
유체 운동의 실시간 시뮬레이터 개발을 위한 적합직교분해 및 인공신경망 기반의 차수축소모델 구성 프레임워크 개발

2022 9th OpenFOAM Korea Users' Community Conference

2022.9.22.

(주) 넥스트폼 이웅현



개요

■ 연구 목표

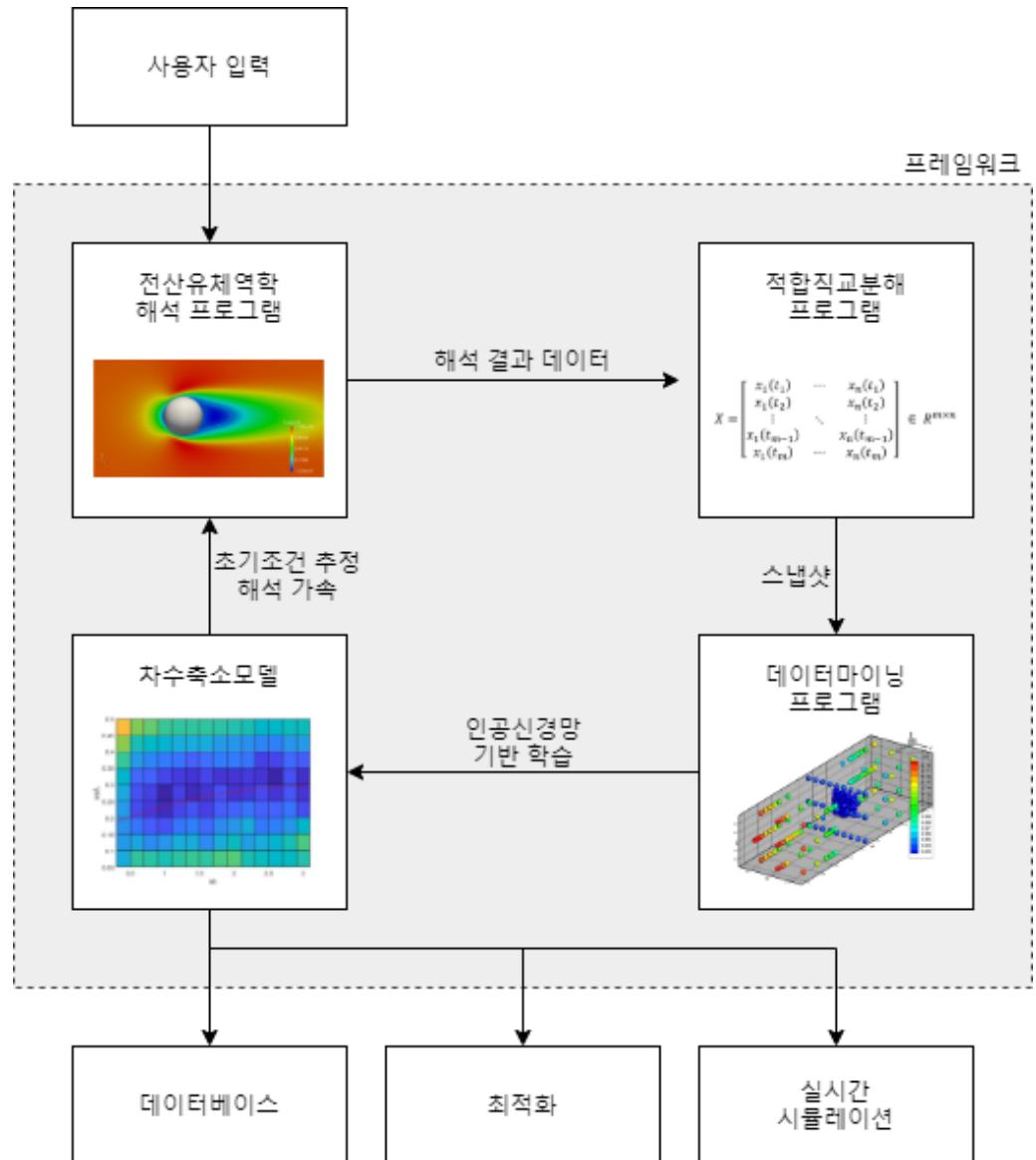
- 반복해석 소요 시간 저감
- 기존에 있는 구성 요소 프로그램을 결합, 자동화된 프레임워크 개발

■ 키워드

- 전산유체역학 (CFD)
- 적합직교분해 (POD)
- 데이터마이닝
- 차수축소모델 (ROM)
- 인공신경망 (ANN)
- 데이터베이스
- 최적화
- 실시간 시뮬레이션

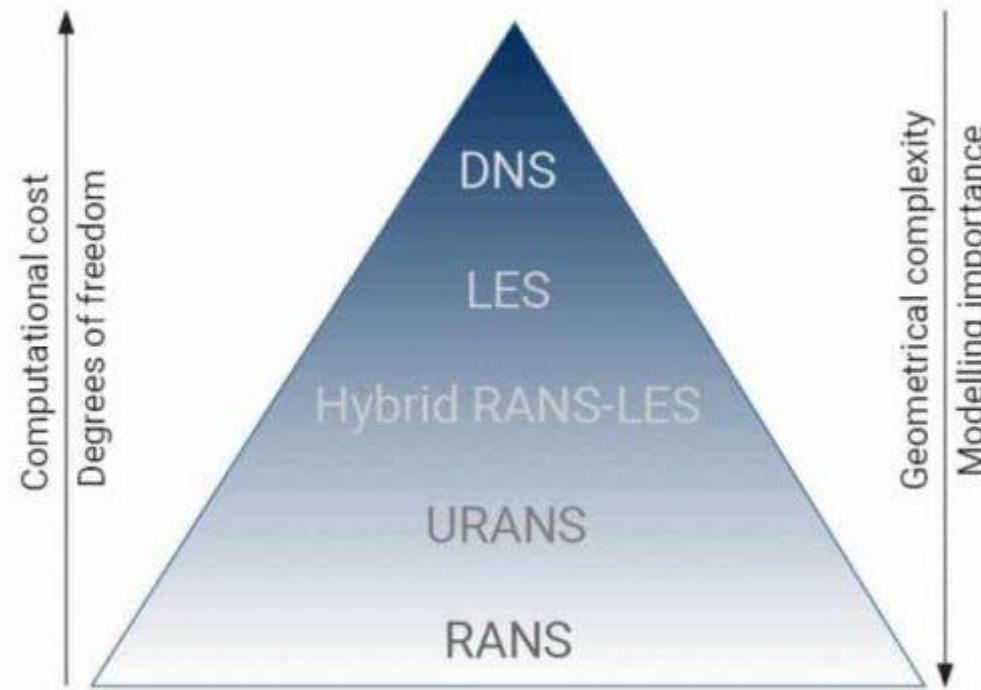
■ 목차

- (1) 적합직교분해
- (2) DAKOTA toolkit
- (3) 프레임워크 개발



적합직교분해 (POD)

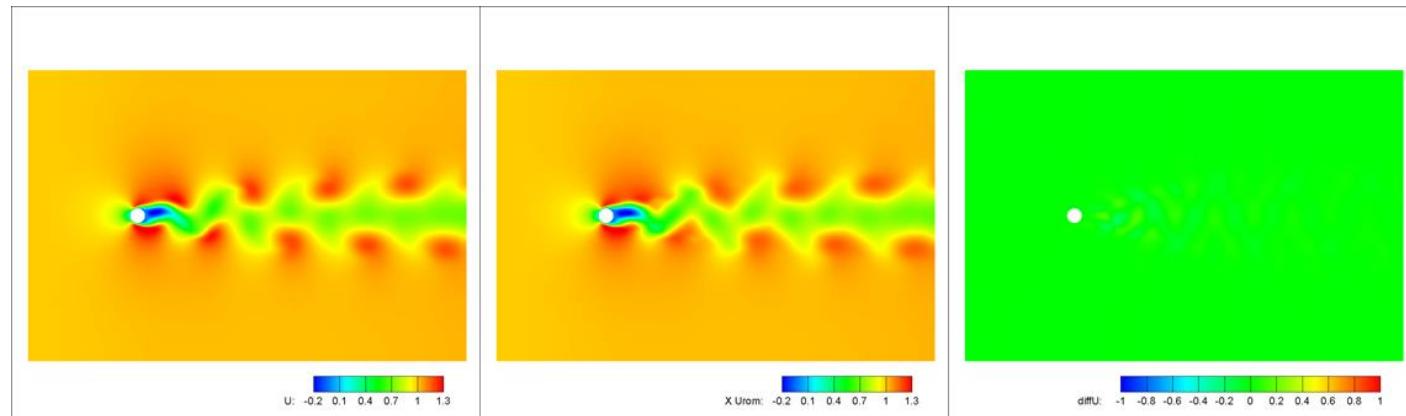
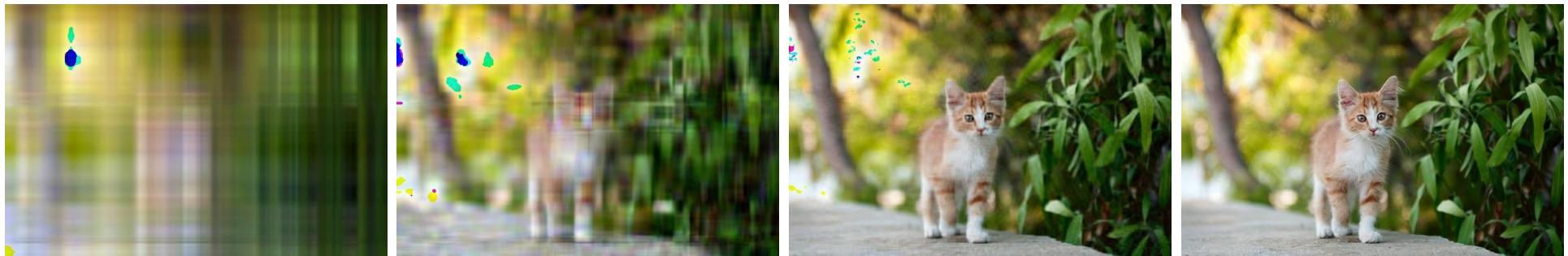
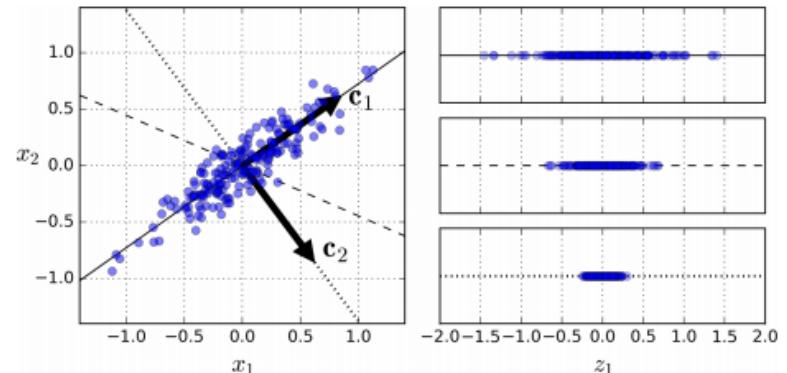
- 기 존재하는 전산유체역학 해석 결과 사용, 모드 추출
- 신규 조건에 대한 해석 결과를 예측



POD, Flow angle, Panel method, etc.

적합직교분해 (POD)

- 주성분 분석 (PCA) 기법 기반
 - 특이치 분해 (SVD)
 - 이미지 압축 등 광범위한 분야에 사용
 - CFD 해석 결과에 적용 가능

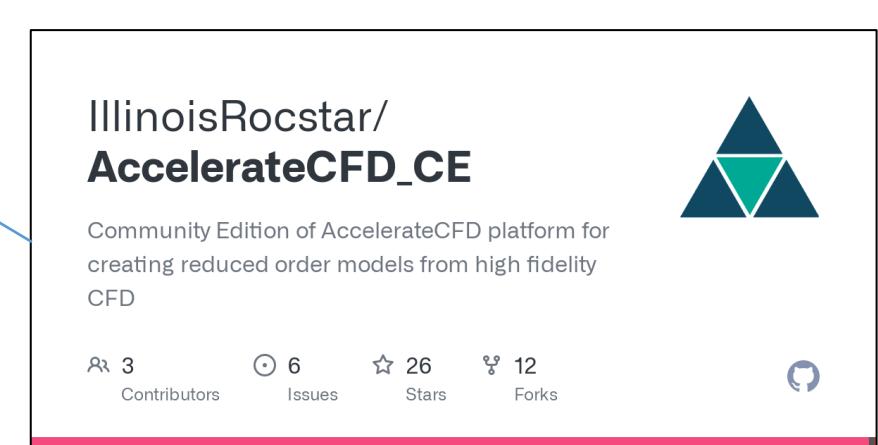
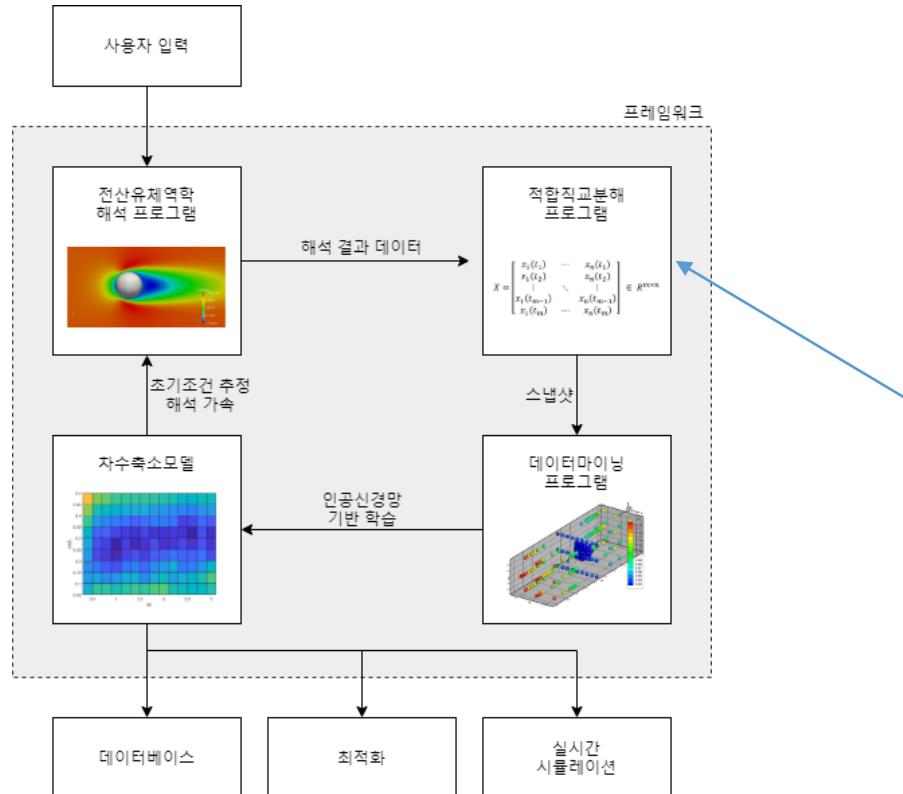


적합직교분해 (POD)

- 고전 적합직교분해
 - 연속적인 시간에 대해 적합직교분해 수행
 - 1967, Lumley, J. L., “The structure of inhomogeneous turbulent flows”, Atmosphere and its Influence on Radio Wave Propagation. pp. 166-178, Moskow: Nauka.
- 스냅샷 적합직교분해 *
 - 특정 시점들에 대한 해석 결과만으로 적합직교분해 수행
 - 1987, Sirovich, L., “Turbulence and the dynamics of coherent structures. Part I: coherent structures”, Applied mathematics, vol.45, no.3, pp.561-571.

적합직교분해 (POD)

- 프레임워크 구성 요소 프로그램 선정 (AccelerateCFD)
 - 공개 소스 (GNU GPL v3), 절차 지향적 코드
 - OpenFOAM utility 형태
 - 계산 결과 입출력 및 추가 기능 개발 용이



적합직교분해 (POD)

■ 절차

ir/AccelerateCFD_CE	
☰	README.md
There are five modules to this software.	
podBasisCalc	<ul style="list-style-type: none">This application calculates POD basis for velocities in CFD case directory and gives information about flow energy contained in each POD basis. It solves eigen value problem to identify most important flow characteristics from the flow.
podPrecompute	<ul style="list-style-type: none">This application calculates gradient and tensor terms as well as some tensor innerproducts for velocity and POD basis vectors. All the pre-computed data which is essential for computing the reduced order model (ROM) is written in the case directory. The file "prevVals.csv" contains time varying coefficients obtained using proper orthogonal decomposition.
podROM	<ul style="list-style-type: none">This application uses all the data written out from podPrecompute application and calculates the time varying coefficients for spatial POD basis to construct reduced order model. It writes values of time varying coefficients in the case directory which are used for reconstructing the velocity field.
podReconstruct	<ul style="list-style-type: none">This application reads in the values of time varying coefficients and reconstructs velocities. These reconstructed velocities are automatically written into their respective time directories of CFD case for ease of visualization.
podPostProcess	<ul style="list-style-type: none">This application allows users to obtain additional information from reduced order as well as full order models for comparison and reference purposes. Right now this utility supports calculation of time varying coefficients from full order model that can serve as a reference to reduced order time coefficients calculated using podROM utility. This utility operates based on command line arguments. All available arguments are explained later in this guide.

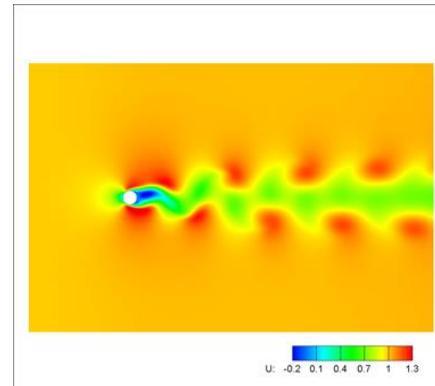
02_SVD	<ul style="list-style-type: none">공분산행렬(Covariance Matrix) 생성Eigenvalue & Eigenvectors 추출 ✓ Left Eigenvector = POD mode (Basis function)	$R = X^T X$ $X = \begin{bmatrix} x_1(t_1) & \cdots & x_n(t_1) \\ \vdots & \ddots & \vdots \\ x_1(t_m) & \cdots & x_n(t_m) \end{bmatrix} \in R^{m \times n}$ $R\beta_m = \lambda_m \beta_m$
03_POD	<ul style="list-style-type: none">Calculate the basis function(ϕ_m)Normalize basis function($\bar{\phi}_m$)Calculate POD's expansion coefficient(a_m)	$\phi_m = X^T \cdot \beta_m$ $\bar{\phi}_m = \phi_m / \sqrt{\phi^2}$ $a_m = X^T \cdot \bar{\phi}_m$
04_Interpolation_Coefficient	<ul style="list-style-type: none">POD's expansion coefficient(a_m)을 Design space의 변수에 따라 보간 또는 회귀모델 생성	<ul style="list-style-type: none">Linear interpolation using SciPy module (1D)또는 RBF 모델 사용(2D, <u>Isight</u> 사용)
05_Reconstruction	<ul style="list-style-type: none">보간/회귀모델로 생성된 POD's expansion coefficient과 basis function을 이용하여 Flow filed를 재생성	$X^{new} = a_m \cdot \bar{\phi}_m$

적합직교분해 (POD)

■ 절차

- (1) 해석 결과 (snapshot) 확보

```
Documents/code/211215_pod_test$ ll
...
0 -> ./211214_stl_sonicFoam/motorBike_5/500/
10 -> ./211214_stl_sonicFoam/motorBike/500/
5 -> ./211214_stl_sonicFoam/motorBike_5/500/
8 -> ./211214_stl_sonicFoam/motorBike_8/500/
constant/
system/
...
Documents/code/211215_pod_test$
```



- 입력 parameter 조건 별 해석 결과
 - AOA, BETA, Mach #, geometric parameter 등 종류 무관
 - 많을수록 정확도 ↑
- AccelerateCFD 구동을 위해 단일 OpenFOAM 케이스 폴더로 취합
- N개 격자 * M개 시점 해석 결과가 취합된 $N \times M$ 행렬 Y

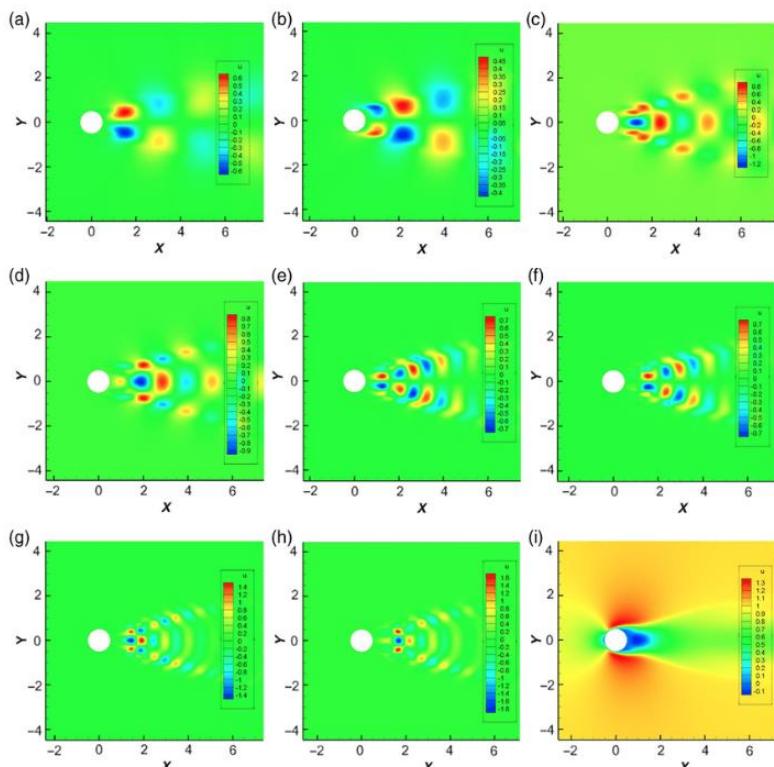
$$\vec{y}_j = \begin{bmatrix} y_{1j} \\ y_{2j} \\ \dots \\ y_{Nj} \end{bmatrix} \quad Y = \begin{bmatrix} \vec{y}_1 & \vec{y}_2 & \dots & \vec{y}_M \end{bmatrix}$$

적합직교분해 (POD)

■ 절차

■ (2) 모드 (basis) 추출

- 해석 결과의 기저 벡터
 - 선형결합으로 유동장 재현



POD modes of the oscillating cylinder: (a) POD mode #1, (b) POD mode #2, (c) POD mode #3, (d) POD mode #4, (e) POD mode #5, (f) POD mode #6, (g) POD mode #7, (h) POD mode #8 and (i) average field.

02 SVD

- 공분산 행렬(Covariance Matrix) 생성
 - Eigenvalue & Eigenvectors 추출
 - ✓ Left Eigenvector = POD mode
(Basis function)

$$R = X^T X$$

$$R = Y^T Y \quad R\vec{v}_i = \sigma_i^2 \vec{v}_i, \quad \vec{u}_i = \frac{\vec{Y}\vec{v}_i}{\sigma_i}$$

```

201     forAll(timeDirs, timei)
202     {
203         n = 0;
204         forAll(timeDirs, timej)
205         {
206             Cmn(m, n) = 0.0;
207             // applying symmetry
208             if (n < m)
209             {
210                 Cmn(m, n) = Cmn(n, m);
211                 n++;
212                 continue;
213             }
214
215             volScalarField U1dotU2(generateCustomField(runTime, mesh, "U1dotU2"),
216                                     (vels[timei]&vels[timej])*cellVolume);
217
218             Cmn(m,n) = Cmn(m,n) + gSum(U1dotU2);
219             n++;
220         }
221         m++;
222     }
223
224     // Normalization of correlation matrix by dividing with total number of
225     Cmn = Cmn/nDim;
226
227     // Self Adjoint Eigen Solver is used here to solve for eigenvalue problem
228     Info<< "Solving eigenvalue problem" << nl;
229
230     Eigen::SelfAdjointEigenSolver<Eigen::MatrixXcd> es(Cmn);

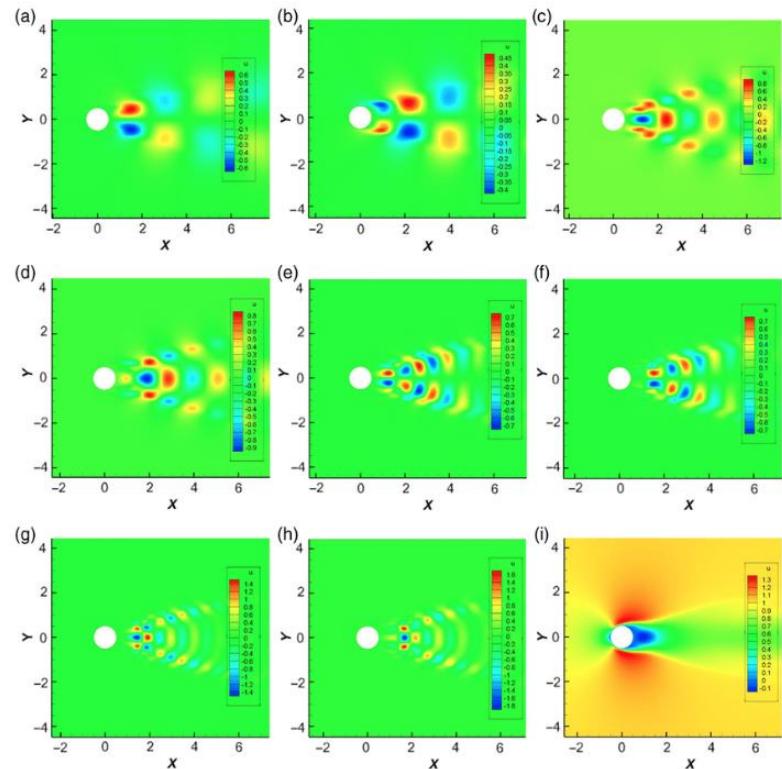
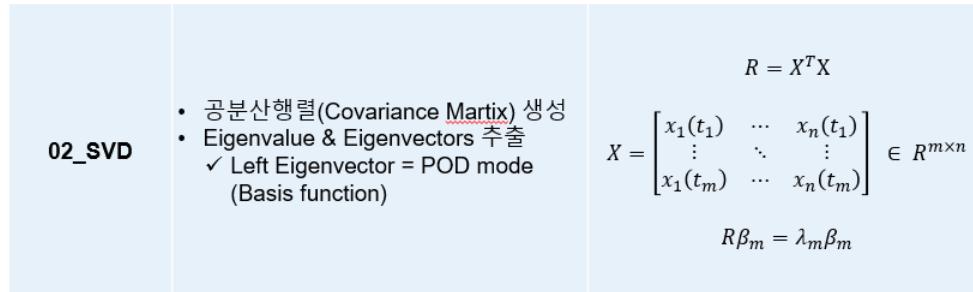
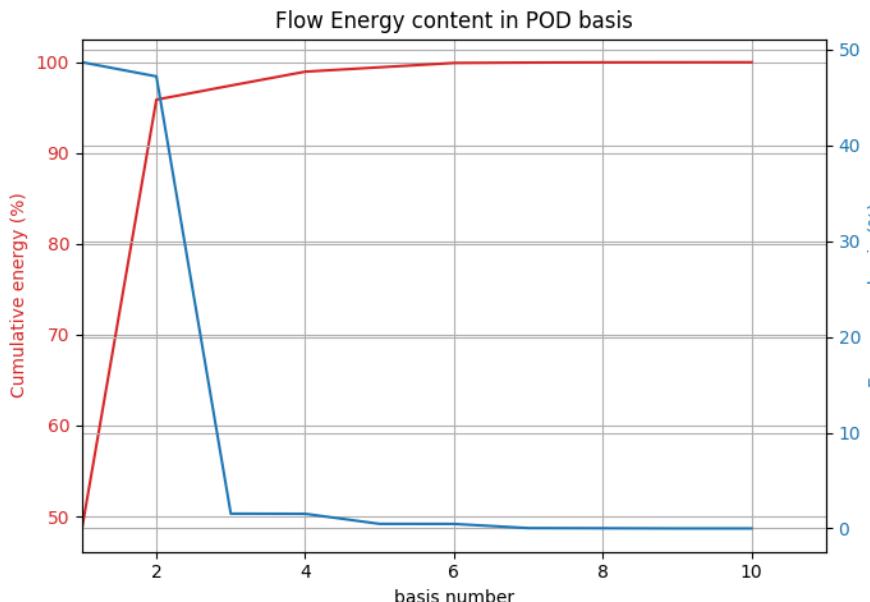
```

적합직교분해 (POD)

■ 절차

■ (2) 모드 (basis) 추출

- 해석 결과의 기저 벡터
- 선형결합으로 유동장 재현
- 5~10개 모드에 99.9% 에너지 분포
- 전체 해석 결과 요약/예측 가능

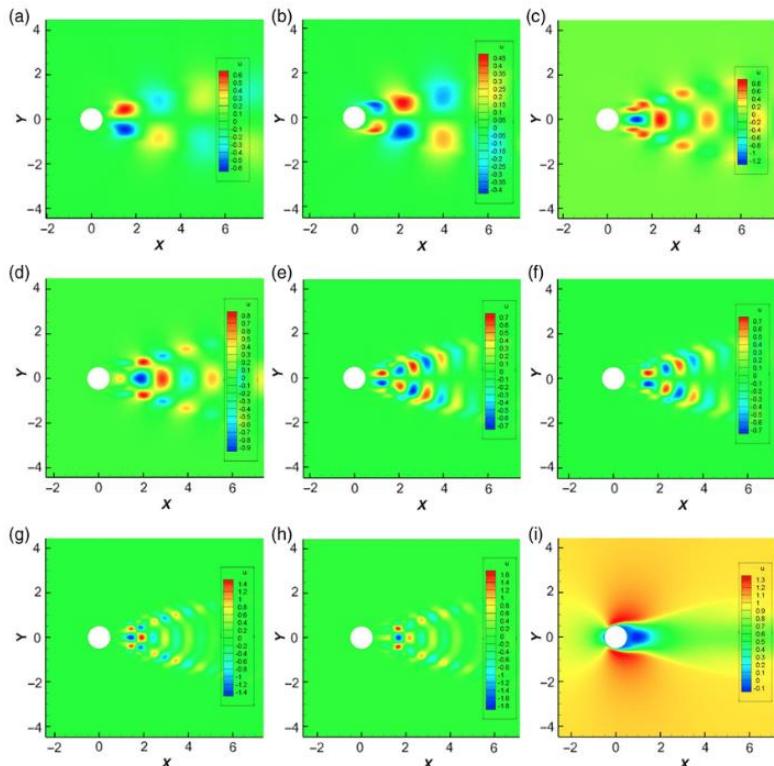


POD modes of the oscillating cylinder: (a) POD mode #1, (b) POD mode #2, (c) POD mode #3, (d) POD mode #4, (e) POD mode #5, (f) POD mode #6, (g) POD mode #7, (h) POD mode #8 and (i) average field.

적합직교분해 (POD)

■ 절차

- (3) expansion coefficient 계산
 - 각 snapshot에 포함된 basis의 비중



POD modes of the oscillating cylinder: (a) POD mode #1, (b) POD mode #2, (c) POD mode #3, (d) POD mode #4, (e) POD mode #5, (f) POD mode #6, (g) POD mode #7, (h) POD mode #8 and (i) average field.

03_POD

- Calculate the basis function(ϕ_m)
- Normalize basis function($\bar{\phi}_m$)
- Calculate POD's expansion coefficient(a_m)

$$\phi_m = X^T \cdot \beta_m$$

$$\bar{\phi}_m = \phi_m / \sqrt{\phi^2}$$

$$a_m = X^T \cdot \bar{\phi}_m$$

```

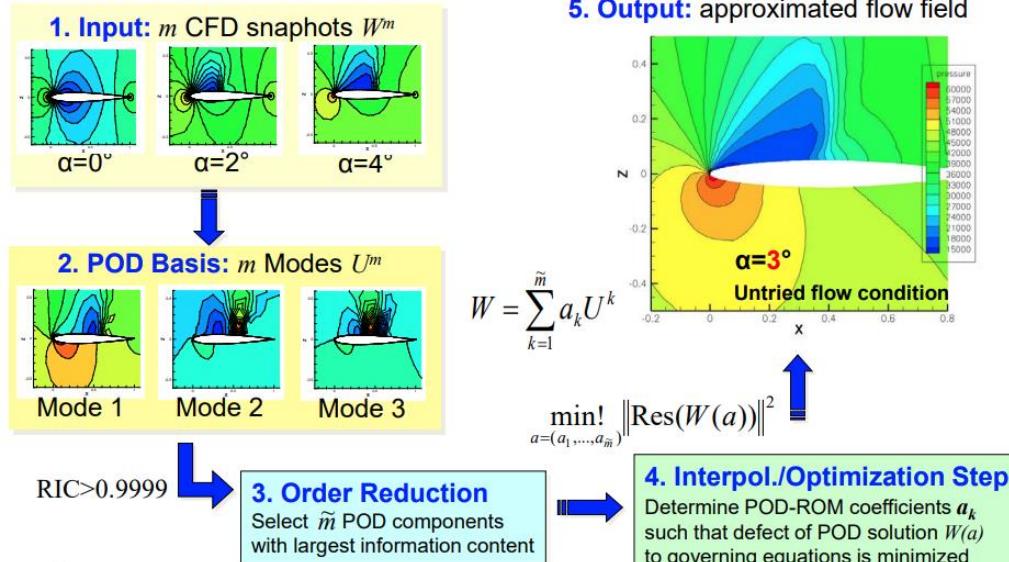
271   for (int i=0; i<nDim; i++) {
272     std::string uGradSigName;
273     uGradSigName = "uGradSigName_" + std::to_string(i);
274     volVectorField uGradSig(generateCustomField(runTime, mesh, uGradSigName),
275                             UMean&gradSigs[i]);
276
277     uGradSigs.push_back(uGradSig);
278
279     std::string sigGradUName;
280     sigGradUName = "sigGradUName_" + std::to_string(i);
281     volVectorField sigGradU(generateCustomField(runTime, mesh, sigGradUName),
282                           sigs[i]&gradU);
283     sigGradUs.push_back(sigGradU);
284   }
285
286   for (int i=0; i<nDim; i++) {
287     for (int j=0; j<nDim; j++) {
288       std::string sigGradSigName;
289       sigGradSigName = "sigGradSigName_" + std::to_string(i+nDim*j);
290       volVectorField sigGradSig(generateCustomField(runTime, mesh, sigGradSigName),
291                               sigs[i]&gradSigs[j]);
292       sigGradSigs[i+nDim*j] = sigGradSig;
293     }
294   }
295
296   std::vector<double> constant(nDim, 0.0);
297   std::vector<std::vector<double>> linear(nDim, std::vector<double>(nDim, 0.0));
298   std::vector<std::vector<std::vector<double>>> quadratic(nDim, std::vector<std::vector<double>>(nDim, std::vector<double>(nDim, 0.0)));
299
// this loop calculates the Galerkin system matrices Q L C for the ROM equation.
// constant term, linear term, and quadratic term
300   for (int k=0; k<nDim; k++) {
301     constant[k] = -1*innerProductPOD(sigs[k], UgradU, cellVolume) + (nu+nu_tilda)*innerProductPOD(sigs[k], laplUMean, cellVolume);
302     for (int m=0; m<nDim; m++) {
303       linear[k][m] = -1*innerProductPOD(sigs[k], uGradSigs[m], cellVolume) - innerProductPOD(sigs[k], sigGradUs[m], cellVolume) + (nu+nu_tilda)*innerProductPOD(sigs[k], laplSigs[m], cellVolume);
304       for (int n=0; n<nDim; n++) {
305         quadratic[k][m][n] = -1*innerProductPOD(sigs[k], sigGradSigs[m+nDim*n], cellVolume);
306       }
307     }
308   }
309
310
311 }
```

적합직교분해 (POD)

■ 절차

■ (4) expansion coefficient 보간

- 입력 매개변수 (AOA, etc.)
- POD's expansion coefficient
- Snapshot으로부터 보간 / 대체모델 생성



$$\vec{y}_j \approx \sum_{i=1}^d a_{ij} \vec{w}_i$$

04_Interpolation Coefficient

- POD's expansion coefficient(a_m)을 Design space의 변수에 따라 보간 또는 회귀모델 생성
- Linear interpolation using SciPy module (1D)
- 또는 RBF 모델 사용(2D, lsight 사용)

```

04_Interpolation_Coefficient
int writeSteps = 0;

if(writeFreq == 0){
    writeSteps = nSteps/numDirs;
}
else{
    writeSteps = writeFreq;
}

std::ofstream afiles;
afiles.open("avals.csv");

for (int t=0; t<nSteps+1; t++){
    cout << "t = " << timeElapsed << endl; // Case progress info in terminal
    for (int i=0; i<nDim; i++) {
        double da = constant[i];
        for (int j=0; j<nDim; j++) {
            da += linear[i+j*nDim]*prevAvals[j];
            for (int k=0; k<nDim; k++) {
                da += quadratic[i+j*nDim+k*nDim*nDim]*prevAvals[k]*prevAvals[j];
            }
        }
        avals[i] = prevAvals[i] + da*dt;
    }

    for (int i=0; i<nDim; i++){
        prevAvals[i] = avals[i]; // updating previous avals
        avalues[t][i] = avals[i]; // storing a values for in 2D vector for later
    }

    if(t%writeSteps == 0){
        double tcol = startTime + dt*t;
        afiles << tcol << ",";
        for (int j=0; j<nDim; j++){
            afiles << std::fixed << std::setprecision(16) << avalues[t][j] << ",";
        }
        afiles << endl << std::flush;
    }

    timeElapsed += dt;
}
afiles.close();

```

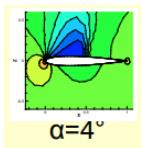
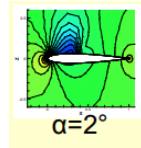
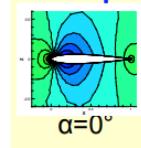
적합직교분해 (POD)

■ 절차

■ (4) expansion coefficient 보간

- 입력 매개변수 (AOA, etc.)
- POD's expansion coefficient
- Snapshot으로부터 보간 / 대체모델 생성

Snapshot
(Known)



$$U(0) = a_1(0) \sigma_1 + a_2(0) \sigma_2 + a_3(0) \sigma_3 \quad \text{const.}$$

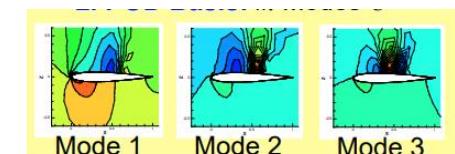
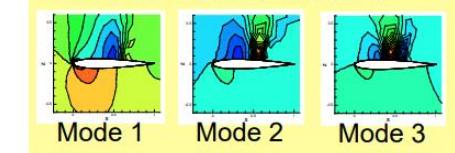
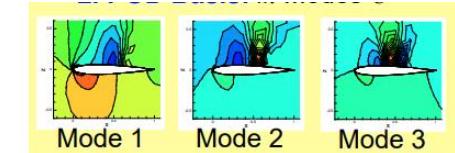
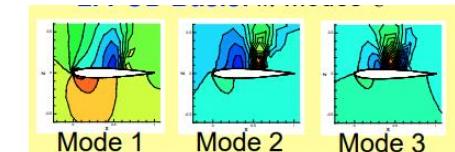
$$U(2) = a_1(2) \sigma_1 + a_2(2) \sigma_2 + a_3(2) \sigma_3$$

$$U(4) = a_1(4) \sigma_1 + a_2(4) \sigma_2 + a_3(4) \sigma_3$$

$$U(3) = a_1(3) \sigma_1 + a_2(3) \sigma_2 + a_3(3) \sigma_3$$

04_Interpolation Coefficient

- POD's expansion coefficient(a_m)을 Design space의 변수에 따라 보간 또는 회귀모델 생성
- Linear interpolation using SciPy module (1D)
- 또는 RBF 모델 사용(2D, lsight 사용)



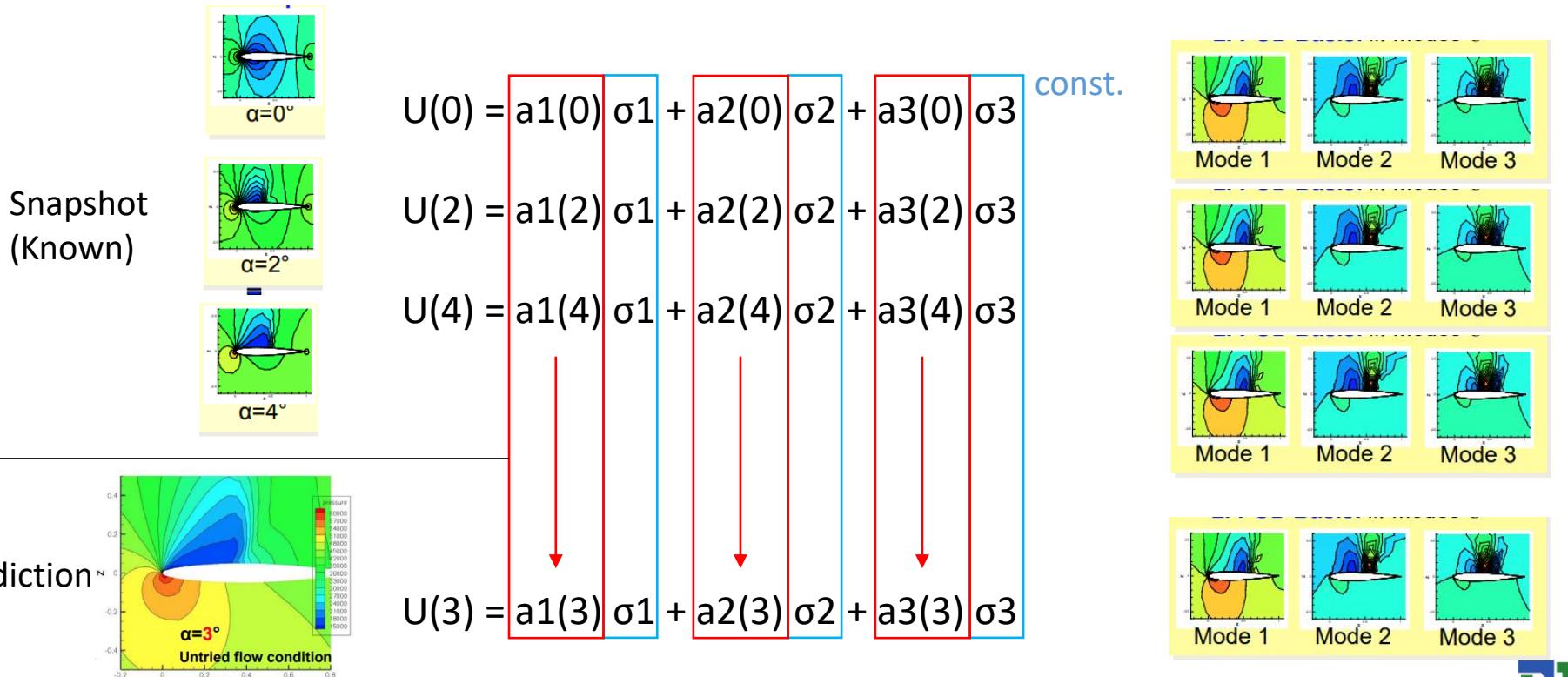
적합직교분해 (POD)

■ 절차

■ (5) 유동장 재생성

- Basis × Expansion coefficient

	<p>05_Reconstruction</p> <ul style="list-style-type: none"> 보간/회귀모델로 생성된 POD's expansion coefficient과 basis function을 이용하여 Flow field를 재생성 	$X^{new} = a_m \cdot \bar{\phi}_m$
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적합직교분해 (POD)

■ 절차

■ (5) 유동장 재생성

- Basis × Expansion coefficient

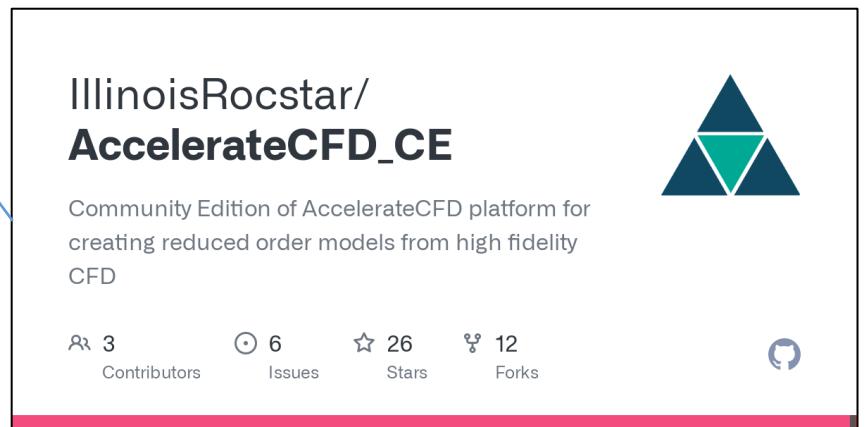
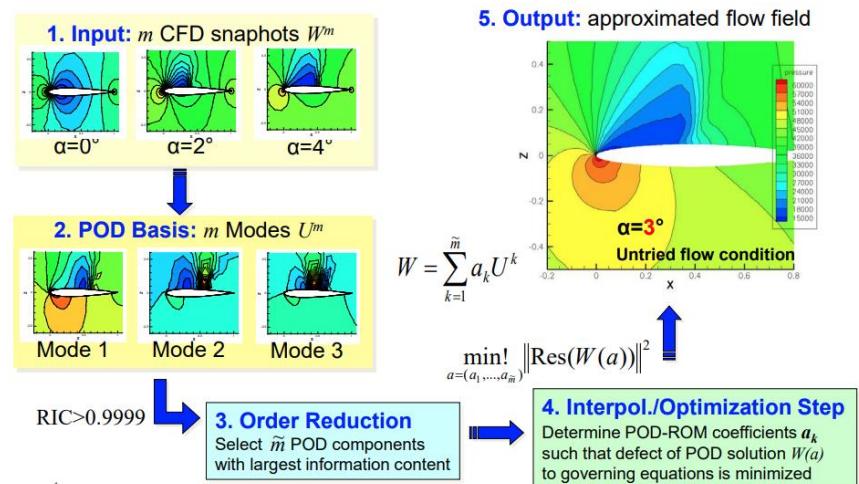
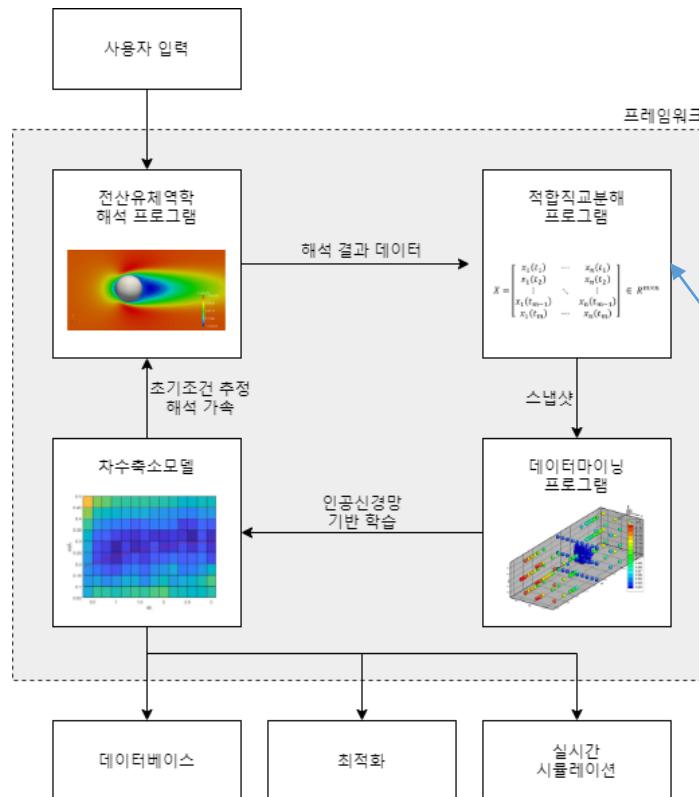
	<p>05_Reconstruction</p> <ul style="list-style-type: none">보간/회귀모델로 생성된 POD's expansion coefficient과 basis function을 이용하여 Flow field를 재생성	$X^{new} = a_m \cdot \bar{\phi}_m$

```
188 // this loops through all time directories in case and writes reconstructed..
189 // velocity (Urom)
190 forAll(timeDirs,timei)
191 {
192     std::vector<double> aVect = aVals[i];
193     Info << "t = " << timeDirs[timei].value() << nl; // Case progress info in terminal
194     runTime.setTime(timeDirs[timei],0);
195
196     std::string UName;
197     UName = "Urom";
198     volVectorField Urom(generateCustomField(runTime,mesh,UName),UMean);
199
200     for (int i=0; i<nDim; i++)
201     {
202         Urom += aVect[i]*sigs[i];
203
204         Urom.write(); // writes Urom in every time directory
205         i++;
206     }
207 }
```

적합직교분해 (POD)

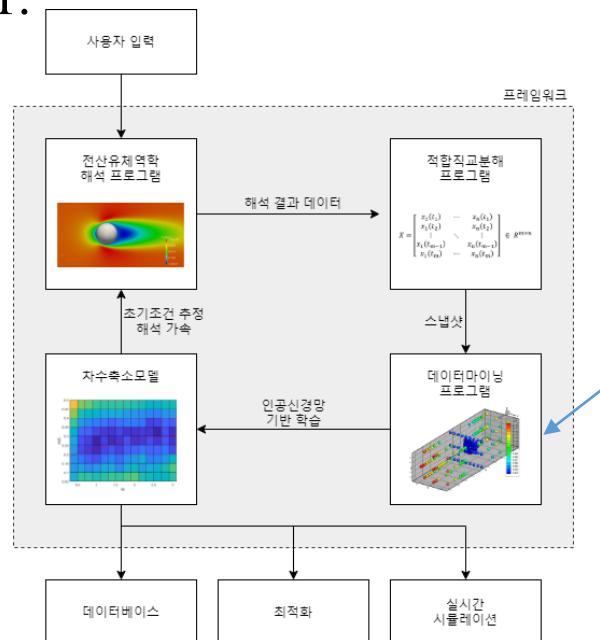
■ 요약

- 고차원 CFD 해석 결과를 저차원 basis들의 선형결합으로 압축 가능
- 프레임워크 구성 프로그램 선정
 - OpenFOAM
 - AccelerateCFD



데이터마이닝 (DAKOTA toolkit)

- 설계 및 데이터마이닝 툴박스
 - 각종 기법 알고리즘 수백 가지 내장
 - 키워드 input (ex. Surrogate, kriging, ann) 으로 간단한 실행
- GNU Lesser General Public License (versions 5.0+)
- M. S. Eldred et al., "DAKOTA : a multilevel parallel object- oriented framework for design optimization, parameter estimation, uncertainty quantification, and sensitivity analysis," E. United States. Department Of, Ed., ed: Sandia National Laboratories, 2011.



Mathematical and statistical methods to help scientists and engineers assess and improve the accuracy of computational models



데이터마이닝 (DAKOTA toolkit)

■ 작동 방식

■ 실행 목적

- 매개 변수 sweep
- 최적화
- 민감도 분석
- ...

■ 입력 변수 조작

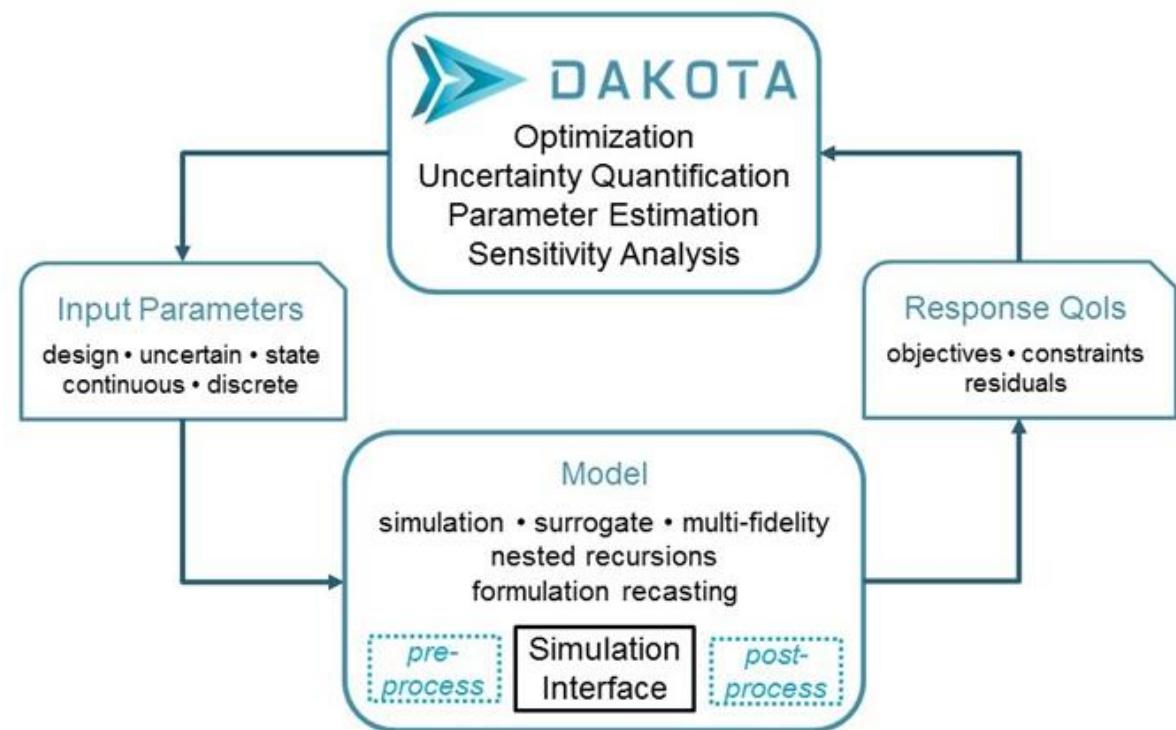
- 임의 형식 파일 대응 가능
 - Line / Column 지정

■ 솔버 실행

- Python 인터페이스 모듈 지원

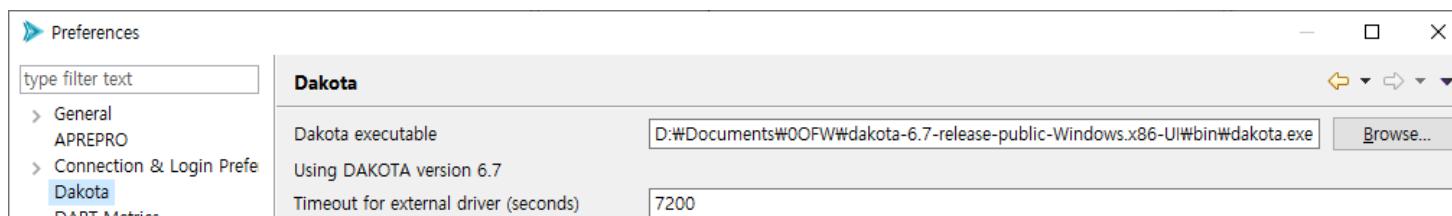
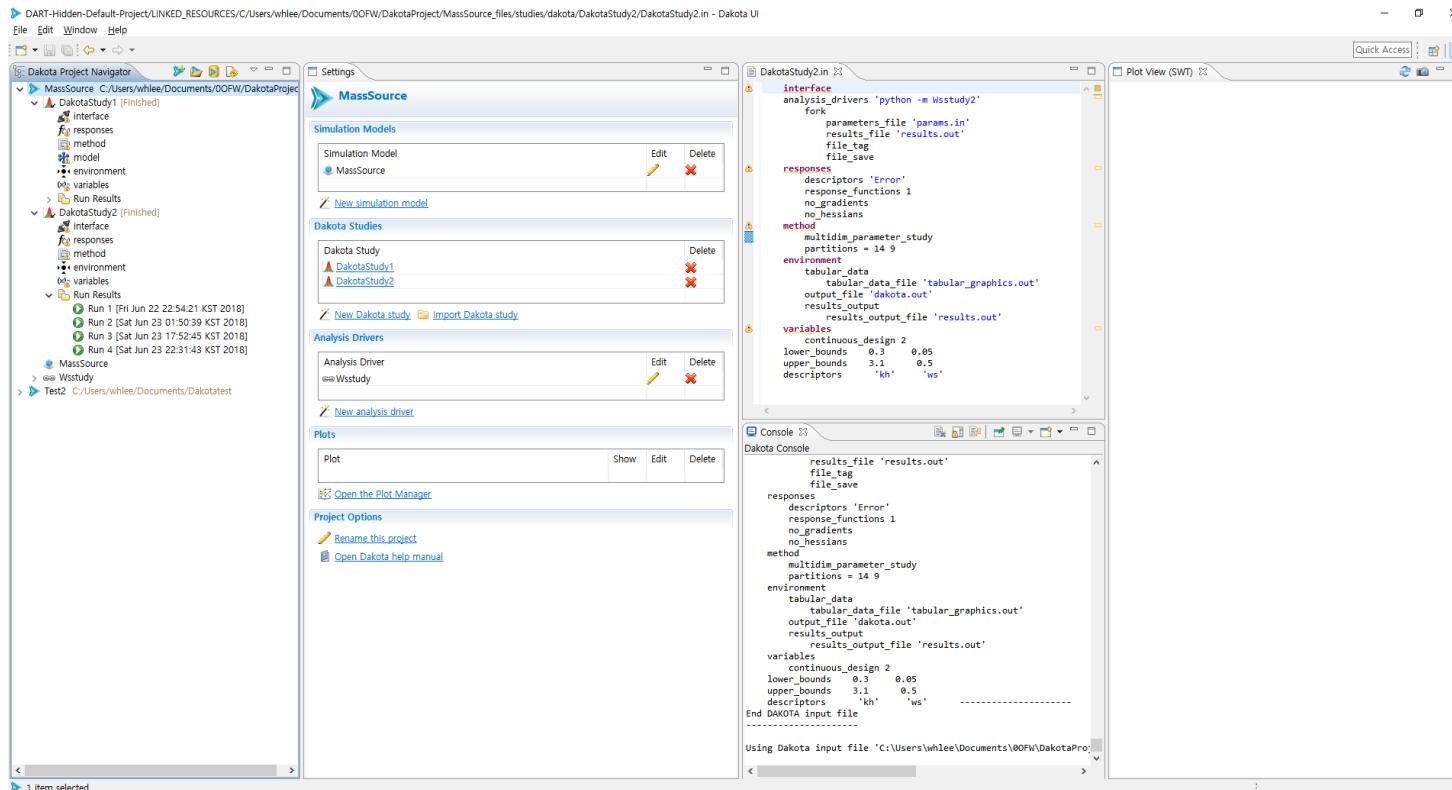
■ 출력 변수 인식

- 임의 형식 파일 혹은 stdout 인식 가능
 - Line / Column 지정



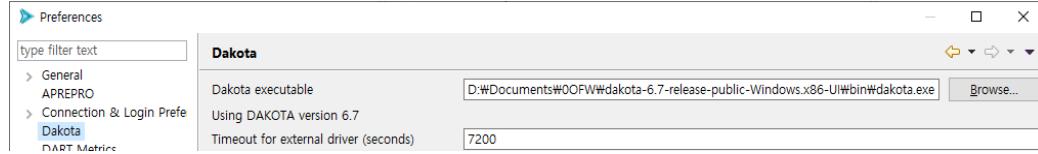
데이터마이닝 (DAKOTA toolkit)

■ CLI 프로그램 + GUI 지원



데이터마이닝 (DAKOTA toolkit)

- CLI 프로그램 + GUI 지원
 - GUI만으로는 구동 불가
 - Dakota Executable 연동 필요



- Linux binary는 RHEL만 제공
- Ubuntu 등에서 사용 시 직접 빌드 필요
 - mkdir build && cd build
 - cmake ..
 - 3.15 or higher
- 빌드 시 Python module 제공
 - Binary 다운로드 시 제공 X

Description

Dakota supports a library-linked interface to Python, but it must be explicitly enabled when compiling Dakota from source.

download.html

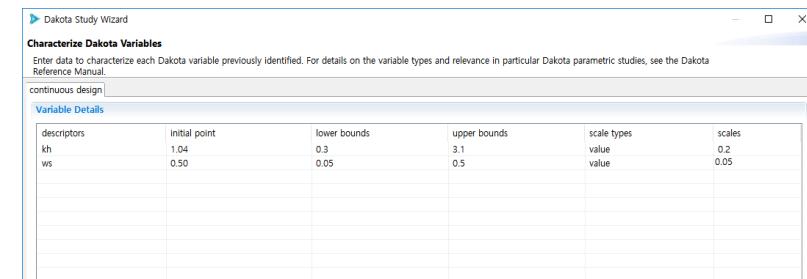
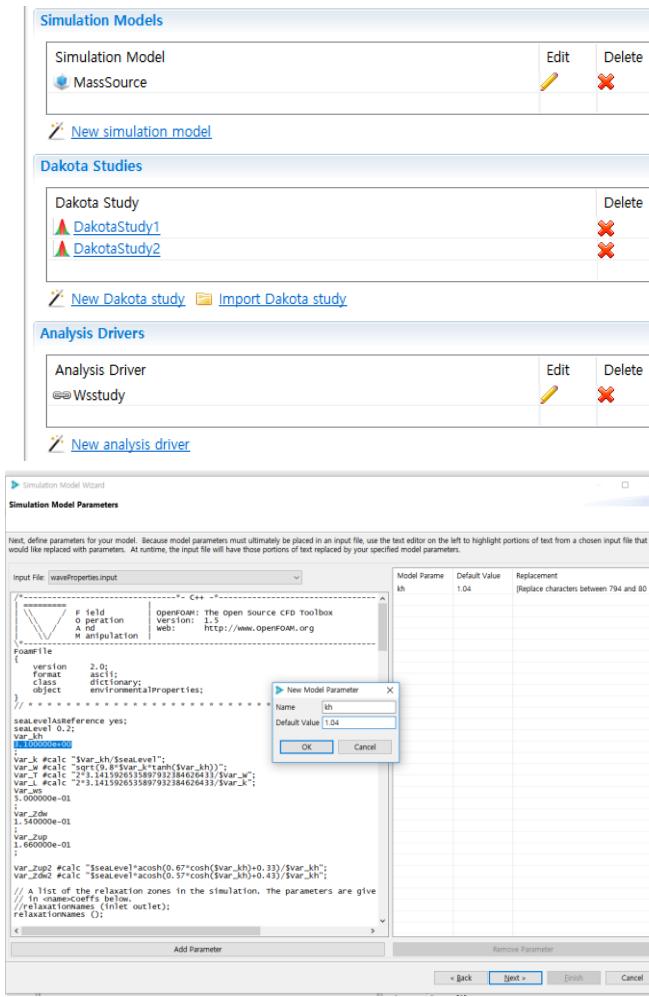
Public

Release 6.16.0 (*Release Notes*) — Released Mon, 05/

Platform	Link
Linux	dakota-6.16.0-release-public-Linux.x86_64-gui.tar.gz
Mac OS X	dakota-6.16.0-release-public-Darwin.x86_64-gui.tar.gz
Mac OS X	dakota-6.16.0-public-darwin.Darwin.x86_64-gui_cli.tar.gz
Mac OS X	dakota-6.16.0-public-darwin.Darwin.x86_64-cli.tar.gz
Windows	dakota-6.16.0-release-public-Windows.x64-gui.zip
Windows	dakota-6.16.0-public-windows.Windows.x64-gui_cli.zip
Windows	dakota-6.16.0-public-windows.Windows.x64-cli.zip
Source	dakota-6.16.0-release-public-src-gui.zip
Linux (RHEL7)	dakota-6.16.0-public-rhel7.Linux.x86_64-gui_cli.tar.gz
Linux (RHEL7)	dakota-6.16.0-public-rhel7.Linux.x86_64-cli.tar.gz
Source (Windows)	dakota-6.16.0-public-src-cli.zip
Source (Unix/OS X)	dakota-6.16.0-public-src-cli.tar.gz

데이터마이닝 (DAKOTA toolkit)

- 초기 사용자 GUI 지원
 - 솔버 실행 스크립트 자동 작성 지원



The screenshot shows a Notepad window with Python code for a Dakota study. The code is divided into three main sections:

- Pre-processing:** Substitutes values from the Dakota parameter file into the model's input file. It reads parameters from a file like "waveProperties.input" and replaces specific values (e.g., "kh" with 1.04).
- Execution:** Runs the computational model as a separate process. It uses subprocess.Popen to execute the model executable ("bash.exe") with specific arguments.
- Post-processing:** Finds quantities of interest in the model output and writes them to the Dakota output file. It reads the output file "rmseerror.output" and extracts specific error values.

```
# Pre-processing: Substitute values from the Dakota parameter file to the model's input file
input_map = readParamFile(sys.argv[1])
sub_file_input_map = {}
for file_key in mdl_in_map:
    in_file = open(os.path.join(sim_path, file_key), "r")
    in_file_sub = substituteModelInput(file_key, input_map, in_file)
    sub_file_input_map[file_key] = subName(file_key)
    #writeInput(sub_file_input_map[file_key], in_file_sub)
    writeInput("C:/meetingpoint/3imp", in_file_sub)

# Execution: Run the computational model as a separate process.
#model_exec_args = model_exec_args.replace(" "+model_name+" ", " "+os.path.join(sim_path, model_name)+" ")
model_exec_args = model_exec_args.replace(model_name, " "+os.path.join(sim_path, model_name))
for file_key in mdl_in_map:
    new_file_key = sub_file_input_map[file_key]
    model_exec_args = model_exec_args.replace(" "+file_key+" ", " "+new_file_key+" ")
#model_exec = subprocess.Popen(shlex.split(model_exec_args.replace('\\','/'),999), stdout=PIPE)
#model_exec = subprocess.Popen(shlex.split('C:/Windows/System32/bash.exe'), stdout=PIPE, stderr=PIPE)
model_exec = subprocess.Popen(shlex.split("bash -i"), stdout=PIPE, stderr=PIPE, shell=True)
stdout, stderr = model_exec.communicate()

if stderr:
    print "ERROR WAS ENCOUNTERED:", stderr

# Post-processing: Find quantities of interest in the model output, and write to the Dakota output file
for file_key in mdl_out_map:
    output_map = None
    if file_key == "stdout":
        output_map = findModelOutput(file_key, stdout)
    else:
        out_file_str = ""
        with open(os.path.join(sim_path, file_key), "r") as out_file:
            out_file_str=out_file.read()
        output_map = findModelOutput(file_key, out_file_str)
    writeOutput(sys.argv[2], output_map)
```

데이터마이닝 (DAKOTA toolkit)

- 숙련 사용자 CLI 실행
 - 간소한 양식의 input file로 용이한 실행

The screenshot shows a terminal window titled "rosen_opt_sbo.in" containing the Dakota input file "rosen_opt_sbo.in". The file defines an optimization problem using a neural network surrogate model. The terminal output on the right shows the execution of the Dakota toolkit, displaying progress messages, function evaluations, and statistical results for the objective function.

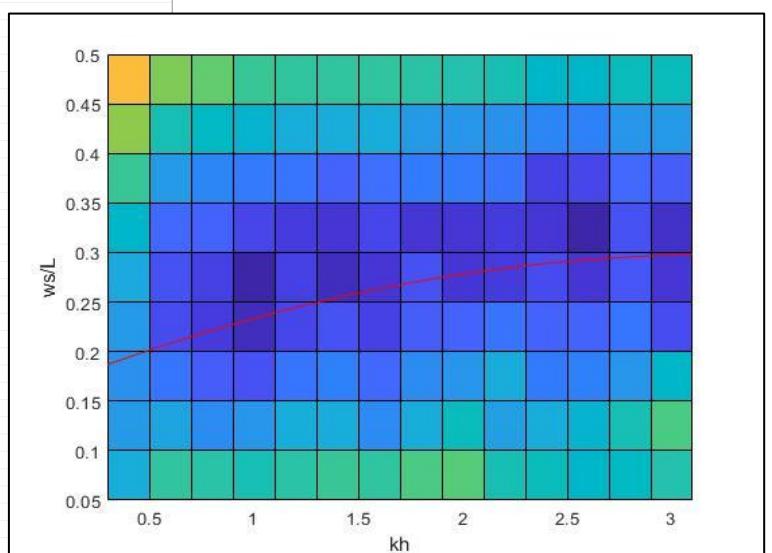
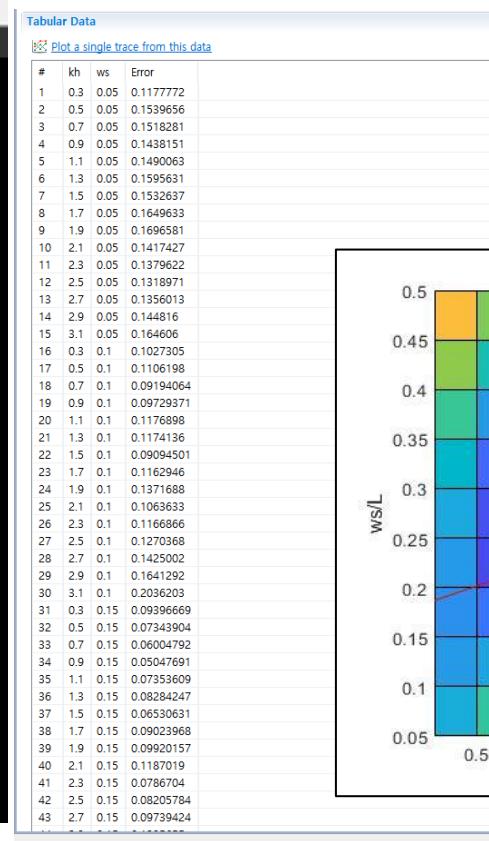
```
# Dakota Input File: rosen_opt_sbo.in
environment
tabular_data
tabular_data_file = 'rosen_opt_sbo.dat'
top_method_pointer = 'SBLO'
method
id_method = 'SBLO'
surrogate_based_local
model_pointer = 'SURROGATE'
method_pointer = 'NLP'
max_iterations = 500
trust_region
initial_size = 0.10
minimum_size = 1.0e-6
contract_threshold = 0.25
expand_threshold = 0.75
contraction_factor = 0.50
expansion_factor = 1.50
method
id_method = 'NLP'
conmin_frcg
max_iterations = 50
convergence_tolerance = 1e-8
model
id_model = 'SURROGATE'
surrogate_global
correction additive zeroth_order
polynomial quadratic
dace_method_pointer = 'SAMPLING'
responses_pointer = 'SURROGATE_RESP'
variables
continuous_design = 2
initial_point -1.2 1.0
lower_bounds -2.0 -2.0
upper_bounds 2.0 2.0
descriptors 'x1' 'x2'

whlee@DESKTOP-ECSUKD1: /mnt/c/Users/whlee/Documents/code/220704_dakota_suropt_test/src
<===== global_neural_network approximation builds completed.
Beginning Approximate Fn Evaluations...
<===== Function evaluation summary (APPROX_INTERFACE_1): 100 total (100 new, 0 duplicate)
<===== Function evaluation summary (SimulationInterface): 10 total (10 new, 0 duplicate)
<===== Best parameters =
      5.1892411019e+01 x1
      7.7085409127e-01 x2
<===== Best objective function =
      -7.3934829807e+01
<===== Best evaluation ID: 37
-----
Statistics based on 100 samples:
Sample moment statistics for each response function:
      Mean   Std Dev   Skewness   Kurtosis
obj_fn  2.4296102525e+03  1.9546250171e+03  5.0011856105e-01 -8.9182874568e-01
95% confidence intervals for each response function:
      LowerCI_Mean   UpperCI_Mean   LowerCI_StdDev   UpperCI_StdDev
obj_fn  2.04177024932e+03  2.8174502618e+03  1.7161741452e+03  2.2706395157e+03
Simple Correlation Matrix among all inputs and outputs:
      x1          x2          obj_fn
x1  1.000000e+00
x2  2.08607e-02  1.000000e+00
obj_fn  2.23949e-02  8.15550e-01  1.000000e+00
Partial Correlation Matrix between input and output:
      obj_fn
x1  9.19860e-03
x2  8.15465e-01
Simple Rank Correlation Matrix among all inputs and outputs:
      x1          x2          obj_fn
x1  1.000000e+00
x2  2.26463e-02  1.000000e+00
obj_fn  3.69757e-02  8.19694e-01  1.000000e+00
Partial Rank Correlation Matrix between input and output:
```

데이터마이닝 (DAKOTA toolkit)

- DAKOTA input 예시 (parametric study)
 - \$ dakota DakotaStudy.in

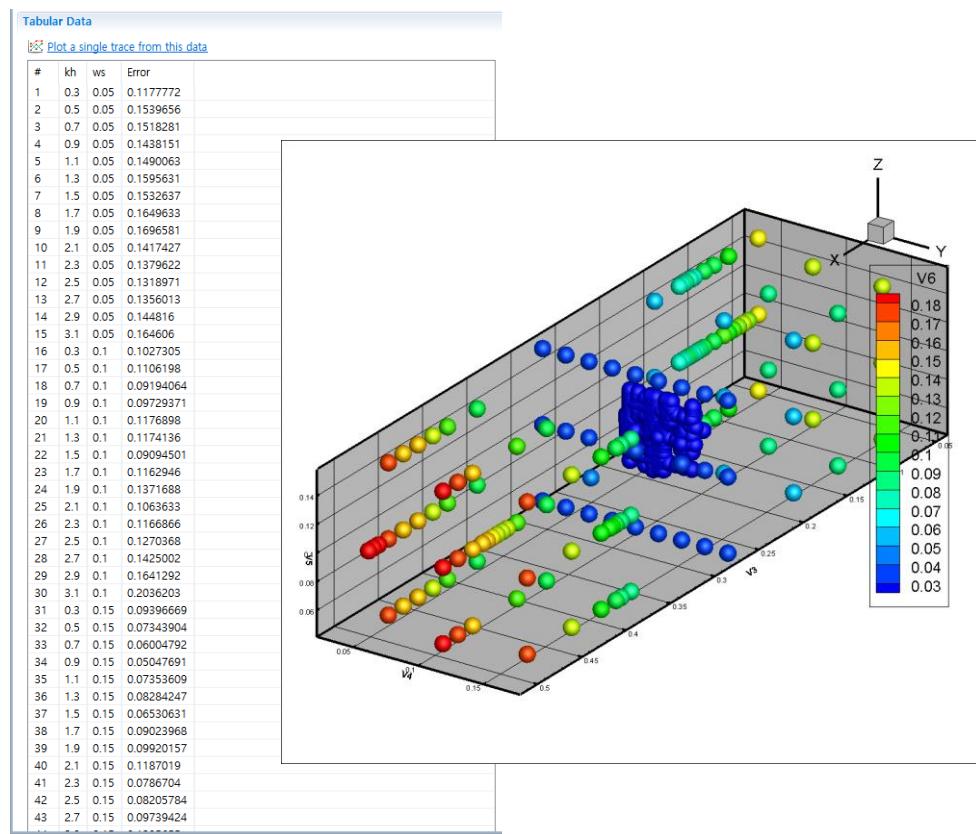
```
DakotaStudy2.in
1  interface
2      analysis_drivers 'python -m Wsstudy2'
3          fork
4              parameters_file 'params.in'
5                  results_file 'results.out'
6                  file_tag
7                  file_save
8  responses
9      descriptors 'Error'
10     response_functions 1
11     no_gradients
12     no_hessians
13 method
14     multidim_parameter_study
15     partitions = 14 9
16 environment
17     tabular_data
18         tabular_data_file 'tabular_graphics.out'
19         output_file 'dakota.out'
20         results_output
21             results_output_file 'results.out'
22 variables
23     continuous_design 2
24     lower_bounds    0.3      0.05
25     upper_bounds    3.1      0.5
26     descriptors      'kh'     'ws'
```



데이터마이닝 (DAKOTA toolkit)

- DAKOTA input 예시 (최적화)
 - \$ dakota DakotaStudy.in

```
DakotaStudy3.in
1      interface
2          analysis_drivers 'python -m Wsstudy3'
3              fork
4                  parameters_file 'params.in'
5                  results_file 'results.out'
6                  file_tag
7                  file_save
8      responses
9          descriptors 'Error'
10         objective_functions 1
11             sense 'min'
12             numerical_gradients
13             numerical_hessians
14     method
15     coliny_direct
16         max_function_evaluations 999999
17     model
18         single
19     environment
20         tabular_data
21             tabular_data_file 'tabular_graphics.out'
22             output_file 'dakota.out'
23             results_output
24                 results_output_file 'results.out'
25     variables
26         continuous_design 3
27             initial_point 0.30 0.154 0.166
28             lower_bounds 0.05 0.02 0.02
29             upper_bounds 0.5 0.18 0.18
30                 descriptors 'ws' 'zd' 'zu'
```



데이터마이닝 (DAKOTA toolkit)

■ DAKOTA input 예시 (대체모델 기반 최적화)

D:\#Documents#\OOFW#\dakota-6.7-release-public-Windows.x86-UI#\examples#\users#\rosen_opt_sbo.in - Notepad++

파일(F) 폴더(O) 찾기(O) 보기(V) 인코딩(N) 언어(S) 설정(I) 도구(O) 매크로 실행 플러그인 정 관리 ?

罗斯顿_优化.sbo.in

Dakota Input File: rosen_opt_sbo.in

environment

tabular_data

tabular_data_file = 'rosen_opt_sbo.dat'

top_method_pointer = 'SBLO'

method

id_method = 'SBLO'

surrogate_based_local

model_pointer = 'SURROGATE'

method_pointer = 'NLP'

max_iterations = 500

trust_region

initial_size = 0.10

minimum_size = 1.0e-6

contract_threshold = 0.25

expand_threshold = 0.75

contraction_factor = 0.50

expansion_factor = 1.50

method

id_method = 'NLP'

conmin_frcg

max_iterations = 50

convergence_tolerance = 1e-8

model

id_model = 'SURROGATE'

surrogate_global

correction additive zeroth_order

polynomial quadratic

dace_method_pointer = 'SAMPLING'

responses_pointer = 'SURROGATE_RESP'

variables

continuous_design = 2

initial_point -1.2 1.0

lower_bounds -2.0 -2.0

upper_bounds 2.0 2.0

descriptors 'x1' 'x2'

responses

id_responses = 'SURROGATE_RESP'

objective_functions = 1

numerical_gradients

method_source dakota

interval_type central

fd_step_size = 1.e-6

no_hessians

method

id_method = 'SAMPLING'

sampling

samples = 10

seed = 531

sample_type lhs

model_pointer = 'TRUTH'

model

id_model = 'TRUTH'

single

interface_pointer = 'TRUE_FN'

responses_pointer = 'TRUE_RESP'

interface

id_interface = 'TRUE_FN'

analysis_drivers = 'rosenbrock'

direct

deactivate evaluation_cache restart_file

responses

id_responses = 'TRUE_RESP'

objective_functions = 1

no_gradients

no_hessians

대체모델
생성 기법

length : 1,617 lines : 77 Ln : 77 Col : 1 Pos : 1,618 Windows (CR LF) UTF-8 INS

데이터마이닝 (DAKOTA toolkit)

- DAKOTA input 예시 (Multi-fidelity surrogate model)

The screenshot shows a Notepad++ interface with two tabs open. The left tab is titled 'rosenrock_sbo_hierarchical.in' and contains the following configuration file content:

```
1 environment,
2     ## Activate Dakota's legacy X Windows-based graphics on Unix systems
3     ## Consider newer capabilities in the Dakota GUI
4     graphics
5     tabular_data
6     top_method_pointer = 'SBLO'
7
8 method,
9     id_method = 'SBLO'
10    surrogate_based_local
11    model_pointer = 'SURROGATE'
12    method_pointer = 'NLP'
13    trust_region
14        initial_size = 0.10
15        contract_threshold = 0.25
16        expand_threshold = 0.75
17        contraction_factor = 0.50
18        expansion_factor = 1.50
19
20 method,
21     id_method = 'NLP'
22     ## (NPSOL requires a software license; if not available, try
23     ## conmin_frcg or optpp_newton instead)
24     npsol_sqp
25         max_iterations = 50
26         convergence_tolerance = 1e-10
27
28 model,
29     id_model = 'SURROGATE'
30     surrogate_hierarchical
31     ordered_model_fidelities = 'LOFI' 'HIFI'
32     correction additive second_order
33
34 variables,
35     continuous_design = 2
36     initial_point -1.2 1.0
37     lower_bounds -2.0 -2.0
38     upper_bounds 2.0 2.0
39     descriptors 'x1' 'x2'
40
41 responses,
```

The right tab is also titled 'rosenrock_sbo_hierarchical.in' and contains the following configuration file content:

```
41
42 responses,
43     objective_functions = 1
44     analytic_gradients
45     analytic_hessians
46
47 model,
48     id_model = 'LOFI'
49     single
50     interface_pointer = 'LOFI_FN'
51
52 interface,
53     id_interface = 'LOFI_FN'
54     analysis_drivers = 'lf_rosenbrock'
55     direct
56     deactivate restart_file
57
58 model,
59     id_model = 'HIFI'
60     single
61     interface_pointer = 'HIFI_FN'
62
63 interface,
64     id_interface = 'HIFI_FN'
65     analysis_drivers = 'rosenbrock'
66     direct
67     deactivate restart_file
68
```

데이터마이닝 (DAKOTA toolkit)

■ DAKOTA 스크립트 예시

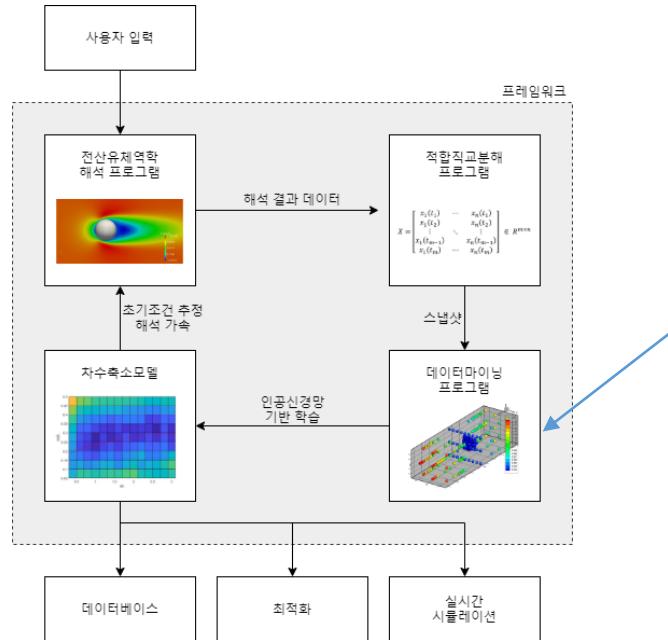
- dakota.interfacing 모듈을 통해 연동
 - PYTHONPATH 환경변수에 Dakota 설치 경로 추가
- 솔버 실행 / 입출력 순서 임의 구성 가능

```
1 #!/usr/bin/env python3
2 import numpy as np
3 import dakota.interfacing as di
4
5 params, results = di.read_parameters_file()
6
7 write_input(params)
8
9 run_solver()
10
11 results = read_output()
12
13 results.write()
```

데이터마이닝 (DAKOTA toolkit)

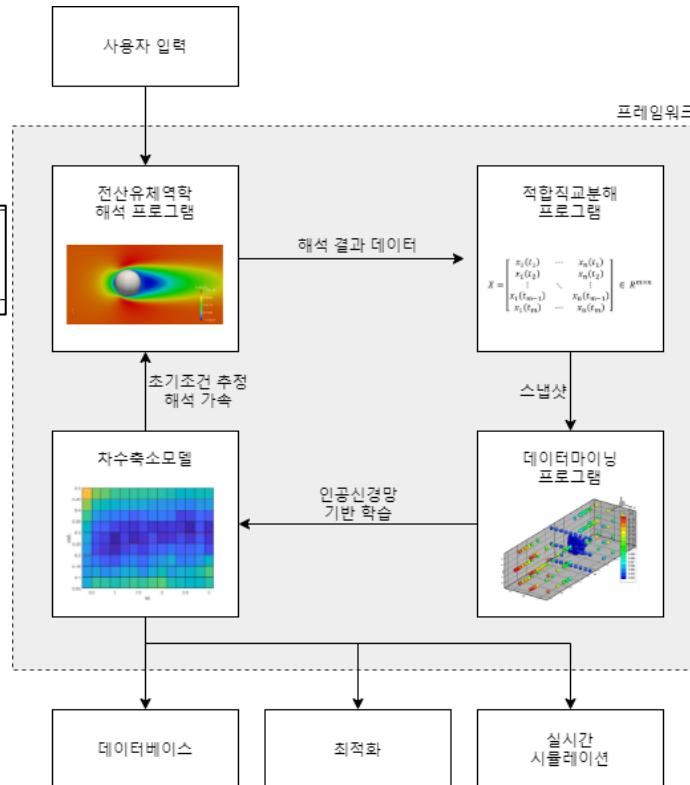
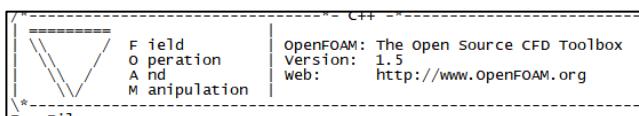
■ 요약

- 수백개의 기법을 간단한 input 파일 작성으로 사용 가능한 데이터마이닝 툴박스 DAKOTA
- 프레임워크 구성 요소 프로그램으로 선정



프레임워크 개발

- 구성 요소 프로그램
 - OpenFOAM
 - AccelerateCFD
 - DAKOTA



**IllinoisRocstar/
AccelerateCFD_CE**

Community Edition of AccelerateCFD platform for creating reduced order models from high fidelity CFD

3 Contributors 6 Issues 26 Stars 12 Forks



Mathematical and statistical methods to help scientists and engineers assess and improve the accuracy of computational models



DAKOTA
Explore and predict with confidence

프레임워크 개발

■ AccelerateCFD 보완

■ 기존

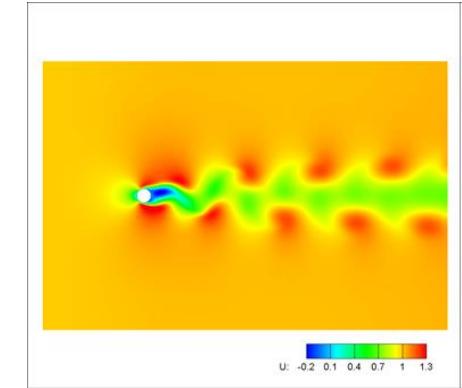
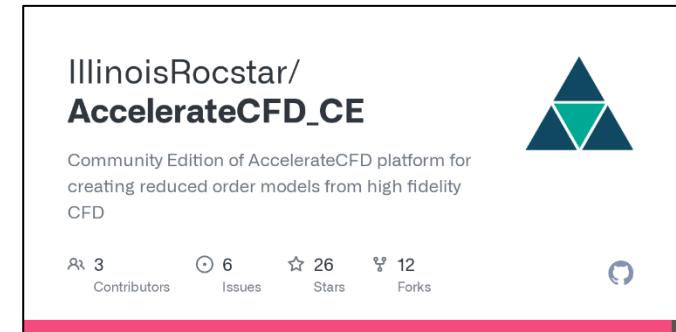
- Umean + U(t)
- Vector U(t) 의 시간에 따른 POD 추정 용도

■ 개량

- Umean 등 시간 평균 물리량 없이 POD를 수행하도록 수정
- Scalar P에 대한 POD 추가
- podROM 이후

신규 입력조건에 대한 expansion coefficient를
추가 행으로 직접 기록하는 절차 추가

- => podFlowReconstruct에서 신규 유동장 생성
- Expansion coefficient 보간에 DAKOTA 사용



```
85 12.5991600000000012,-0.0693057512687179,-0.1339006029651310,-0.0207532644376687,-0.013049781112034,0.0011800165798592,  
86 12.7491500000000002,0.04592808961045100,-0.1440396562984638,-0.0128330907700062,0.0223013963593660,-0.0008483570761670,  
87 12.8991400000000009,0.1332179382399867,-0.0683375036788400,0.0221005212119751,0.0128224325816151,-0.0002058570163162,  
88 13.0491300000000017,0.1390199634468887,0.0469109259700145,0.0115742263361113,-0.0222412342683103,0.0010644760384889,  
89 13.1991200000000006,0.0599394305805869,0.1311692048711383,-0.0224389413060522,-0.0109656204400389,-0.0013391637849161,  
90 13.3491100000000014,-0.0554414918077339,0.1330694382196505,-0.0095945744277698,0.0225072508508146,0.0009611004502626,  
91 13.4991000000000003,-0.1366694682455949,0.0512435328522021,0.0232928128399615,0.0084756140662863,-0.0002845926286241,  
92 13.6490900000000011,-0.1338251272003634,-0.0641111796107720,0.0075689723628032,-0.0232688799541422,-0.0010006680928169,  
93 13.7990800000000018,-0.0489346767104756,-0.1419509074480909,-0.0237595245725301,-0.0058962798581446,0.0011893254758908,  
94 13.9490700000000007,0.0658592268507926,-0.1350081499591214,-0.0056720592052062,0.0250283322828671,-0.0012461203150463,  
95 14.0990600000000015,0.1405168875164147,-0.0477514676189012,0.024614912836608,0.0055390426245586,0.0004149955424102,  
96 14.2490500000000004,0.1292046263280497,0.0664293679184010,0.0043421317940559,-0.0246470606172846,0.0005651070976550,  
97 14.3990400000000012,0.0391072649817660,0.1377133430606408,-0.0244819198335083,-0.0037042434384433,-0.0012108524841335,  
98 14.5490300000000019,-0.0745809908107926,0.1226416020205486,-0.002279283069106,0.0242403232407572,0.0012326165070807,  
99 14.6690200000000009,-0.1423793349210730,0.0301806476744027,0.024674815199811,0.00123952092013030,-0.0008729662446150,  
100 14.8490100000000016,-0.1226078935902394,-0.0828002942144563,0.0002771795817250,-0.0243053187103875,-0.0004372786646340,  
101 14.999000000000006,-0.0277564215228415,-0.1468467960749114,-0.0245856912556688,0.0014876802015931,0.0009411264980396,  
102
```

프레임워크 개발

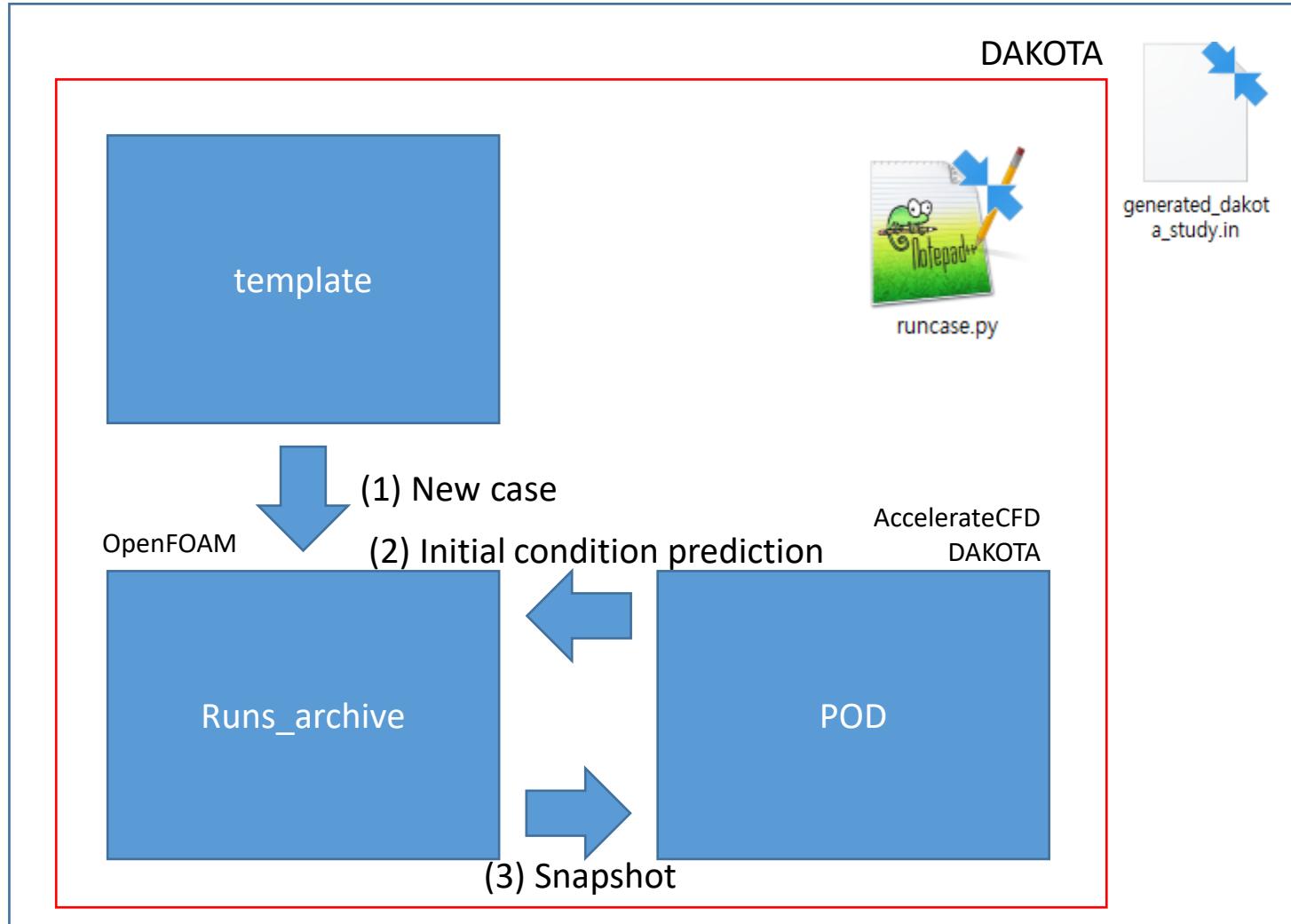
■ 구성



- Template : 해석 케이스 템플릿
- Runs_archive : 입력 조건 별 해석이 진행/완료된 케이스 폴더 모음
- POD : 해석 결과를 취합하여 AccelerateCFD로 POD를 수행하는 폴더
- Runcase.py : POD로 신규 케이스 초기조건 생성 및 해석 수행
- Framework.py : runcase.py를 driver script로 하여 DAKOTA 실행
- INPUT.INP : framework.py의 input으로 프레임워크 실행 목적/조건 지정

프레임워크 개발

■ 작동 순서

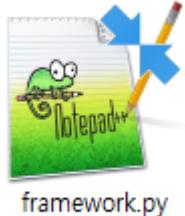


INPUT.INP

프레임워크 개발

- \$ python3 framework.py

- Template 폴더 내 input 파일 경로, input 개수, 위치, 범위
- 실행 명령어
- 출력 파일 경로 및 출력값 위치
- 실행 목적 (DB 생성 / 최적화)



framework.py



INPUT.INP

```
INPUT.INP
1 INPUT_FILE_LOCATION
2 INPUT.INP
3
4 #_OF_INPUT_VARIABLE
5 2
6
7 LOCATION_OF_INPUT_VARIABLE (Line / column)
8 7 1
9 8 1
10
11 BOUND
12 -20 20
13 -20 20
14
15 RUN_COMMAND
16 run.sh
17
18 OUTPUT_FILE_LOCATION
19 postProcessing/forces/0/forceCoeffs.dat
20
21 LOCATION_OF_OUTPUT_VARIABLE (Line / column)
22 1009 1
23
24 OBJECTIVE (1: DB, 2: OPTIMIZATION)
25 1|
```



```
rosen_opt_sbo.in
1 # Dakota Input File: rosen_opt_sbo.in
2
3 environment
4   tabular_data
5     tabular_data_file = 'rosen_opt_sbo.dat'
6     top_method_pointer = 'SBLO'
7
8 method
9   id_method = 'SBLO'
10  surrogate_based_local
11    model_pointer = 'SURROGATE'
12    method_pointer = 'NLP'
13    max_iterations = 500
14  trust_region
15    initial_size = 0.10
16    minimum_size = 1.0e-6
17    contract_threshold = 0.25
18    expand_threshold = 0.75
19    contraction_factor = 0.50
20    expansion_factor = 1.50
21
22 method
23   id_method = 'NLP'
24   connmin_frcg
25     max_iterations = 50
26     convergence_tolerance = 1e-8
27
28 model
29   id_model = 'SURROGATE'
30   surrogate_global
31     correction additive zeroth_order
32     polynomial quadratic
33     dace_method_pointer = 'SAMPLING'
34     responses_pointer = 'SURROGATE_RESP'
35
36 variables
37   continuous_design = 2
38     initial_point -1.2 1.0
39     lower_bounds -2.0 -2.0
40     upper_bounds 2.0 2.0
41     descriptors 'x1' 'x2'
42
```

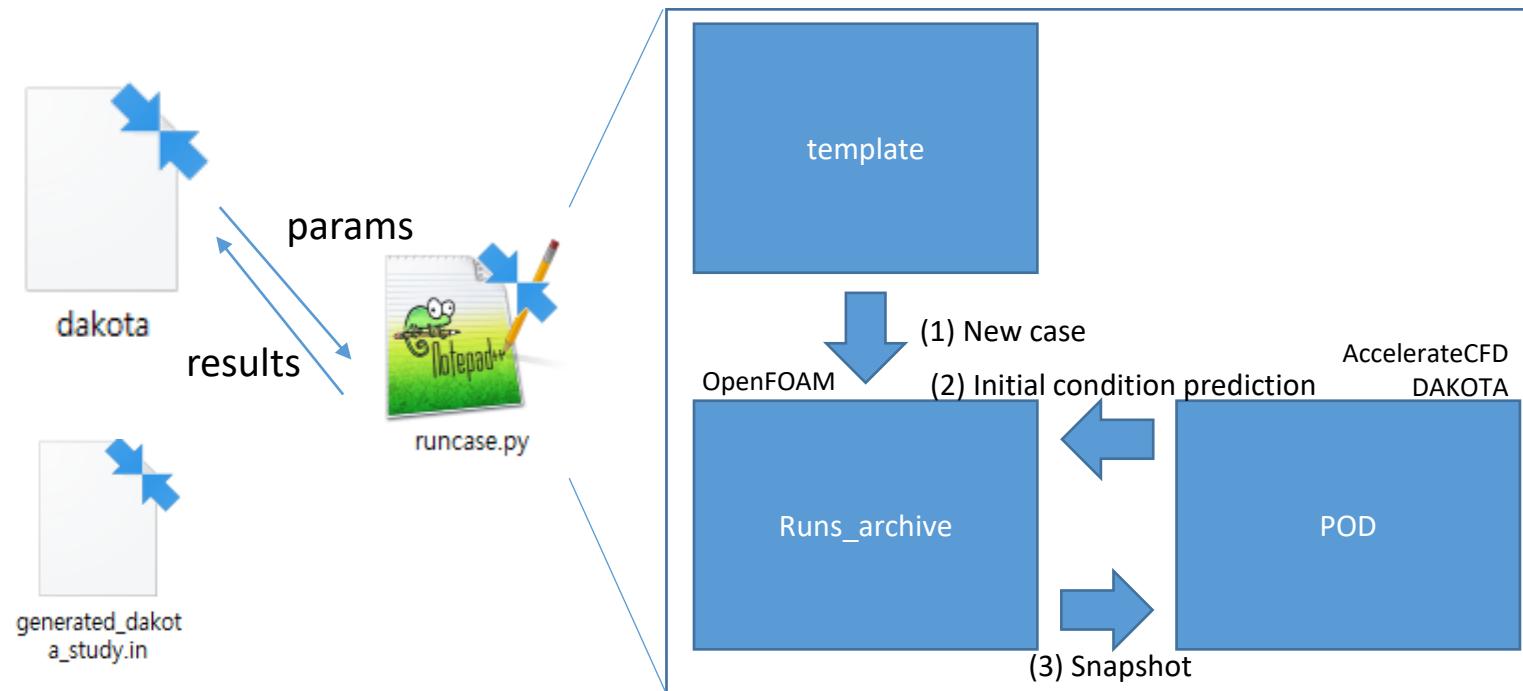


generated_dakota_study.in

프레임워크 개발

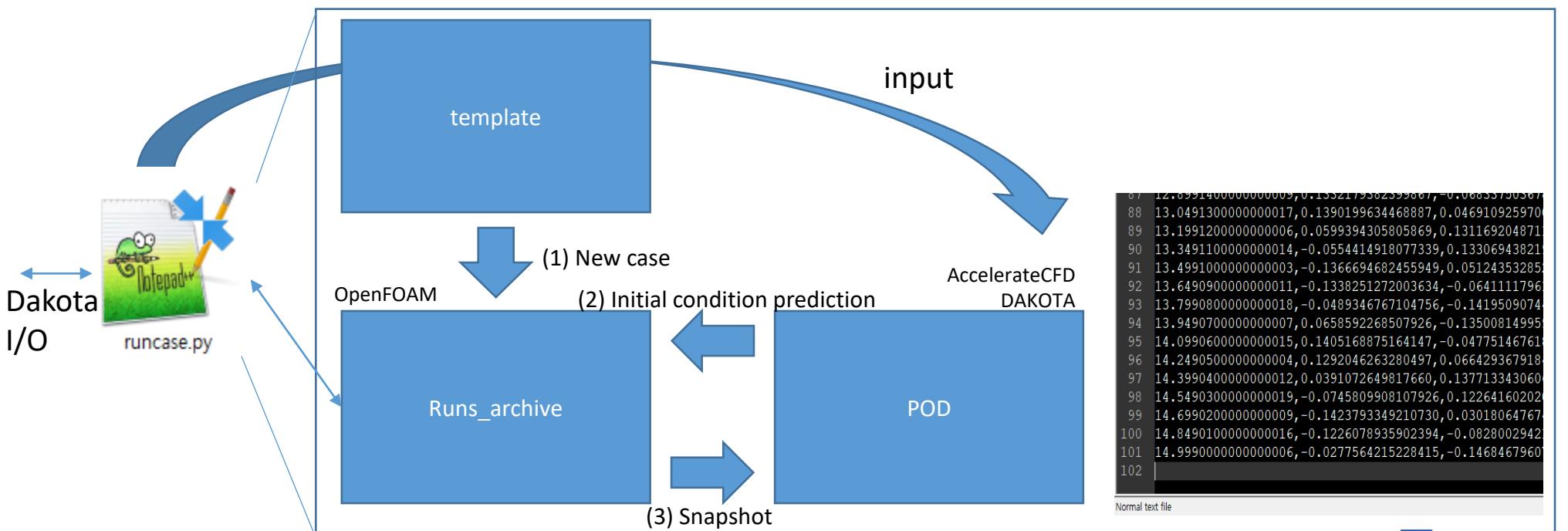
- (하위 프로세스 실행) \$ dakota generated_dakota_study.in
 - dakota.interfacing 모듈을 통해 runcase.py와 입출력 교환
 - DAKOTA가 시도할 입력 매개 변수를 전달하면 runcase.py가 솔버 실행
 - 해석이 완료되면 runcase.py가 결과값을 읽어 DAKOTA에 전달
 - 입출력을 반복하며 DAKOTA의 실행 목적 (DB/최적화) 달성

```
rosen_opt_sbo.in
1 # Dakota Input File: rosen_opt_sbo.in
2
3 environment
4   tabular_data
5     tabular_data_file = 'rosen_opt_sbo.dat'
6   top_method_pointer = 'SBLO'
7
8 method
9   id_method = 'SBLO'
10  surrogate_based_local
11    model_pointer = 'SURROGATE'
12    method_pointer = 'NLP'
13    max_iterations = 500
14  trust_region
15    initial_size = 0.10
16    minimum_size = 1.0e-6
17    contract_threshold = 0.25
18    expand_threshold = 0.75
19    contraction_factor = 0.50
20    expansion_factor = 1.50
21
22 method
23   id_method = 'NLP'
24   commin_frcg
25   max_iterations = 50
26   convergence_tolerance = 1e-8
27
28 model
29   id_model = 'SURROGATE'
30   surrogate_global
31   correction additive zeroth_order
32   polynomial quadratic
33   dace_method_pointer = 'SAMPLING'
34   responses_pointer = 'SURROGATE_RESP'
35
36 variables
37   continuous_design = 2
38   initial_point -1.2 1.0
39   lower_bounds -2.0 -2.0
40   upper_bounds 2.0 2.0
41   descriptors 'x1' 'x2'
42
```



프레임워크 개발

- (하위 프로세스 반복 실행) \$ python3 runcase.py
 - Template 폴더를 복제하여 runs_archive에 신규 케이스 생성
 - POD 폴더에 input 매개변수를 전달 및 DAKOTA를 실행
 - Expansion coefficient를 DAKOTA로 보간하여 해석 초기조건 예측
 - 예측된 초기조건을 runs_archive로 복사하고 그대로 해석하여 빠른 수렴
 - Runcase.py에 결과값 전달 및 POD snapshot 추가

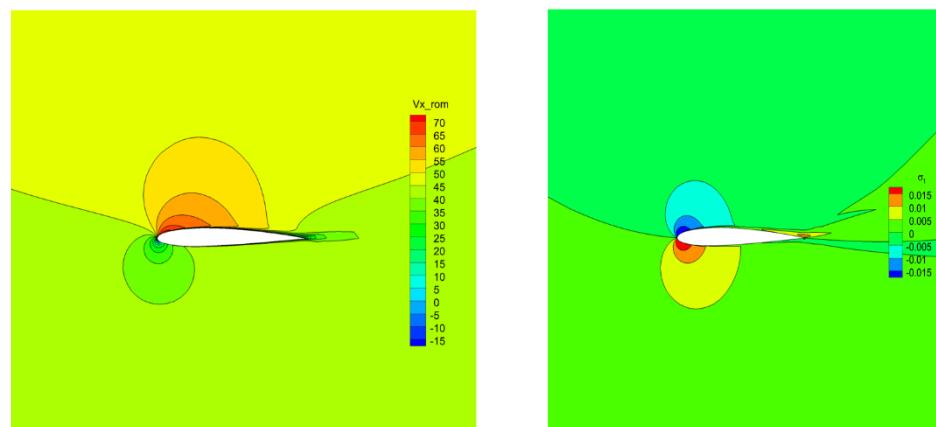
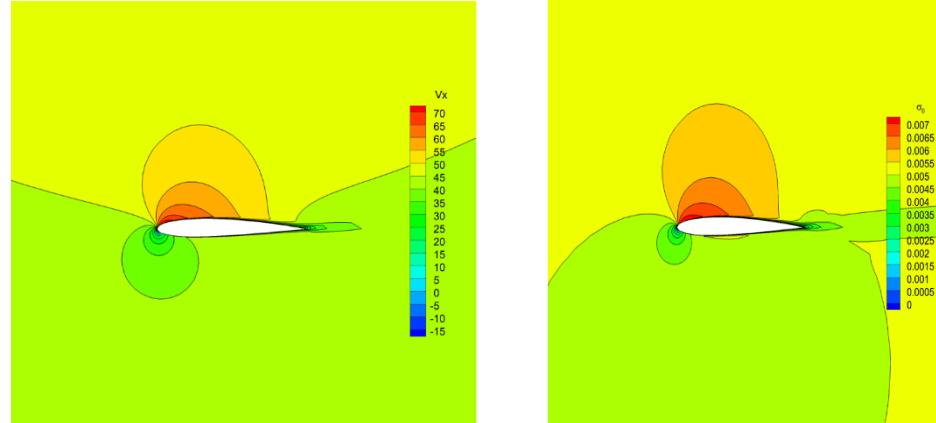
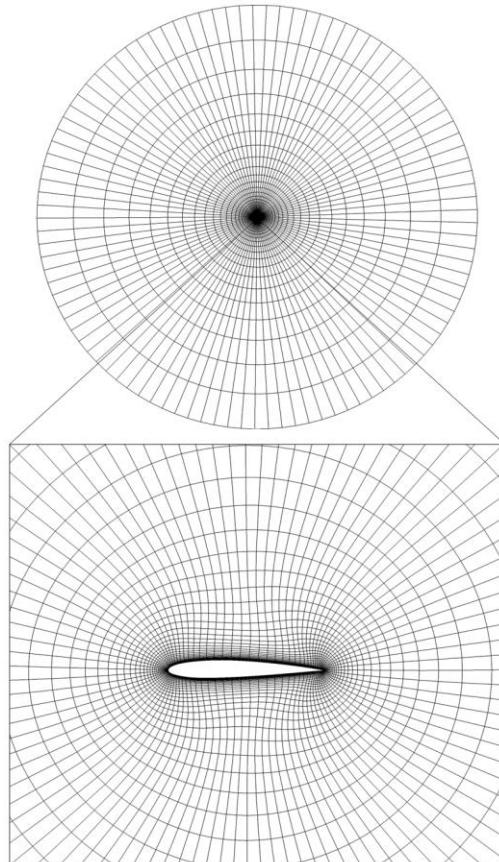


프레임워크 개발

■ 정확도 검증

■ NACA2412 익형 아음속 정상 유동 해석

- 입력 변수 : 자유류 받음각, 속도
- 5개 스냅샷 데이터로부터 신규 조건에 대한 해석 결과 추정
- 해석 격자, CFD / POD+ANN 비교, POD 모드 (2개 모드에 에너지 99.999% 분포)



프레임워크 개발

■ 정확도 검증

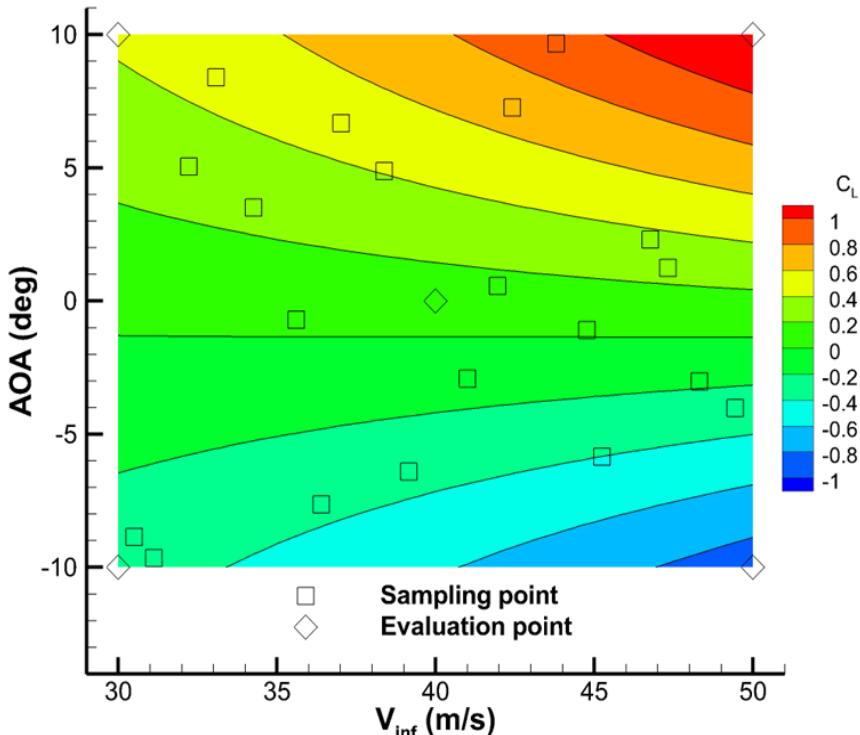
■ NACA2412 익형 아음속 정상 유동 해석

- POD 수행 과정 포함 시 소요 시간 5% 이내
- 기존에 있는 POD 모드로부터 보간만 수행할 경우 추가 단축 가능

Case No.	Method	V_∞ (m/s)	Angle of attack (deg)	Expansion coefficient		Calculati on Time (sec)	Error (%)
1	CFD	30	5	5864.55	-20.0794	5.60	-
2	CFD	40	0	7792.01	654.545		
3	CFD	40	4	7818.70	109.646		
4	CFD	40	10	7787.36	-707.902		
5	CFD	50	5	9774.26	-33.4612		
6	POD + ANN	45	7	8791.20	-302.488	0.27	4.98

프레임워크 개발

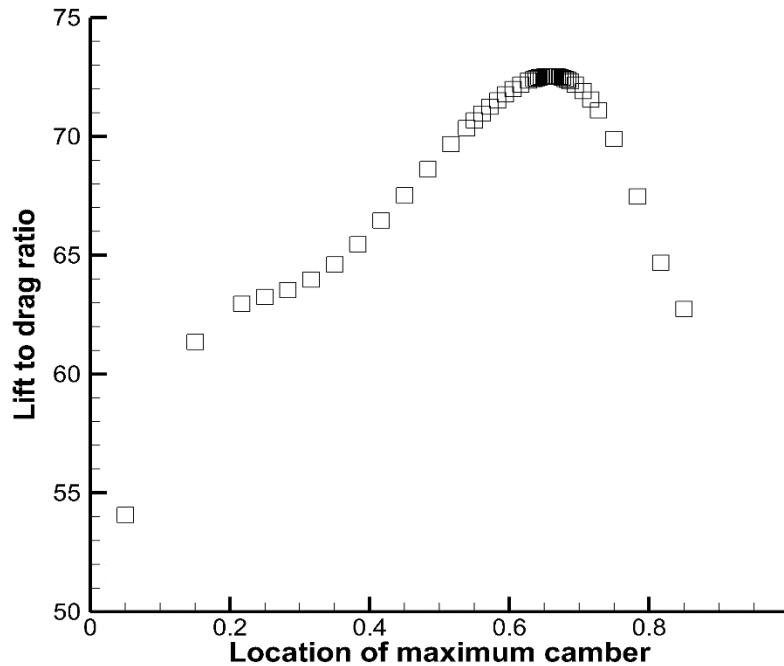
- 데이터베이스 생성 성능 검증
 - NACA2412 익형 아음속 정상 유동 해석
 - 자유류 속도 $30\text{m/s} \sim 50\text{m/s}$
 - 받음각 $-10^\circ \sim 10^\circ$
 - 20개의 입력 매개 변수 표본 사용
 - 5개 해석 결과 누적 시 POD 수행, 이후 해석 결과 초기 조건 추정
 - 소요 시간 : 1개 케이스 해석 소요 시간의 6.55배



Case No.	V_∞ (m/s)	AOA (deg)	Lift coefficient		Error (%)
			CFD	ANN	
1	30	-10	-0.321	-0.321	0.005
2	30	10	0.433	0.433	0.008
3	40	0	0.0965	0.0965	0.022
4	50	-10	-0.900	-0.908	0.878
5	50	10	1.220	1.211	0.790

프레임워크 개발

- 최적화 성능 검증
 - NACA2412 익형 아음속 정상 유동 해석
 - 자유류 속도 30m/s, 밭음각 8°
 - 양항비가 최대가 되는 최대 캠버 위치 탐색 (모핑된 격자 사용)
 - 소요 시간 : 1개 케이스 해석 소요 시간의 11배



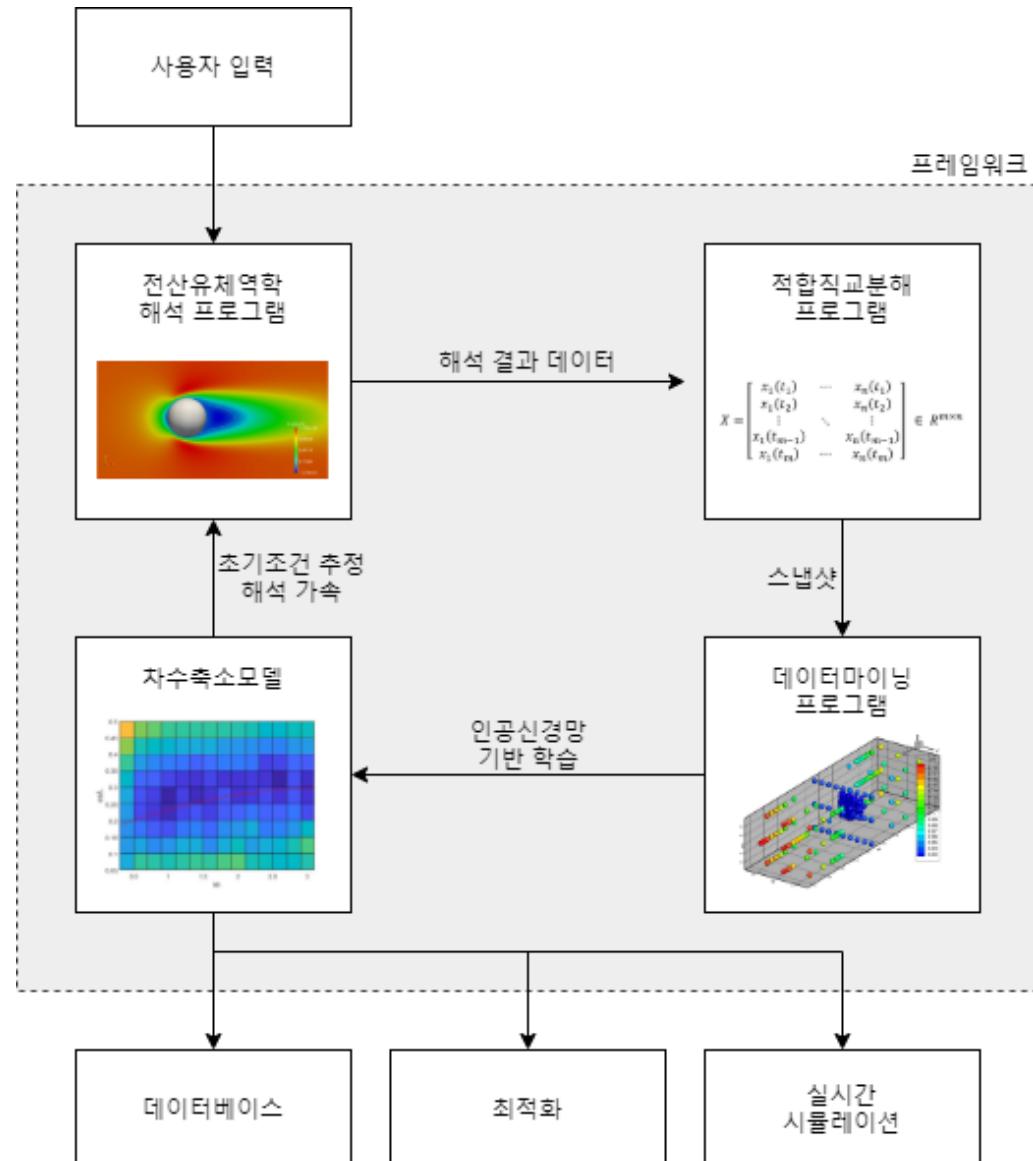
요약

■ 연구 성과

- 3개 프로그램을 연동/실행하는 프레임워크 개발 완료
- 프레임워크 정확도 및 성능 검증 완료
- 반복해석에 필요한 계산량을 용이하게 저감할 수 있는 기반 확보

■ 향후 진행 예정 사항

- 비정상 유동해석 결과에 대한 POD 수행
 - ITHACA-FV로 교체 검토
- 실시간 시뮬레이션 예제 구현



감사합니다