



Overview of the Stability workflow

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Numerical tools: Overview

- **IMAS:**

- Integrated Modelling & Analysis Suite

- **LIGKA[1]:**

- Linear gyrokinetic eigenvalue code

- **HELENA/CHEASE[2]:**

- MHD equilibrium solver

- **EP-Stability-WF:**

- Energetic Particle Stability Workflow (Python)

- **Models form hierarchy of fidelity, complexity:**

- Use local solvers to have an overview of the scenario before attempting global, more expensive runs.
- Use global solver to validate the results obtained by the local, faster runs.

Numerical tools: LIGKA



- **LIGKA:**

- Solves the linearized gyrokinetic equations -> eigenvalues and eigenfunctions (frequency, damping, mode structure).

- Models used in this work are (among others):

>1 s/mode

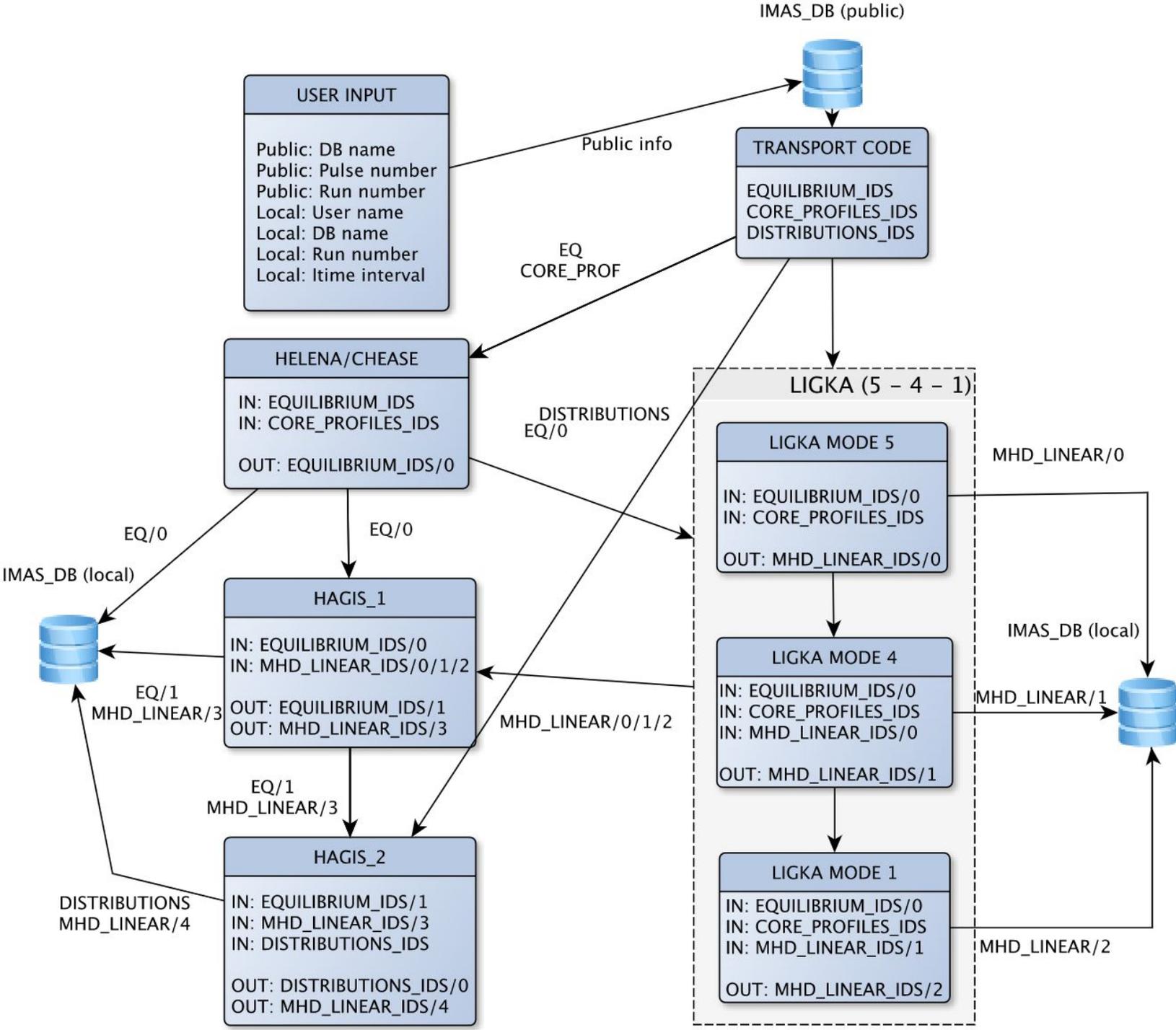
- Model 5: local analytical estimates of various basic AEs properties: frequency, estimated mode structure, rational surface, next and previous gap informations.

~ 10 s/mode

- Model 4: based on model 5 results, the local analytical dispersion relation for each mode is calculated. Determines the starting point for global calculations.

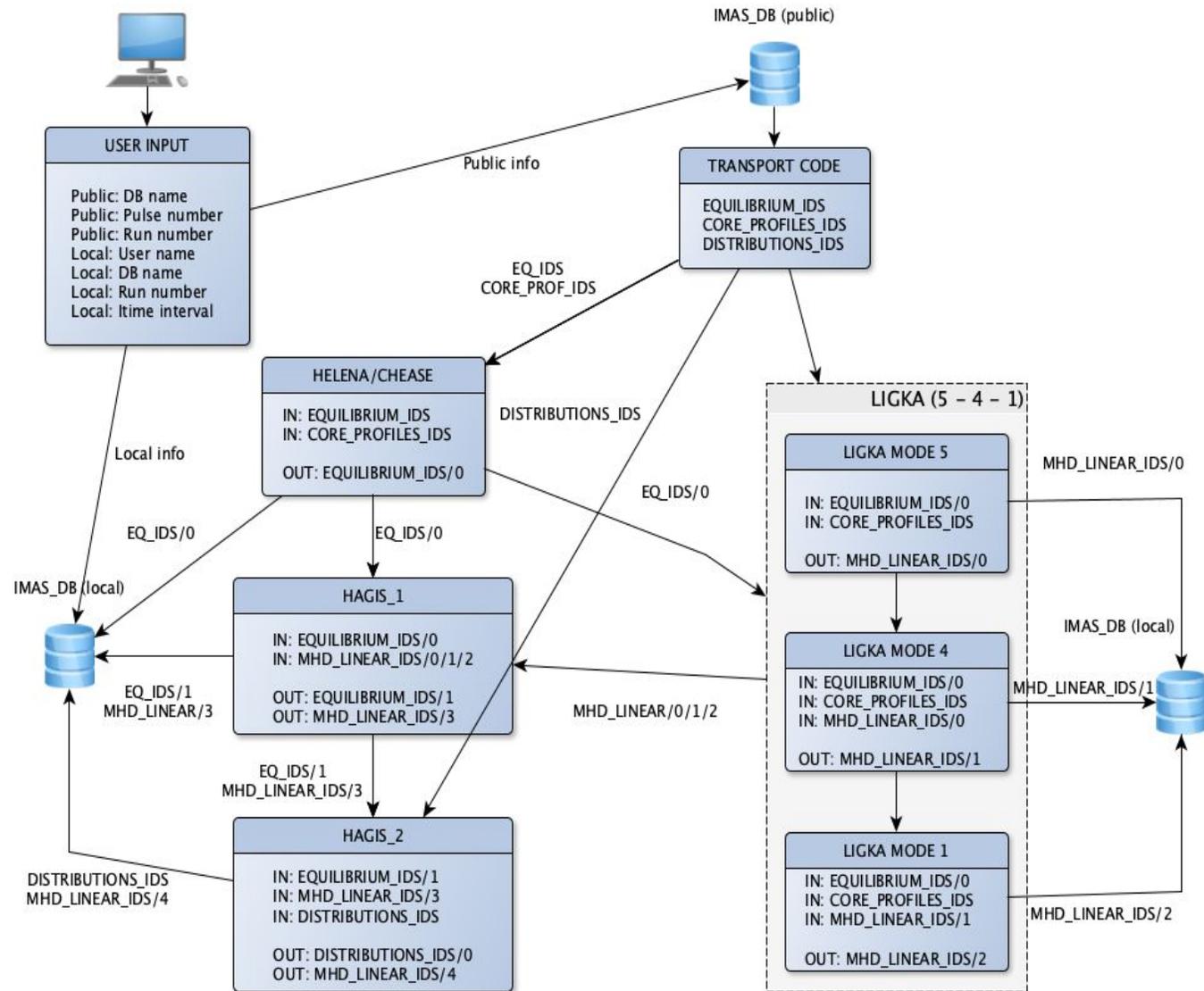
~ 30 min/mode

- Model 1: performs a frequency scan throughout the gap to find global linear properties of the modes.



Numerical tools: Energetic Particles Stability Workflow

- The aim of the WF is to perform an automated linear stability analysis on different time slices of a projected scenario or reconstructed experimental equilibrium.
- First time - dependent workflow which makes use of the IMAS infrastructure and various codes.
- **Scope:**
 - Connect the numerical tools with the data infrastructure (IMAS).
 - Facilitates retrieving/saving data from the DB through XML files.
 - Fast configuration of numerical tools.
 - Complete data analysis suite integrated in the interface. (to be completed)



Flow of FC2K actors in the WF



```
def actor_settings(actor):
    actor_params = {}
    input_ids = {}
    output_ids = {}
    if actor == "Chease":
        actor_params["entrypoint_actor"] = True
        actor_params["wrapper"] = chease_actor_wf_wrapper
        actor_params["config_file_name"] = "chease_input_choices.xml"
        input_ids = {"equilibrium": 0, "core_profiles": 0}
        output_ids = {"equilibrium": 2, "core_profiles": 0}
    if actor == "Helena":
        actor_params["entrypoint_actor"] = True
        actor_params["wrapper"] = helena_actor_wf_wrapper
        actor_params["config_file_name"] = "helena.xml"
        input_ids = {"equilibrium": 0, "core_profiles": 0}
        output_ids = {"equilibrium": 0, "core_profiles": 0}
    if actor == "Ligka_m5":
        actor_params["entrypoint_actor"] = False
        actor_params["wrapper"] = ligka_actor_wf_wrapper
        actor_params["config_file_name"] = "z_ligka.xml"
        input_ids = {"equilibrium": 0, "core_profiles": 0}
        output_ids = {"mhd_linear": 0}
    if actor == "Ligka_m4":
        actor_params["entrypoint_actor"] = False
        actor_params["wrapper"] = ligka_actor_wf_wrapper
        actor_params["config_file_name"] = "z_ligka.xml"
        input_ids = {"equilibrium": 0, "core_profiles": 0, "mhd_linear": 0}
        output_ids = {"mhd_linear": 1}
    if actor == "Ligka_m1":
        actor_params["entrypoint_actor"] = False
        actor_params["wrapper"] = ligka_actor_wf_wrapper
        actor_params["config_file_name"] = "z_ligka.xml"
        input_ids = {"equilibrium": 0, "core_profiles": 0, "mhd_linear": 1}
        output_ids = {"mhd_linear": 2}
    if actor == "Ligka_m6":
        actor_params["entrypoint_actor"] = False
        actor_params["wrapper"] = ligka_actor_wf_wrapper
        actor_params["config_file_name"] = "z_ligka.xml"
        input_ids = {"equilibrium": 0, "core_profiles": 0, "mhd_linear": 0}
        output_ids = {"mhd_linear": 5}
    if actor == "Ligka_m2":
        actor_params["entrypoint_actor"] = False
        actor_params["wrapper"] = ligka_actor_wf_wrapper
        actor_params["config_file_name"] = "z_ligka.xml"
        input_ids = {"equilibrium": 0, "core_profiles": 0, "mhd_linear": 2}
        output_ids = {"mhd_linear": 6}
    if actor == "Ligka_m3":
        actor_params["entrypoint_actor"] = False
        actor_params["wrapper"] = ligka_actor_wf_wrapper
        actor_params["config_file_name"] = "z_ligka.xml"
        input_ids = {"equilibrium": 0, "core_profiles": 0, "mhd_linear": 0}
        output_ids = {"mhd_linear": 7}
    if actor == "Hagis_1":
        actor_params["entrypoint_actor"] = False
        actor_params["wrapper"] = hagis1_actor_wf_wrapper
        actor_params["config_file_name"] = "hagis1.xml"
        input_ids = {"equilibrium": 0, "mhd_linear": 0}
        output_ids = {"equilibrium": 1, "mhd_linear": 3}
    if actor == "Hagis_2":
        actor_params["entrypoint_actor"] = False
        actor_params["wrapper"] = hagis2_actor_wf_wrapper
        actor_params["config_file_name"] = "hagis2.xml"
        input_ids = {"equilibrium": 1, "mhd_linear": 3, "core_profiles": 0}
        output_ids = {"distributions": 0, "mhd_linear": 4}
    if actor == "Finder":
        actor_params["entrypoint_actor"] = False
        actor_params["wrapper"] = finder_actor_wf_wrapper
        actor_params["config_file_name"] = "finder_input.xml"
        input_ids = {"equilibrium": 1}
        output_ids = {"distributions": 1}
    if actor == "Falcon":
        actor_params["entrypoint_actor"] = False
        actor_params["wrapper"] = falcon_actor_wf_wrapper
        actor_params["config_file_name"] = "falcon_input.xml"
        input_ids = {"equilibrium": 2}
        output_ids = {"mhd_linear": 5}

    actor_params["input_ids"] = input_ids
    actor_params["output_ids"] = output_ids

    return actor_params
```

Fc2k actors:

- 1) Have callable functions (wrappers to the fortran, or python code (Falcon))
- 2) Input only IDSs
- 3) Output only IDSs
- 4) Easier to manage input/output and decide order of execution

EP WORKFLOW

WORKFLOW PARAMETERS

user: public
 machine: ITER
 shot_nr: 130012
 run_in: 2
 machine_out: test_DB
 run_out: 10
 itime: 15-17,19

FURTHER SETTINGS

ligka_541
 ligka_5412
 pulse_list
 fast_particles
 hdf5
 mpi_processes: 8

Save Configuration Save and Run
 Save Configuration as Load Configuration
 Restore Default

Scenario Summary Choice

ACTOR SELECTION

Equilibrium_code_chease: 0
 Equilibrium_code: Helena
 Distributions_1: 0
 Distributions_2: 0
 Orbit_Finder: 0
 Stability_code: Ligka_m5

CHEASE Parameters
 HELENA Parameters
 LIGKA Parameters
 HAGIS 1 Parameters
 HAGIS 2 Parameters
 FINDER Parameters
 Species Settings
 SCENARIO Parameters
 IDS Merge

SPECIES SETTINGS

Bulk Ions

H: 0.02
 D: 0.02
 T: 0.02

Impurities

Be: 0.02
 Ne: 0.02
 He4: 0.02
 C: 0.02
 Tu: 0.02
 Ar: 0.02

Fast Ions

H: 0.001
 D: 0.001
 He4: 0.001

Save Species Configuration

SCENARIO PARAMETERS (m)

n_e: 1
 n_H: 1
 n_D: 1
 n_T: 1
 n_Be: 1
 n_C: 1
 n_Ne: 1
 n_He4_ash: 1
 n_He4_EP: 1
 T_e: 1
 T_H: 1
 T_D: 1
 T_T: 1
 T_Be: 1
 T_C: 1
 T_Ne: 1
 T_He4_ash: 1
 T_He4_EP: 1

Save SCENARIO Configuration

IDS Merge

Inputs

user_in_1: public
 machine_in_1: ITER
 shot_in_1: 130012
 run_in_1: 2
 HDF5_1:
 user_in_2: public
 machine_in_2: ITER
 shot_in_2: 130012
 run_in_2: 2
 HDF5_2:

Settings

itime: 15-17,19
 Equilibrium_copy:
 ne:
 Te:
 ni_H:
 Ti_H:
 ni_T:
 Ti_T:
 ni_D:
 Ti_D:
 ni_Be:
 Ti_Be:
 ni_C:
 Ti_C:
 ni_Ne:
 Ti_Ne:

Output

machine_out: TEST_IDS_MERGE
 shot_out: 130012
 run_out: 89
 HDF5_out:

Save IDS_MERGE Configuration

LIGKA PARAMETERS

modus: 5
 min_n_tor: 10
 max_n_tor: 10
 min_m: 11
 max_m: 11
 sidebands: 5
 sidebands_asy: 2
 mode_type: 1
 even: 0
 cocp: 1
 start_pos: 1
 force_m: false
 npsi_out: 256
 kr_read: 0.0d0
 q0: 0.0d0
 rad_start: 0.0d0
 rad_end: 1.0d0
 offset_d: 0.0d0

Save LIGKA Configuration



Extra features



SPECIES SETTINGS

Bulk Ions

H

D

T

Impurities

Be

Ne

He4

C

Tu

Ar

Fast Ions

H

D

He4

Save Species Configuration

SCENARIO PARAMETERS

SCENARIO PARAMETERS (multipliers)

n_e	<input type="text" value="1"/>
n_H	<input type="text" value="1"/>
n_D	<input type="text" value="1"/>
n_T	<input type="text" value="1"/>
n_Be	<input type="text" value="1"/>
n_C	<input type="text" value="1"/>
n_Ne	<input type="text" value="1"/>
n_He4_ash	<input type="text" value="1"/>
n_He4_EP	<input type="text" value="1"/>
T_e	<input type="text" value="1"/>
T_H	<input type="text" value="1"/>
T_D	<input type="text" value="1"/>
T_T	<input type="text" value="1"/>
T_Be	<input type="text" value="1"/>
T_C	<input type="text" value="1"/>
T_Ne	<input type="text" value="1"/>
T_He4_ash	<input type="text" value="1"/>
T_He4_EP	<input type="text" value="1"/>

Save SCENARIO Configuration

Scenario Selector

ulse	Run	Database	Reference	Ip[MA]	B0[T]	Fuelling	Confinement	Workflow
100002	1	ITER	ITER-half-field-H	-7.5	-2.65	H	L-mode	METIS
100001	2	ITER	ITER-full-field-H	-15.0	-5.3	H	L-mode	METIS
100003	1	ITER	ITER-third-field-H	-5.0	-1.8	H	L-H-L	METIS
100007	1	ITER	ITER-intermediate-3T-H	-8.5	-3.0	H	L-H-L	METIS
100008	1	ITER	ITER-intermediate-3.3T-H	-9.5	-3.3	H	L-H-L	METIS
100009	1	ITER	ITER-intermediate-4.5T-H	-12.5	-4.5	H	L-mode	METIS
100013	1	ITER	ITER-PFPO1-1.8T-H	-5.0	-1.8	H	L-H-L	METIS
100015	1	ITER	ITER-PFPO2-1.8T-H-0.9*n_GW-NBI_745keV_22.3I	-5.0	-1.8	H	L-H-L	METIS
100014	2	ITER	ITER-PFPO2-1.8T-H-0.5*n_GW-NBI_530keV_9.4M	-5.0	-1.8	H	L-H-L	METIS
100016	1	ITER	ITER-10MA-5.3T-Hydrogen	-10.0	-5.3	H	L-mode	METIS

Development/testing cycle, current and previous versions



- **Maintenance cycle of actors + WF:**
 - **Actors are self contained codes that can act independently or as part of a workflow.**
 - **They are continuously tested and maintained via versions (different modules in sdcc/gw)**
 - **On top of that we have the EP-Stability-WF integrated testing.**
 - **When testing the wf, we also test the integration of LIGKA + HELENA/CHEASE inside the WF (2x testing for actors)**
 - **Testing happens automatically at every push of every piece of code (via automated bamboo tests)**

- **Following runs were performed with:**

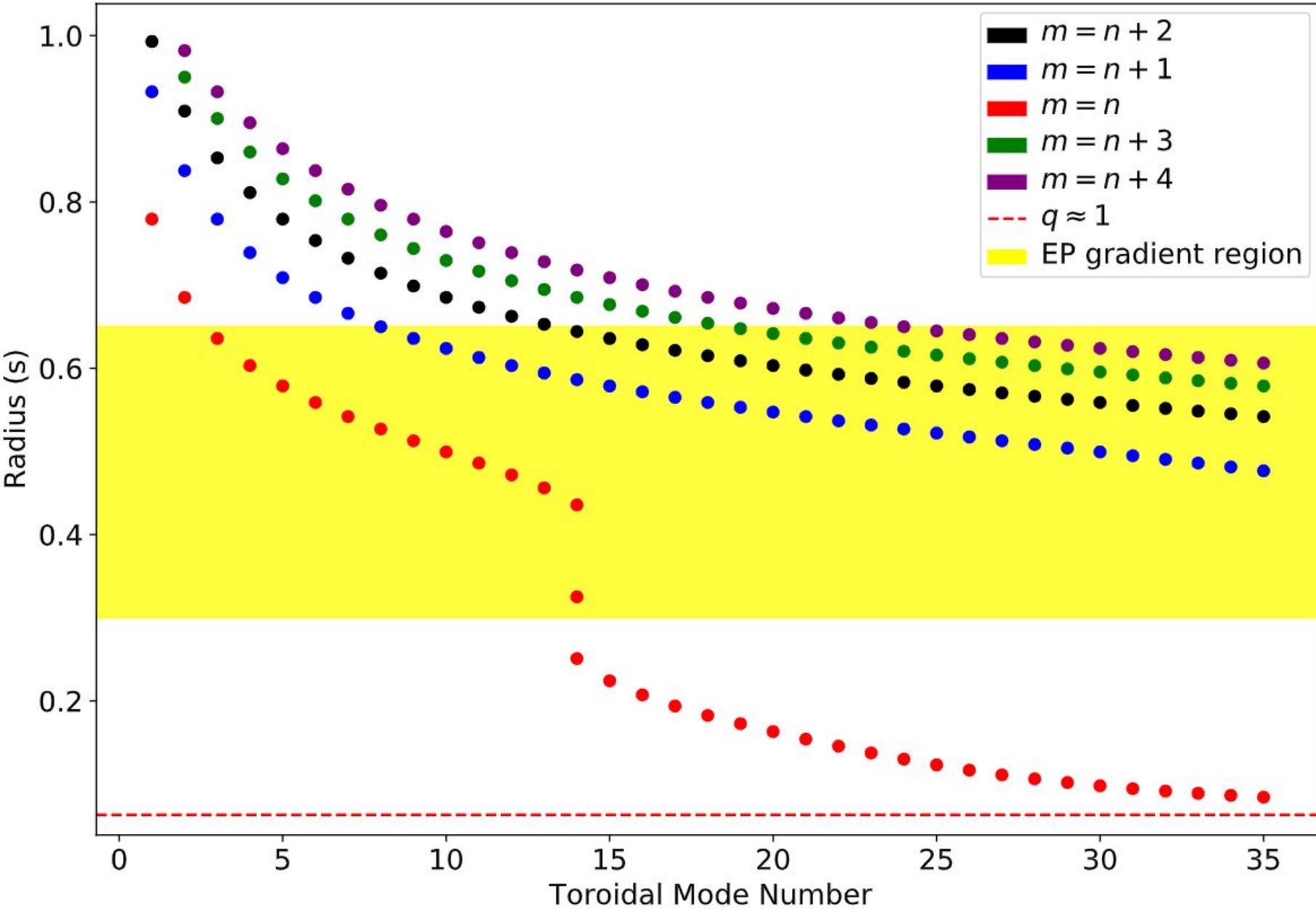
- **HELENA: 2.0.1-intel-2020b-DD-3.35.0**
- **LIGKA: 1.0.1-intel-2020b-DD-3.35.0**
- **CHEASE: 1.0.9-intel-2020b-DD-3.35.0**
- **EP-WF: 1.0.2-intel-2020b-DD-3.35.0**

- **Current status of codes:**

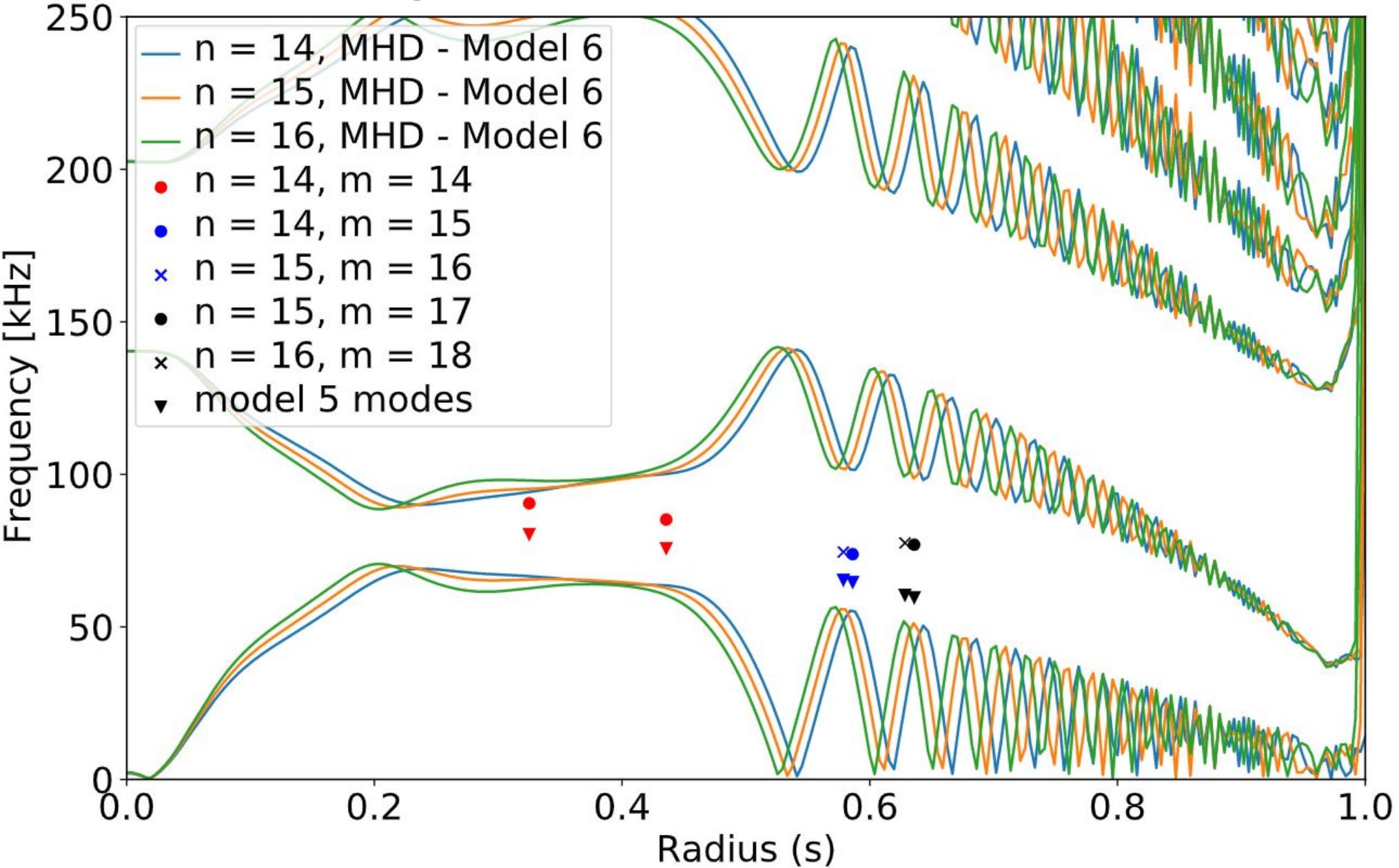
- **HELENA: 2.0.1-intel-2020b-DD-3.37.0**
- **LIGKA: 1.0.5-intel-2020b-DD-3.37.0**
- **CHEASE: 13.1.2-intel-2020b-DD-3.37.0**
- **EP-WF: 1.0.4-intel-2020b-DD-3.37.0**

Things are moving fast!!

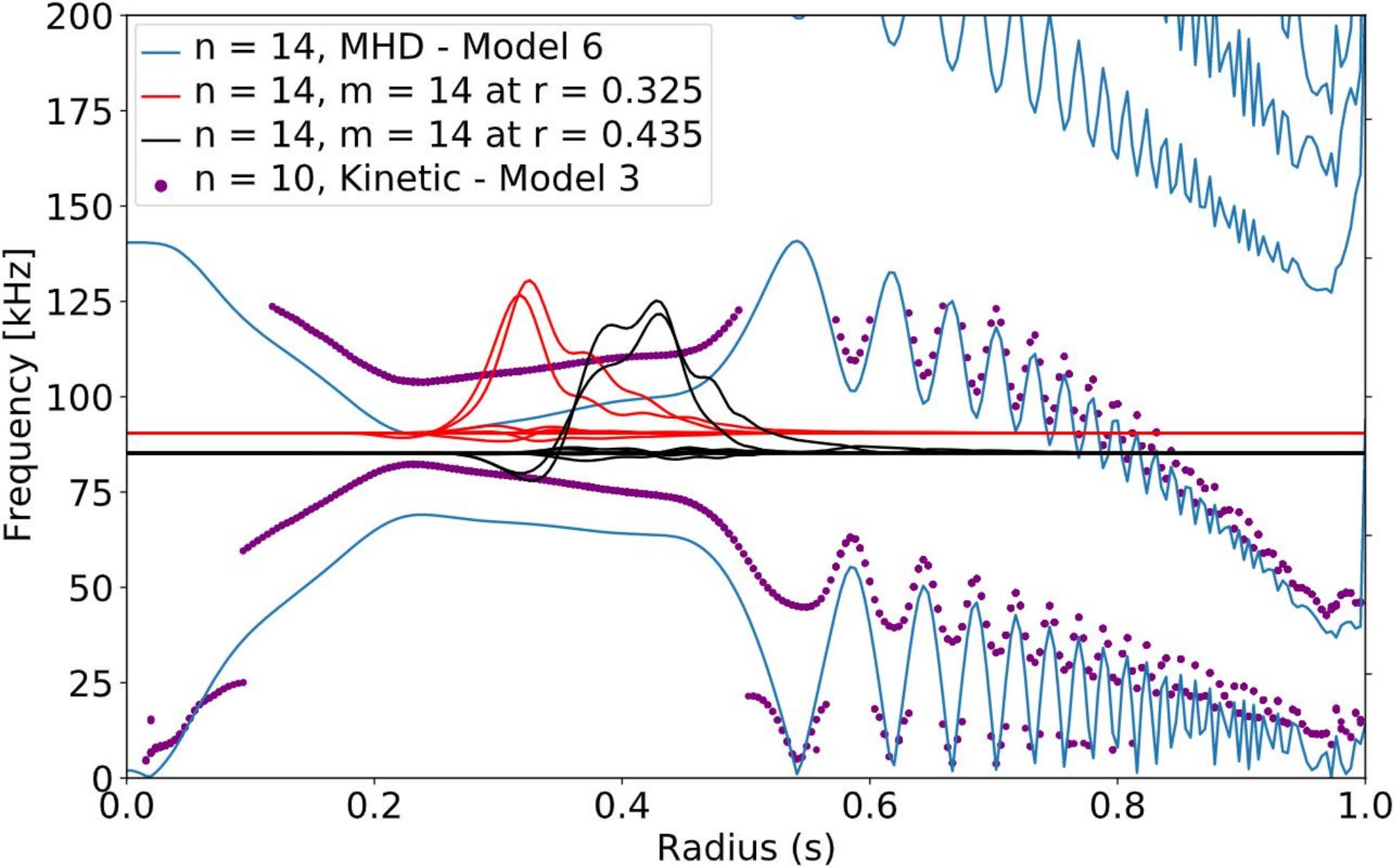
Scenario 1: ASTRA - 131025/34 ITER DB - model 5



Model 6 + comparison between model 5-4-1

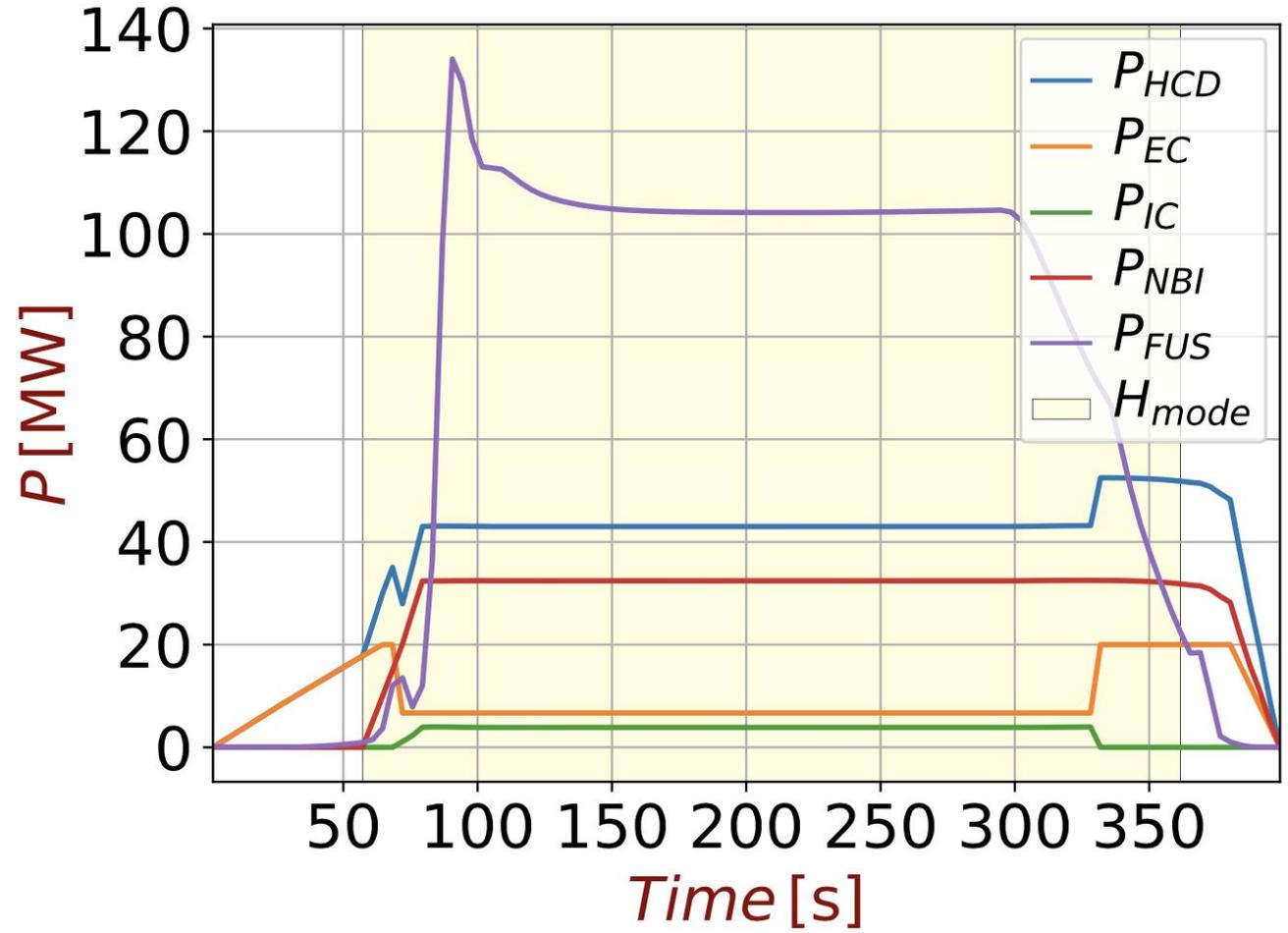
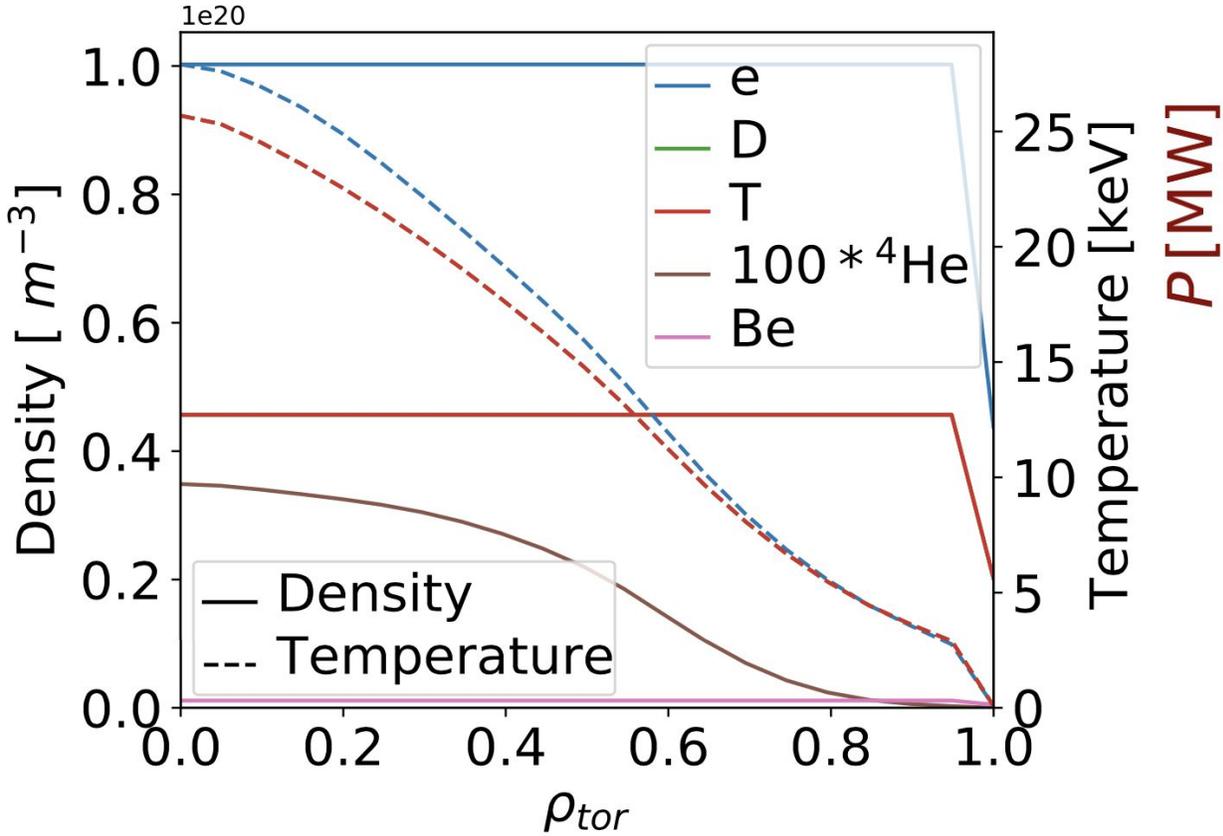


Model 6 + 3 + model 1 global EFs.

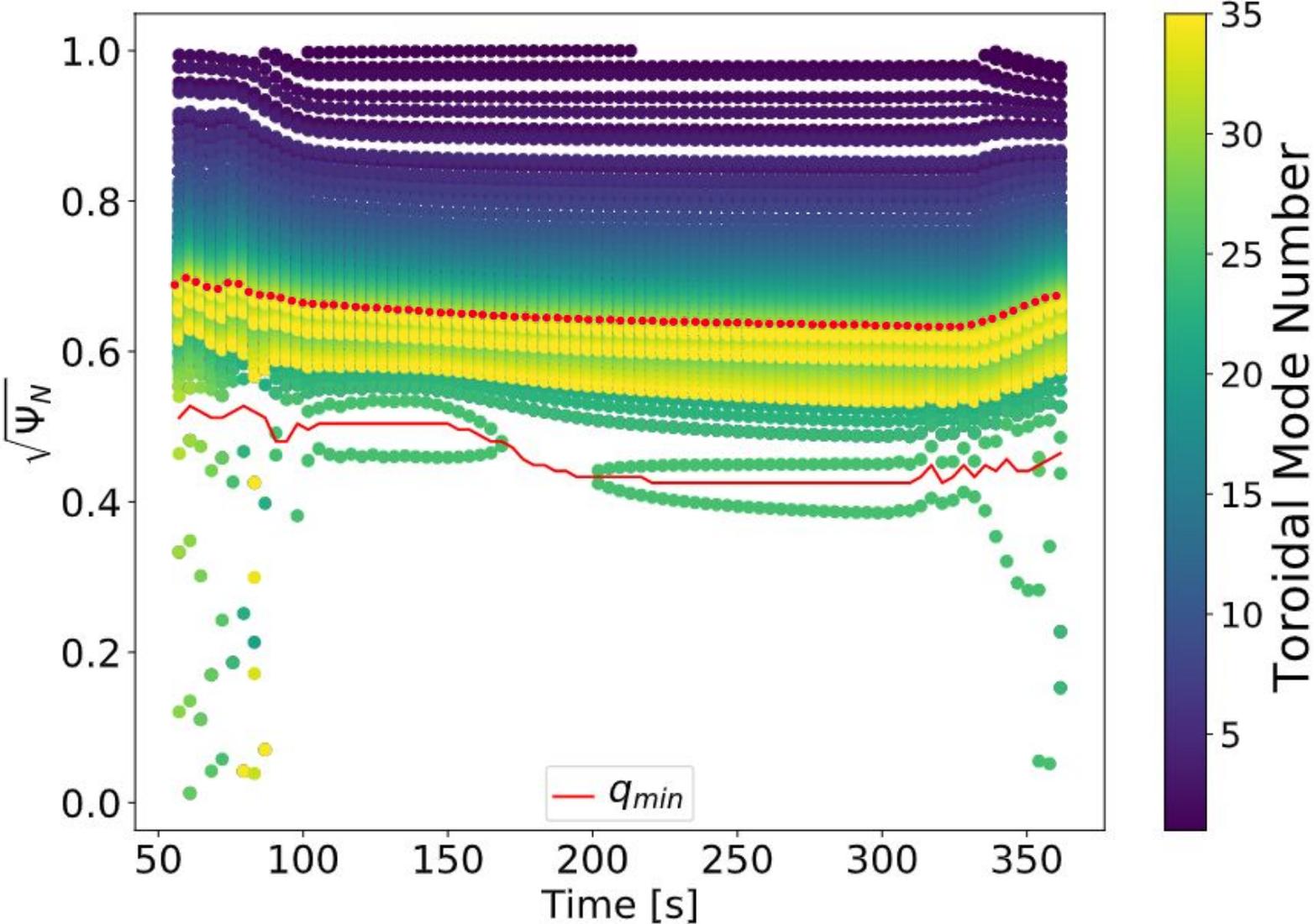


Scenario: METIS time-dependant Q=10 ITER baseline

- D-T plasma, $Q = 10$, 15 MA
- Peak axis $T_{e,0} \cong 27$ keV, 4 keV pedestal top
- $T_{D/T,0} \cong 25$ keV
- $n_{e,0} \cong 10^{20} \text{ m}^{-3}$, $n_{D/T,0} \cong 0.45 \times 10^{20} \text{ m}^{-3}$

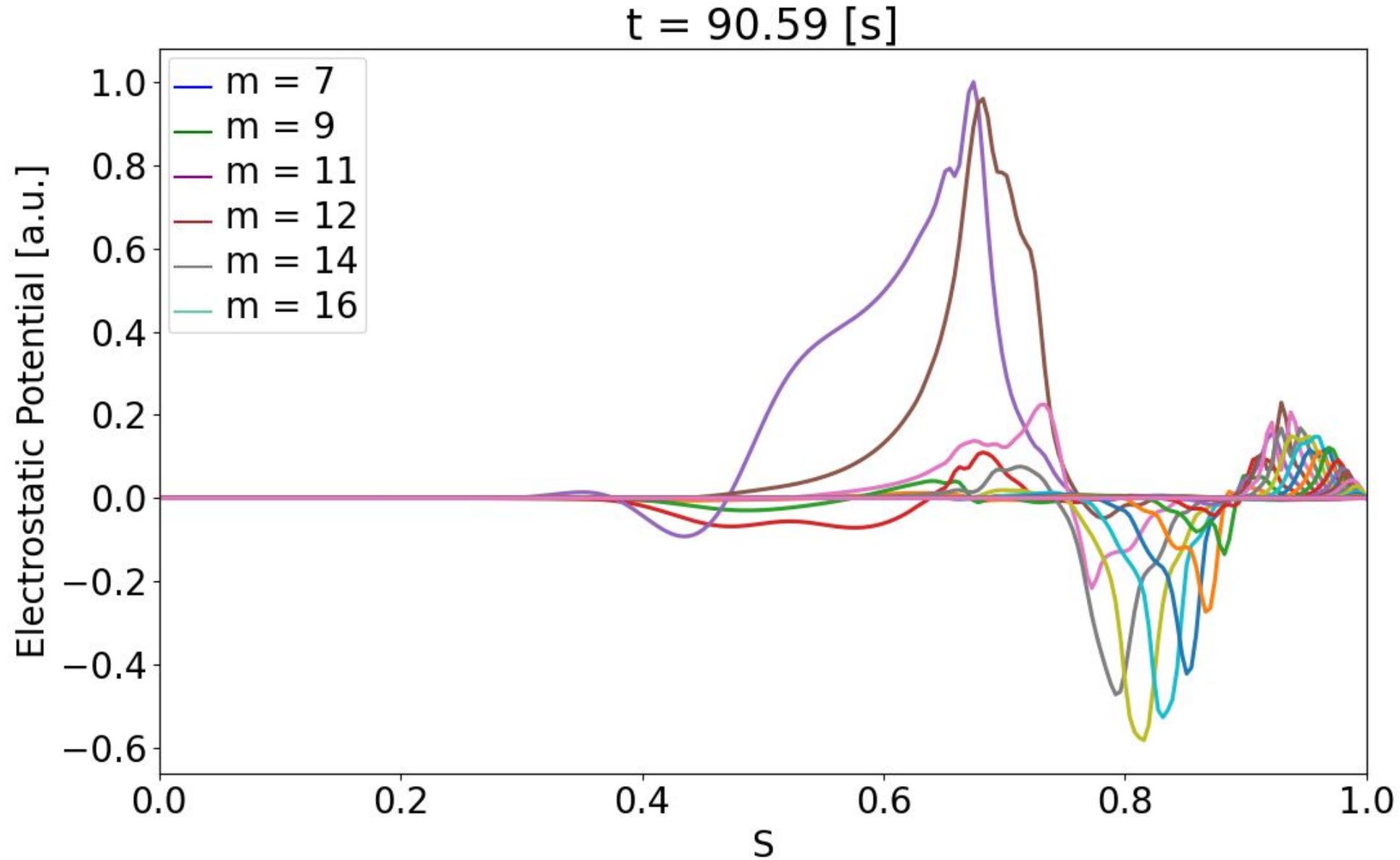


Results: Scan over the entire Radial/Temporal domain

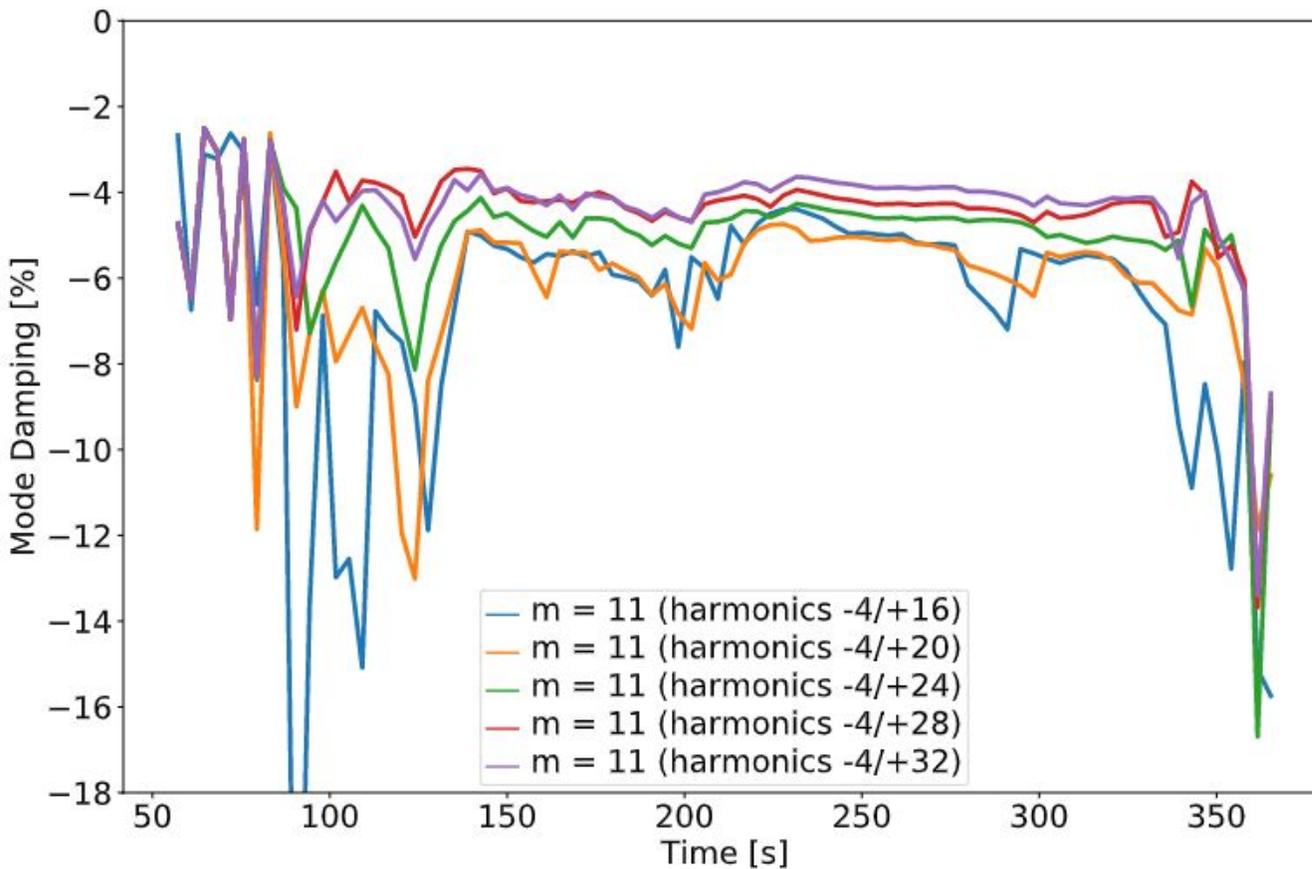


- $n = 1 \dots 35$
- $m = n + 1 \dots m = n + 5$
- 12014 total number of modes, 137/time-point
- With red $n = 10, m = 11$ mode
- Slightly inverted q -profile around $s = 0.5$ leads to two different TAE branches with the same mode numbers.
- After 90s no more TAEs are found in the core, due to small magnetic shear in the core assumed by the METIS-given equilibrium.

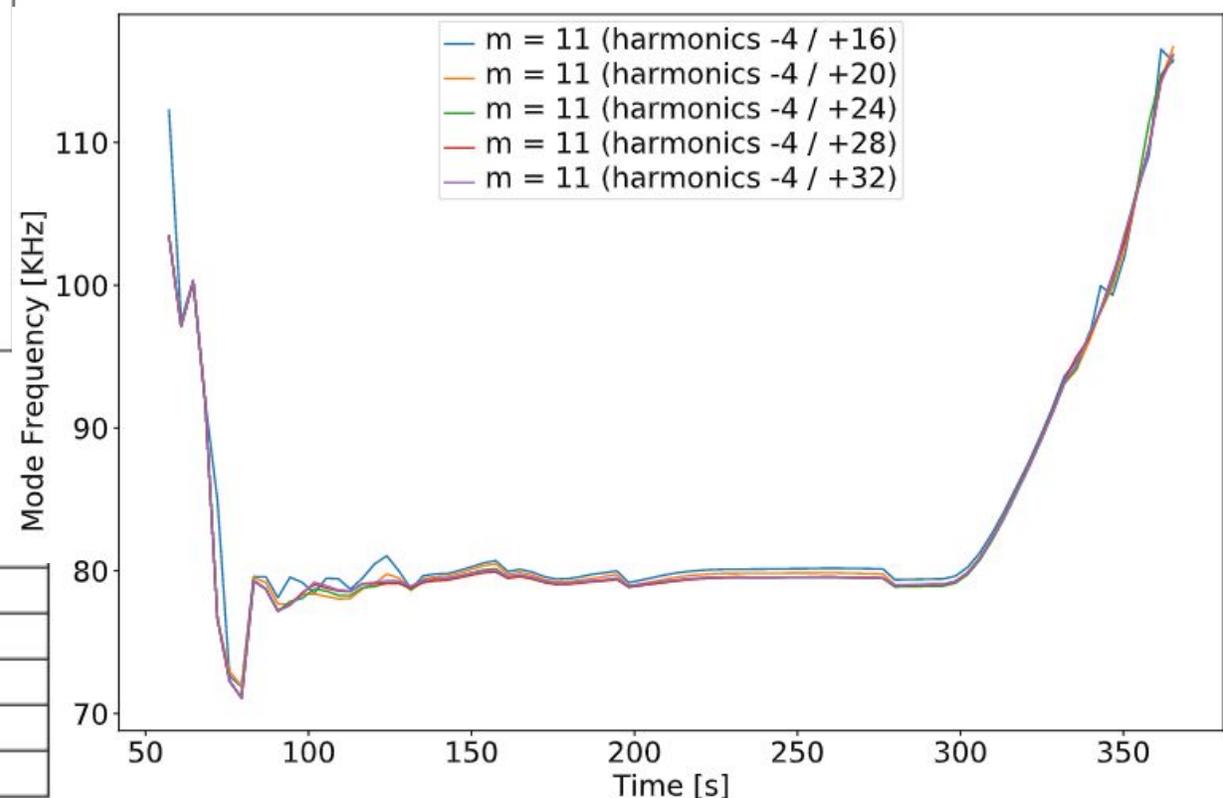
Results: TAE $n = 10$, $m = (11,12)$ global mode structure



Results: Convergence test (n = 10)

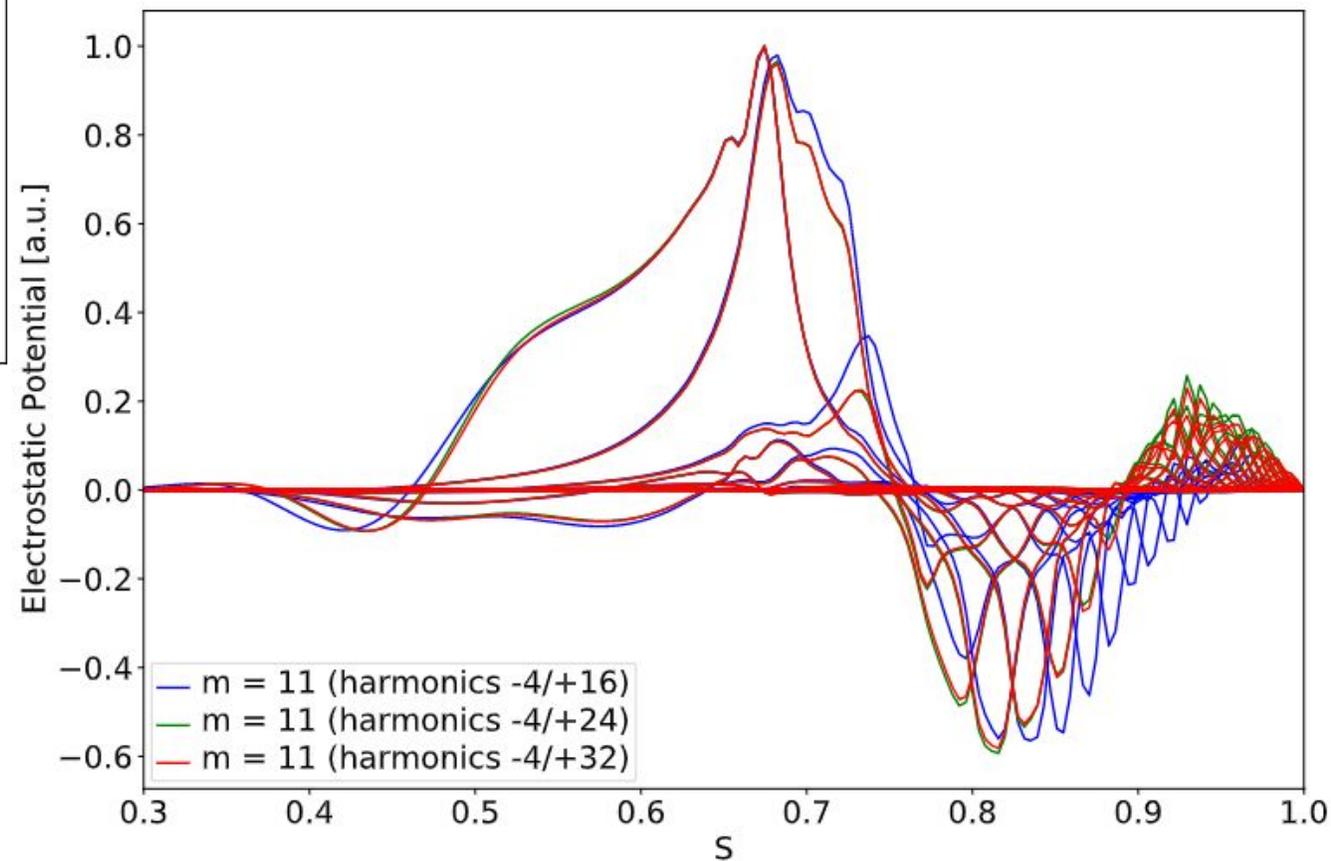
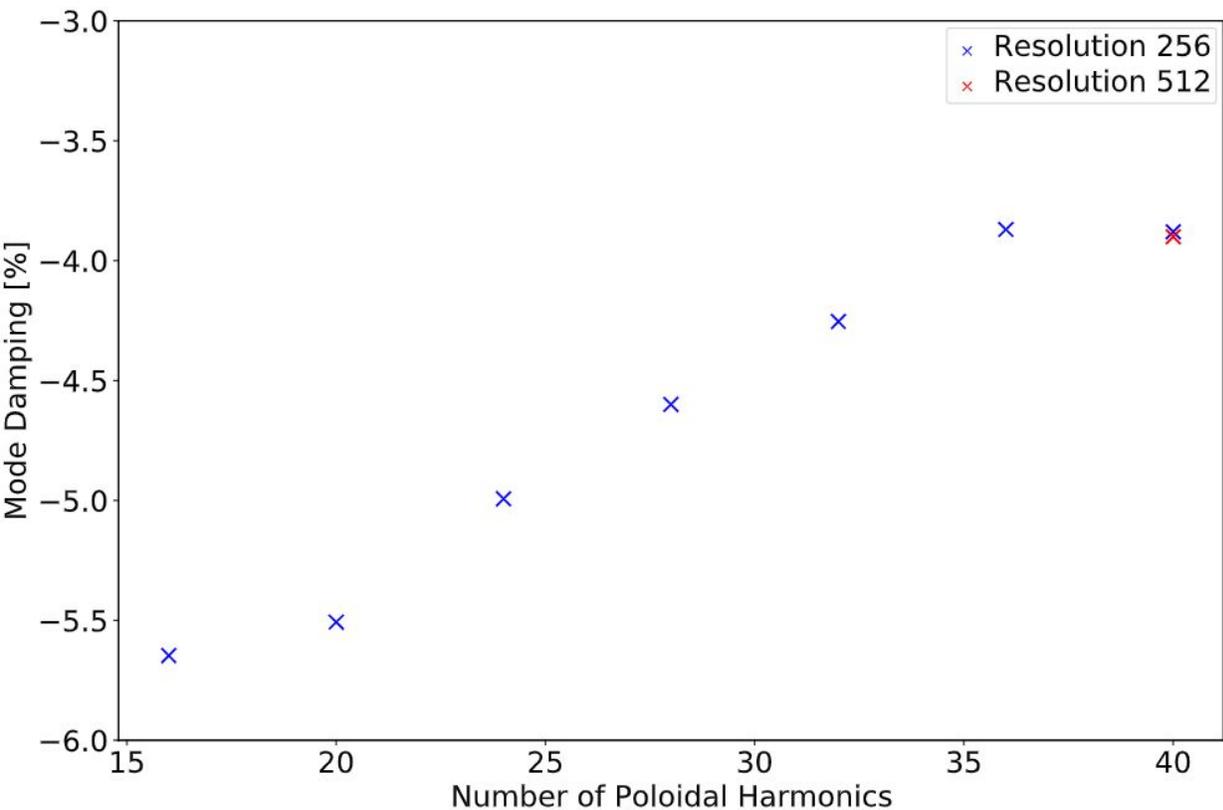


METIS	local scan	106	12014	1
METIS	local + global (-4/+16)	106	1	local = 1, global = 8
METIS	local + global (-4/+20)	106	1	local = 1, global = 8
METIS	local + global (-4/+28)	106	1	local = 1, global = 8
METIS	local + global (-4/+32)	106	1	local = 1, global = 8



METIS	local scan	01:00:24	00:00:34	<1s
METIS	local + global (-4/+16)	04:36:41	00:02:36	-
METIS	local + global (-4/+20)	06:20:02	00:03:35	-
METIS	local + global (-4/+28)	10:31:06	00:05:56	-
METIS	local + global (-4/+32)	13:00:35	00:07:23	-

Results: Convergence test - mode structure/extra points



Results: EP influence on TAE, n = 10 mode

