrogates_3d{1}
```

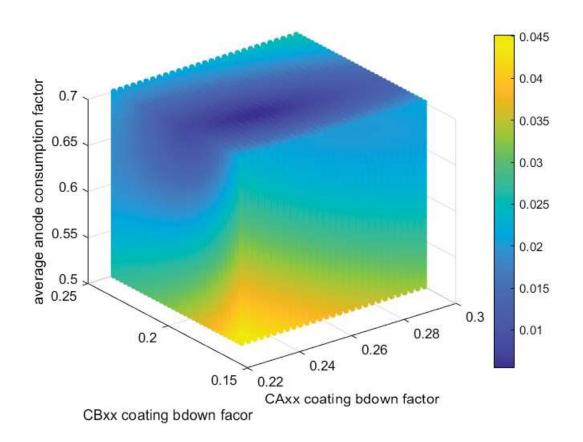
ans = struct with fields:

ModelTerms: [10×3 double]

```
Coefficients: [-629.0825 -202.6927 -13.1925 526.6198 662.3681 -464.8554 23.8937 93.4226 28.4219 -1.0791e+03]
   ParameterVar: [2.9296e+04 1.9555e+05 5.4551e+04 2.7400e+03 4.6670e+05 2.8728e+05 1.1682e+04 1.6283e+04 3.5925e
   ParameterStd: [171.1604 442.2061 233.5625 52.3451 683.1581 535.9806 108.0826 127.6067 59.9375 4.9720]
            p: [0.0063 0.6589 0.9563 8.1140e-06 0.3607 0.4110 0.8306 0.4850 0.6480 2.2735e-16]
            R2: 0.9959
     AdjustedR2: 0.9912
          RMSE: 1.8967
       VarNames: {'X1' 'X2' 'X3'}
calib_dir = fullfile(root_folder1, 'year_20', 'Calibration_data');
%meas dir1 = 'C:\Users\msapkota\EXPERIMENT\DOE nd data generation\Model updated nonlinear\Par
files_name = 'BU_TimeStepped_01_20';
%meas_dict = py.BEASY_IN_OUT1.get_output_data_for_IDs_from_simulation_folder(meas_dir, files
calib data IDs = {IPs IDs1, IDs current density};
calib data type = response data type(1:2);
%meas_data = convert_pydict2data(meas_dict,0);
calib_data_file_err_inc = strcat('data_with_error_',strjoin(response_data_type, '_'), '.xlsx'
%all_position_dict = py.BEASY_IN_OUT2.get_output_data_for_IDs_from_simulation_folder(calib_di
if ~isfile(fullfile(calib_dir, calib_data_file_err_inc))
    all_position_dict = py.BEASY_IN_OUT2.get_output_data_for_IDs_from_simulation_folder(calib
        files name, py.list(calib data type), py.list({py.list(IPs IDs), py.list(IDs current
    all_position_data = convert_pydict2data(all_position_dict,0);
    introduce_error_and_write_file( {IPs_IDs.', IDs_current_density.'},all_position_data, ...
        calib_dir, calib_data_file_err_inc,1);
end
Warning: Added specified worksheet.
%model out = output from surrogates([2.0, 3.0], surrogates, [17,6]);
calib data inc error = data from tables(fullfile(calib dir, calib data file err inc), calib d
%if we know one of the value we can fix it
figure;
ax = gca;
DOE range20 = [0.23 \ 0.3; \ 0.17 \ 0.25; \ 0.5, 0.7];
[plot data, min value, min out pos20] = plot objective with surrogates(ax, ...
    DOE_range20, surrogates_3d, calib_data_inc_error, 'nmsq', calib_data_type, ...
    cellfun('length',IDs_mat_arr),[1,0.5,0,0], [0.0025,0.0025, 0.0025]);
```

```
sol_output_from_surrogate = output_from_surrogates(min_out_pos20, surrogates_3d,cellfun('leng
    IDs_mat_arr));

xlabel('CAxx coating bdown factor');
ylabel('CBxx coating bdown facor');
zlabel('average anode consumption factor');
```



## min\_out\_pos20

```
min_out_pos20 = 1×3
0.2625 0.2025 0.6825
```

```
%min_out_pos20 = [0.2575    0.2000    0.6925]
testing_par_value = min_out_pos20;

sol_output_from_surrogate = output_from_surrogates(testing_par_value, ...
    surrogates_3d,cellfun('length',IDs_mat_arr));

testing_par_value = min_out_pos20(1:2);
root_folder = 'D:\DOE_nd_data_generation\TIme_step\readyToTimeStepUsingV10_B\year_20';
files_name = 'BU_TimeStepped_01_20';
```

```
parameters= {'CA20','CB20'};
simulation seed folder = fullfile(root folder, 'Initial files');
solution_folder = '';
for i = 1:length(parameters)
                        strcat(solution_folder, parameters{i},'_', num2str(testing_par_value(
    solution folder =
        '%.4f'));
    if i~=length(parameters)
        solution_folder = strcat(solution_folder, '_');
end
solution colection dir = fullfile(root folder, 'Solution results');
solution_dir = fullfile(solution_colection_dir, solution_folder);
if ~isfolder(solution dir)
    solution_dict = py.BEASY_IN_OUT2.get_response_data_for_IDs_and_input_parameters(py.list(s
        py.list(parameters), testing_par_value, simulation_seed_folder, solution colection di
    solution_data = convert_pydict2data(solution_dict,1);
else
    solution_dict = py.BEASY_IN_OUT2.get_output_data_for_IDs_from_simulation_folder(solution_
        files_name, py.list(response_data_type), py.list(IDs), py.list(IDs_types));
    solution data = convert pydict2data(solution dict,0);
end
('running_simulation in:', 'D:\\DOE_nd_data_generation\\TIme_step\\readyToTimeStepUsingV10_B\\year_20\\Solution_re
```

```
figure;
%difference_in_bar_chart(ax,simulation_data{1}(1:4:end,:), out_frm_nnet{1}(1:4:end,:),{'simua
%legend cell = strcat('Data from Simulation on year', strsplit(num2str(time period)));
legend cell = {'calibration data', 'solution output from surrogate',
    'simulation output from solution model'};
data_ID = 1;
response in bar chart(ax, solution data{data ID}(1:2:end,:), ...
    {calib_data_inc_error{data_ID}(1:2:end,:), sol_output_from_surrogate{data_ID}(1:2:end,:),
    solution data{data ID}(1:2:end,:)}, legend cell);
xlabel('Internal Points IDs');
%xlabel('Mesh Points IDs');
%xlabel('Data Positional IDs', 'Rotation', 30);
%ylabel('consumption rate (kg/yr)')
%ylabel('Z electric field (micro-V/m)');
%ylabel('Normal current density (mAmp/Sq.m)');
ylim([-1000 -950]);
```

## Internal Points IDs 3282@28243282832832836 3284B284A284B28523285B286B286A32868 -950 Potential difference Ag/Agcl/Sea-water (mV) -955 -960 -965 -970 calibration data -975 solution output from surrogate simulation output from solution model -980 -985 -990 -995 -1000

```
%ylim([15 22]);
%ylabel('anode Current');
%ylim([1200 2000]);
```

```
function average_CONSUMPTION_FACTOR = average_anode_CONSUMPTION_FACTOR(root_folder_address,
    average_CONSUMPTION_FACTOR = zeros(1, length(years_time));
    for i = 1:length(years_time)
        year_t = years_time(i);
    response_folder = fullfile(root_folder_address, strcat('year_', string(year_t)), 'Calibra
        simulation_files = dir(response_folder);
        files_name = simulation_files(end-3).name;
        files_name = strsplit(files_name, '.');
        files_name = files_name{1};

        %anode_file = fullfile(response_folder, strcat(files_name, '.cp_anode_decay'));
        initial_anode_data = py.BEASY_IN_OUT2.get_output_data_for_IDs_from_simulation_folder(response_initial_anode_data = initial_anode_data{1};
        initial_anode_data = initial_anode_data{1};
}
```

```
average CONSUMPTION FACTOR(i) = mean(initial anode data(:,2));
end
end
function anode_data = get_anode_data(response_folder, anode_IDs, data_type)
    simulation_files = dir(response_folder);
    files_name = simulation_files(end-2).name;
    files_name = strsplit(files_name, '.');
    files_name = files_name{1};
    anode_file = fullfile(response_folder,strcat(files_name, '.cp_anode_decay'));
    anode_data = py.BEASY_IN_OUT2.get_output_data_for_IDs_from_simulation_folder(response_fol
    anode_data = convert_pydict2data(anode_data,0);
    anode_data = anode_data{1};
end
function de_normaised_data = reverse_normalization(normalised_data, value_ranges)
de_normaised_data = zeros(size(normalised_data));
for i = 1:size(normalised_data, 2)
    de_normaised_data(:,i) = value_ranges(i,1)+ diff(value_ranges(i,:))/2 * (normalised_data(
end
end
```