

final

December 6, 2025

0.1 Final Project

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[1]: import sympy as sp
import numpy as np
from scipy.linalg import expm
from scipy.io import loadmat
from scipy.integrate import solve_ivp
import matplotlib.pyplot as plt

# =====
# Provided Nominal Values
# =====

# State
xi_g_nom      = 10.0
eta_g_nom     = 0.0
theta_g_nom   = float(sp.pi/2)
xi_a_nom      = -60.0
eta_a_nom     = 0.0
theta_a_nom   = float(-sp.pi/2)
x_nom = np.array([
    xi_g_nom,
    eta_g_nom,
    theta_g_nom,
    xi_a_nom,
    eta_a_nom,
    theta_a_nom
], dtype=float)
n = x_nom.shape[0] # state dimensions

#Input
v_g_nom      = 2.0
phi_g_nom    = float(-sp.pi/18)
v_a_nom      = 12.0
omega_a_nom  = float(sp.pi/25)
u_nom = np.array([
    v_g_nom,
    phi_g_nom,
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    v_a_nom,
    omega_a_nom
], dtype=float)
m = u_nom.shape[0] # input dimensions

L_nom = 0.5

# =====
#      Setup Sympy
# =====

# State and Input
xi_g, eta_g, theta_g, xi_a, eta_a, theta_a = sp.symbols('xi_g eta_g theta_g' +
    'xi_a eta_a theta_a', real=True)
v_g, phi_g, v_a, omega_a = sp.symbols('v_g phi_g v_a omega_a', real=True)

# State and input vectors
x_sym = sp.Matrix([xi_g, eta_g, theta_g, xi_a, eta_a, theta_a])
u_sym = sp.Matrix([v_g, phi_g, v_a, omega_a])

# Parameters
L_sym = sp.symbols('L', positive=True)

# =====
#      Setup Dynamics
# =====

# State
f1 = v_g * sp.cos(theta_g)
f2 = v_g * sp.sin(theta_g)
f3 = (v_g / L_sym) * sp.tan(phi_g)
f4 = v_a * sp.cos(theta_a)
f5 = v_a * sp.sin(theta_a)
f6 = omega_a
f_sym = sp.Matrix([f1, f2, f3, f4, f5, f6])

# Measurements
delta_xi = xi_a - xi_g
delta_eta = eta_a - eta_g
h1 = sp.atan2(delta_eta, delta_xi) - theta_g # Azimuth of UAV from UGV, ↵relative to UGV heading
h2 = sp.sqrt(delta_xi**2 + delta_eta**2) # Distance between UGV/UAV
h3 = sp.atan2(-delta_eta, -delta_xi) - theta_a # Azimuth of UGV from UAV, ↵relative to UAV heading
h4 = xi_a # UAV easting position
h5 = eta_a # UAV northing position

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h_sym = sp.Matrix([h1, h2, h3, h4, h5])
p = h_sym.shape[0] # measurement dimensions

# Symbolic Jacobians for A, B, C
A_sym = f_sym.jacobian(x_sym) # df/dx - 6x6
B_sym = f_sym.jacobian(u_sym) # df/du - 6x4
C_sym = h_sym.jacobian(x_sym) # dh/dx - 5x6

# =====
# Define functions from Sympy
# =====

A_eval = sp.lambdify(
    (x_sym, u_sym, L_sym),
    A_sym,
    "numpy"
)

B_eval = sp.lambdify(
    (x_sym, u_sym, L_sym),
    B_sym,
    "numpy"
)

C_eval = sp.lambdify(
    (x_sym,),
    C_sym,
    "numpy"
)

# Function that will evaluate our NL CT dynamics for given x and u
f_eval = sp.lambdify(
    (x_sym, u_sym, L_sym),
    f_sym,
    "numpy"
)

def f_ct(x_vec, u_vec, L_val=0.5):
    """
    Wrapper for f_eval sympy lambda.
    Computes continuous-time nonlinear dynamics xdot = f(x,u).
    """

    x_vec: shape (6,) numpy array [xi_g, eta_g, theta_g, xi_a, eta_a, theta_a]
    u_vec: shape (4,) numpy array [v_g, phi_g, v_a, omega_a]
    L_val: scalar (e.g. 0.5)

    Returns: shape (6,) numpy array xdot

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"""
# Ensure 1D arrays
x_vec = np.asarray(x_vec, dtype=float).reshape(-1)
u_vec = np.asarray(u_vec, dtype=float).reshape(-1)

xdot = f_eval(x_vec, u_vec, L_val) # this returns a 6x1-like object
return np.array(xdot, dtype=float).reshape(6,)

h_eval = sp.lambdify(
    (x_sym,),
    h_sym,
    "numpy"
)

def h_meas(x_vec):
    """
    Wrapper for h_eval sympy lambda.
    linear measurement model  $y = h(x)$ .

    x_vec: shape (6,) numpy array [xi_g, eta_g, theta_g, xi_a, eta_a, theta_a]

    Returns: shape (5,) numpy array y
    """
    x_vec = np.asarray(x_vec, dtype=float).reshape(-1)

    y = h_eval(x_vec)
    return np.array(y, dtype=float).reshape(5,)

# =====
# Evaluate at nominal operating point
# =====

# Get ready to sub back into our symbolic expressions
subs_nom = {
    xi_g: xi_g_nom,
    eta_g: eta_g_nom,
    theta_g: theta_g_nom,

    xi_a: xi_a_nom,
    eta_a: eta_a_nom,
    theta_a: theta_a_nom,

    v_g: v_g_nom,
    phi_g: phi_g_nom,
    v_a: v_a_nom,
    omega_a: omega_a_nom,
}

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L_sym: L_nom
}

A_nom_sym = A_sym.subs(subs_nom)
B_nom_sym = B_sym.subs(subs_nom)
C_nom_sym = C_sym.subs(subs_nom)

print("\nA_nom (linearized df/dx at nominal point):")
sp pprint(sp.simplify(A_nom_sym))

print("\nB_nom (linearized df/du at nominal point):")
sp pprint(sp.simplify(B_nom_sym))

print("\nC_nom (linearized dh/dx at nominal point):")
sp pprint(sp.simplify(C_nom_sym))

# Optional: turn into numeric matrices (floats) if you want
A = np.array(sp.N(A_nom_sym), dtype=float)
B = np.array(sp.N(B_nom_sym), dtype=float)
C = np.array(sp.N(C_nom_sym), dtype=float)

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A_nom (linearized df/dx at nominal point):
0 0 -2.0 0 0 0
0 0 1.22464679914735e-16 0 0 0
0 0 0 0 0 0
0 0 0 0 0 12.0
0 0 0 0 0 7.34788079488412e-16
0 0 0 0 0 0

B_nom (linearized df/du at nominal point):
6.12323399573677e-17 0 0 0
1.0 0 0 0
-0.35265396141693 4.12436481650305 0 0
0 0 6.12323399573677e-17 0
0 0 -1.0 0
0 0 0 1

```

```
C_nom (linearized dh/dx at nominal point):
 0  0.0142857142857143 -1  0   -0.0142857142857143  0
 1.0      0      0   -1.0      0      0
 0  0.0142857142857143  0  0   -0.0142857142857143 -1
 0      0      0   1      0      0
 0      0      0   0      1      0
```

```
[2]: def discretize_ct_system(A, B, C, dt):
    """
    Given continuous-time linearized matrices A, B, C and a timestep dt,
    return the discrete-time linearized system matrices F, G, H using ZOH
    ↴matrix exponential
    """
    n = A.shape[0]
    m = B.shape[1]

    AB = np.hstack((A, B))
    A_hat = np.vstack((AB, np.zeros((m, n+m))))
    Z = expm(A_hat*dt)

    F = Z[:n, :n]
    G = Z[:n, n:]
    H = C.copy()
    return F, G, H

def discretize_ct_forward_euler(A, B, C, dt):
    n = A.shape[0]
    I = np.eye(n)

    # Forward-Euler discretization
    F = I + dt * A      # State transition matrix
    G = dt * B          # Input matrix
    H = C.copy()         # Measurement matrix (instantaneous, so same form)

    return F, G, H

def check_stability(F, tol=1e-9):
    """
    Print eigenvalues of F and classify stability (discrete-time).
    """
    eigvals = np.linalg.eigvals(F)
    mags = np.abs(eigvals)
```

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print("Eigenvalues of F:")
for lam, mag in zip(eigvals, mags):
    print(f"    = {lam:.6f}, || = {mag:.6f}")

if np.all(mags < 1 - tol):
    print("\nStability: asymptotically stable (all || < 1).")
elif np.any(mags > 1 + tol):
    print("\nStability: unstable (some || > 1).")
else:
    print("\nStability: marginal (at least one || = 1, none clearly > 1).")

def observability_matrix(F, H):
    """
    Build the observability matrix:
    O = [ H
          H F
          H F^2
          ...
          H F^{n-1} ]
    where n is the state dimension.
    """
    n = F.shape[0]
    O_blocks = []
    F_power = np.eye(n)

    for _ in range(n):
        O_blocks.append(H @ F_power)
        F_power = F @ F_power

    return np.vstack(O_blocks)

def check_observability(F, H, tol=1e-9):
    """
    Compute and print the rank of the observability matrix.
    """
    O = observability_matrix(F, H)
    rank = np.linalg.matrix_rank(O, tol=tol)
    n = F.shape[0]

    print(f"Observability matrix shape: {O.shape}")
    print(f"Rank(O) = {rank} (state dimension = {n})")

    if rank == n:
        print("System is locally observable around the linearization point.")
    else:
        print("System is NOT fully observable around the linearization point.")

```

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dt = 0.1 # seconds

# 1) Discretize
F, G, H = discretize_ct_system(A, B, C, dt)

print("F (DT state transition matrix):")
print(F)
print("\nG (DT input matrix):")
print(G)
print("\nH (DT measurement matrix):")
print(H)

# 2) Stability analysis
print("\n==== Stability check ====")
check_stability(F)

# 3) Observability analysis
print("\n==== Observability check ====")
check_observability(F, H)

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F (DT state transition matrix):
[[ 1.0000000e+00  0.0000000e+00 -2.0000000e-01  0.0000000e+00
   0.0000000e+00  0.0000000e+00]
 [ 0.0000000e+00  1.0000000e+00  1.22464680e-17  0.0000000e+00
   0.0000000e+00  0.0000000e+00]
 [ 0.0000000e+00  0.0000000e+00  1.0000000e+00  0.0000000e+00
   0.0000000e+00  0.0000000e+00]
 [ 0.0000000e+00  0.0000000e+00  0.0000000e+00  1.0000000e+00
   0.0000000e+00  1.2000000e+00]
 [ 0.0000000e+00  0.0000000e+00  0.0000000e+00  0.0000000e+00
   1.0000000e+00  7.34788079e-17]
 [ 0.0000000e+00  0.0000000e+00  0.0000000e+00  0.0000000e+00
   0.0000000e+00  1.0000000e+00]]

G (DT input matrix):
[[ 3.52653961e-03 -4.12436482e-02  0.0000000e+00  0.0000000e+00
   1.0000000e-01  2.52544509e-18  0.0000000e+00  0.0000000e+00
  [-3.52653961e-02  4.12436482e-01  0.0000000e+00  0.0000000e+00]
 [ 0.0000000e+00  0.0000000e+00  6.12323400e-18  6.0000000e-02]
 [ 0.0000000e+00  0.0000000e+00 -1.0000000e-01  3.67394040e-18]
 [ 0.0000000e+00  0.0000000e+00  0.0000000e+00  1.0000000e-01]]]

H (DT measurement matrix):
[[ 0.          0.01428571 -1.          0.          -0.01428571  0.          ]
 [ 1.          0.          0.          -1.          0.          0.          ]
 [ 0.          0.01428571  0.          0.          -0.01428571 -1.          ]
 [ 0.          0.          0.          1.          0.          0.          ]]

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[ 0.          0.          0.          0.          1.          0.        ]]

==== Stability check ====
Eigenvalues of F:
= 1.000000, || = 1.000000
= 1.000000, || = 1.000000
= 1.000000, || = 1.000000
= 1.000000, || = 1.000000
= 1.000000, || = 1.000000
= 1.000000, || = 1.000000

Stability: marginal (at least one || < 1, none clearly > 1).

==== Observability check ====
Observability matrix shape: (30, 6)
Rank(0) = 6 (state dimension = 6)
System is locally observable around the linearization point.

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[3]: # =====
# Initial conditions and perturbation
# =====
data = loadmat("cooplocalization_finalproj_KFdata.mat")

# Keep things generalized
tvec = data["tvec"].ravel()
N_steps = tvec.size # 1001
dt = float(tvec[1] - tvec[0]) # 0.1 sec
t_span = (0.0, dt * (N_steps - 1)) # (0, 100)
t_eval = np.linspace(t_span[0], t_span[1], N_steps)

# Randomly chosen initial perturbation
delta_x0 = np.array([
    0.1,                      # +0.1 m in xi_g
    -0.1,                     # -0.1 m in eta_g
    np.deg2rad(1.0),           # +1 deg in theta_g
    0.2,                       # +0.2 m in xi_a
    0.0,                        # 0 m in eta_a
    np.deg2rad(-1.0)           # -1 deg in theta_a
])

# =====
# Nonlinear CT simulation (truth)
# =====
def ode_rhs(t, x_flat):
    return f_ct(x_flat, u_nom)
```

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# Use solve_ivp to get DT solution
x0_nl = x_nom + delta_x0
nl_sim = solve_ivp(
    fun=ode_rhs,
    t_span=t_span,
    y0=x0_nl,
    t_eval=t_eval,
    rtol=1e-9,
    atol=1e-9
)
x_traj_nl = nl_sim.y.T # (N_steps, 6)
y_traj_nl = np.array([h_meas(x) for x in x_traj_nl]) # (N_steps, 5)

# =====
# Nominal Nonlinear simulation
# =====
x_traj_nom = np.zeros((N_steps, n)) # array of column vectors
x_traj_nom[0, :] = x_nom # no perturbation in nominal

# Using forward Euler to get DT results without linearizing
for k in range(N_steps - 1):
    x_traj_nom[k+1, :] = x_traj_nom[k, :] + dt * f_ct(x_traj_nom[k, :], u_nom)
y_traj_nom = np.array([h_meas(x) for x in x_traj_nom]) # Will be used in part 2

# =====
# Linearized DT Simulation
# =====
delta_x_lin = delta_x0.copy()

# Storage for state and measurement trajectories
x_traj_lin = np.zeros_like(x_traj_nom)
y_traj_lin = np.zeros_like(y_traj_nl)

# Precompute time-varying linearization (A_k, C_k, F_k, H_k)
F_list = []
H_list = []
C_list = []

# Loop through each nominal state, linearize, discretize, and evaluate
for k in range(N_steps):
    # Nominal state computed above
    x_nom_k = x_traj_nom[k, :]

    # Linearize around current nominal point
    A_k = np.array(A_eval(x_nom_k, u_nom, L_nom), dtype=float)

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B_k = np.array(B_eval(x_nom_k, u_nom, L_nom), dtype=float)
C_k = np.array(C_eval(x_nom_k), dtype=float)

# Discretize to DT (time-varying F, G, H)
F_k, G_k, H_k = discretize_ct_forward_euler(A_k, B_k, C_k, dt)

# Save these for Part II
F_list.append(F_k)
H_list.append(H_k)
C_list.append(C_k)

# Total linearized state & measurement at this step
x_traj_lin[k, :] = x_nom_k + delta_x_lin
y_traj_lin[k, :] = y_traj_nom[k] + C_k @ delta_x_lin

# Propagate perturbation (no input perturbation)
if k < N_steps - 1:
    delta_x_lin = F_k @ delta_x_lin

# =====
# Plot results
# =====

state_labels = [
    r"\xi_g [m]",
    r"\eta_g [m]",
    r"\theta_g [rad]",
    r"\xi_a [m]",
    r"\eta_a [m]",
    r"\theta_a [rad]",
]
]

fig1, axes1 = plt.subplots(3, 2, figsize=(12, 8))
axes1 = axes1.ravel()

for i in range(6):
    ax = axes1[i]
    ax.plot(t_eval, x_traj_nl[:, i], label="Nonlinear", linewidth=2)
    ax.plot(t_eval, x_traj_lin[:, i], "--", label="Linearized DT", linewidth=2)
    ax.set_ylabel(state_labels[i])
    ax.grid(True)
    if i == 0:
        ax.legend(loc="best")

axes1[-1].set_xlabel("Time [s]")
axes1[-2].set_xlabel("Time [s]")
fig1.suptitle("State Trajectories: Nonlinear vs Linearized DT", fontsize=14)

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fig1.tight_layout(rect=[0, 0.03, 1, 0.95])

meas_labels = [
    r"Azimuth UGV to UAV [rad]",
    r"Range [m]",
    r"Azimuth UAV to UGV [rad]",
    r" $\xi_a$  GPS [m]",
    r" $\eta_a$  GPS [m],
]

fig2, axes2 = plt.subplots(3, 2, figsize=(12, 8))
axes2 = axes2.ravel()

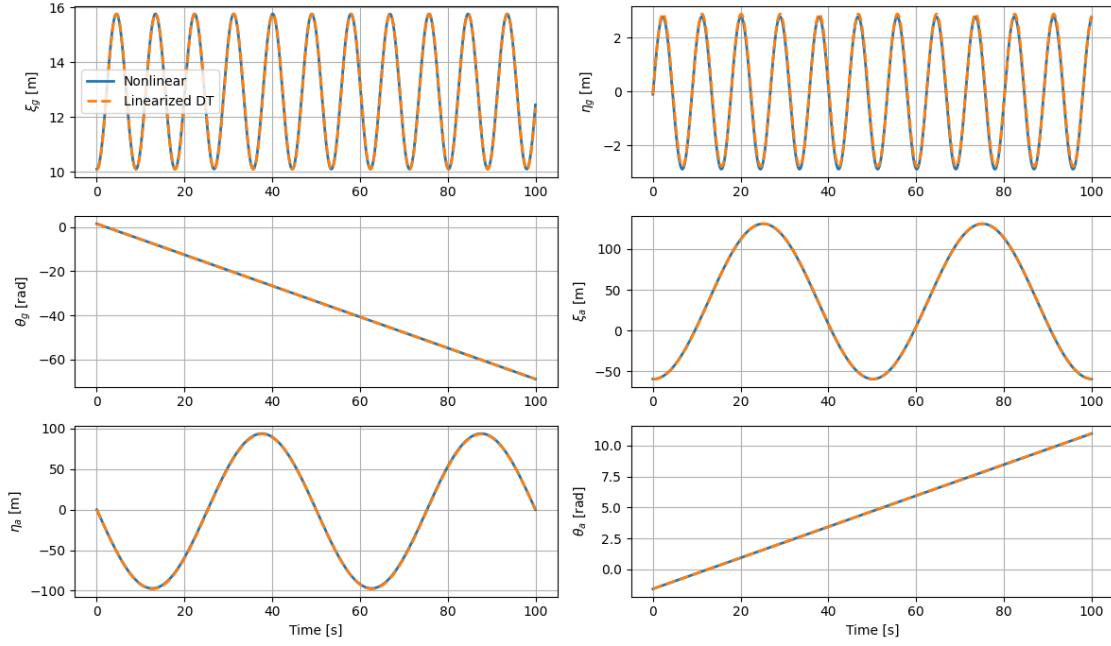
for i in range(5):
    ax = axes2[i]
    ax.plot(t_eval, y_traj_nl[:, i], label="Nonlinear", linewidth=2)
    ax.plot(t_eval, y_traj_lin[:, i], "--", label="Linearized DT", linewidth=2)
    ax.set_ylabel(meas_labels[i])
    ax.grid(True)
    if i == 0:
        ax.legend(loc="best")

axes2[-1].axis("off")
axes2[-2].set_xlabel("Time [s]")
axes2[-3].set_xlabel("Time [s]")
fig2.suptitle("Measurement Trajectories: Nonlinear vs Linearized DT", fontsize=14)
fig2.tight_layout(rect=[0, 0.03, 1, 0.95])

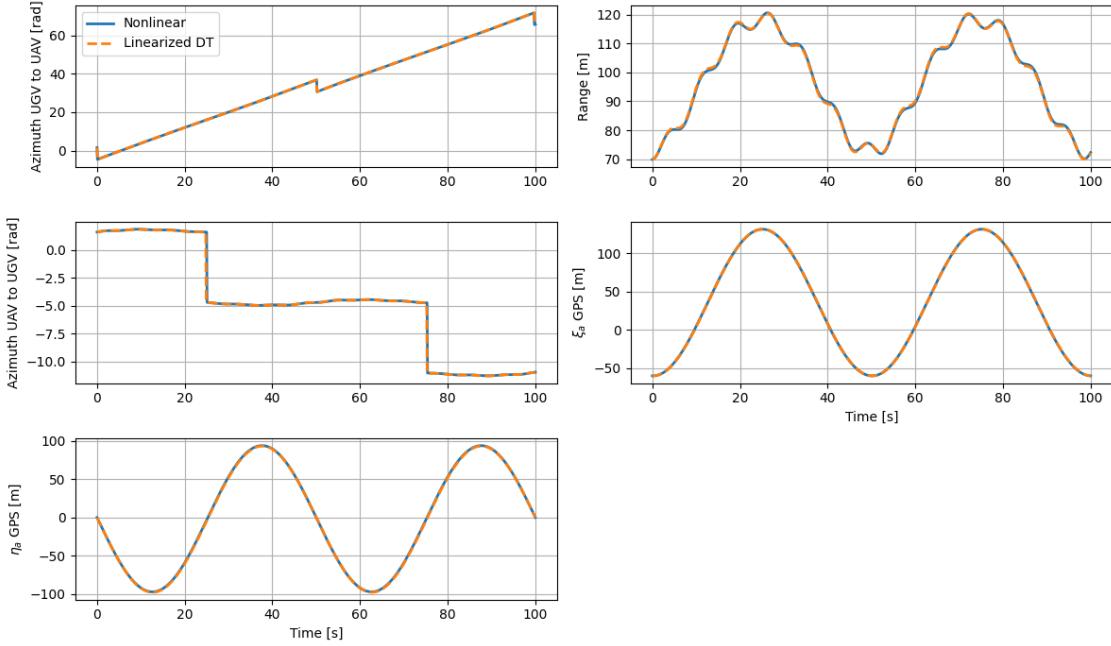
plt.show()

```

State Trajectories: Nonlinear vs Linearized DT



Measurement Trajectories: Nonlinear vs Linearized DT



The Linearized KF that is linearized around a nominal trajectory loops through the nominal tra-

jectory values x_{nom} and linearizes the dynamics around the current nominal state value $x_{nom}(k)$. This contrasts with the EKF, which linearizes around the current state estimate x_k^+

Now we are tasked with implementing a Kalman filter. Recall the equations we use for each prediction step:

$$\begin{aligned}\hat{x}_{k+1}^- &= F\hat{x}_k^+ + Gu_k \\ P_{k+1}^- &= FP_k^+F^T + Q \\ K_{k+1} &= P_{k+1}^- H_{k+1}^T [H_{k+1}P_{k+1}^-H_{k+1}^T + R]^{-1}\end{aligned}$$

When a new measurement comes in we do

$$\begin{aligned}\hat{x}_{k+1}^+ &= \hat{x}_{k+1}^- + K_{k+1}(y_{k+1} - H_{k+1}\hat{x}_{k+1}^-) \\ P_{k+1}^+ &= (I - K_{k+1}H_{k+1})P_{k+1}^-\end{aligned}$$

Since we are doing perturbation mechanics here, all of our x values in the equation above are now δx . For example in our LKF loop, one of the first things we do is $\delta\hat{x}_{k+1}^- = F_k\delta\hat{x}_k^+ + G_k\delta u_k$. Since we have no input/control perturbation, δu_k is 0. So $\delta\hat{x}_{k+1}^- = F_k\delta\hat{x}_k^+$. Our measurement's also use the perturbation dynamics. The equation for \hat{y}_{k+1}^- is $\hat{y}_{k+1}^- = y_{nom}(k) + H_k\delta\hat{x}_k^-$

```
[4]: # =====
#   Initial Setup
# =====

# Load the cov matrices from .mat data
Q_true = data["Qtrue"] # 6x6
R_true = data["Rtrue"] # 5x5

# Non global version of random.seed()
rng = np.random.default_rng(42)

# Filter's initial mean and cov - we are told to keep these the same for each run
# Will use them to generate a random ground truth x(0) on each run
delta_x0_mean = np.zeros(n)
P0_plus = np.diag([
    5.0,      # std dev in xi_g [m]
    5.0,      # std dev in eta_g [m]
    np.deg2rad(5.0), # std dev in theta_g [rad]
    10.0,     # std dev in xi_a [m]
    10.0,     # std dev in eta_a [m]
    np.deg2rad(5.0)  # std dev in theta_a [rad]
])**2 # ^2 makes it variances

def simulate_truth_trajectory(
    Q, R,
    delta_x0_mean, P0_plus,
    x_traj_nom,      # only x_traj_nom[0] is used here
```

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        u_nom, L_nom, dt,
        rng=None
):
    """
    Generate ONE nonlinear truth trajectory + noisy measurements for TMT.

    Returns:
        - x_true: (N_steps, n)
        - y_meas: (N_steps, p)
    """
    if rng is None:
        rng = np.random.default_rng()
    N_steps, n = x_traj_nom.shape
    p = R.shape[0]

    # =====
    # Truth simulation (nonlinear + noises)
    # =====

    # Sample true initial perturbation and build true initial state
    delta_x0_true = rng.multivariate_normal(mean=delta_x0_mean, cov=P0_plus)
    x_true = np.zeros((N_steps, n))
    x_true[0, :] = x_traj_nom[0, :] + delta_x0_true
    w = rng.multivariate_normal(mean=np.zeros(n), cov=Q, size=N_steps)  # ↳process noise
    for k in range(N_steps - 1):
        x_true[k+1, :] = x_true[k, :] + dt * f_ct(x_true[k, :], u_nom, L_nom) + w[k]

    # Simulate noisy measurements from nonlinear h(x) with R_true
    y_meas = np.zeros((N_steps, p))
    y_meas[0, :] = np.nan  # t=0 often unused / NaNs as in ydata
    v = rng.multivariate_normal(mean=np.zeros(p), cov=R, size=N_steps)  # ↳measurement noise
    for k in range(1, N_steps):
        y_meas[k, :] = h_meas(x_true[k, :]) + v[k]

    return x_true, y_meas

def run_LKF_on_truth(
    Q_KF, R_KF,
    delta_x0_mean, P0_plus,
    x_traj_nom, y_traj_nom,
    F_list, H_list,
    x_true, y_meas
):

```

```

"""
Run ONE truth-model test (TMT) + linearized KF pass.

Returns a dict with everything needed for NEES/NIS and plotting:
- x_true, y_meas: nonlinear truth state and noisy measurement trajectories
- x_hat_hist: reconstructed full state estimates
- e_hist: state estimation errors ( $x_{true} - x_{hat}$ )
- P_hist: posterior covariances  $P_k^+$ 
- innov_hist: innovations  $_k$ 
- S_hist: innovation covariances  $S_k$ 
- y_pred_hist: predicted measurements from the filter
"""

N_steps, n = x_traj_nom.shape
p = R_true.shape[0]

# =====
# Linearized KF around nominal trajectory
# =====

# Initial filter state in delta_x coordinates
delta_x_hat_plus = delta_x0_mean.copy() # filter's initial mean
P_plus = P0_plus.copy()

# Storage
delta_x_hat_hist = np.zeros((N_steps, n))
P_hist = np.zeros((N_steps, n, n))
x_hat_hist = np.zeros((N_steps, n))
y_pred_hist = np.zeros((N_steps, p))
innov_hist = np.zeros((N_steps, p))
S_hist = np.zeros((N_steps, p, p))

for k in range(N_steps):
    x_nom_k = x_traj_nom[k, :]
    F_k = F_list[k]
    H_k = H_list[k]

    # Time update (predict)
    delta_x_hat_minus = F_k @ delta_x_hat_plus
    P_minus = F_k @ P_plus @ F_k.T + Q_KF
    S_k = H_k @ P_minus @ H_k.T + R_KF
    K_k = P_minus @ H_k.T @ np.linalg.inv(S_k)

    # Get innovation term
    y_meas_k = y_meas[k, :]
    y_pred_k = y_traj_nom[k] + H_k @ delta_x_hat_minus

```

```

innov_k = np.zeros(p) if k == 0 else y_meas_k - y_pred_k      # t=0 data
↳ is NaNs

# Measurement update (KF)
delta_x_hat_plus = delta_x_hat_minus + K_k @ innov_k
P_plus = (np.eye(n) - K_k @ H_k) @ P_minus

# Reconstruct full state estimate
x_hat_k = x_nom_k + delta_x_hat_plus

# Log everything
delta_x_hat_hist[k, :] = delta_x_hat_plus
P_hist[k, :, :] = P_plus
x_hat_hist[k, :] = x_hat_k
y_pred_hist[k, :] = y_pred_k
innov_hist[k, :] = innov_k
S_hist[k, :, :] = S_k

# Save state estimation errors
e_hist = x_true - x_hat_hist # shape (N_steps, 6)
sigma_hist = np.sqrt(np.stack([np.diag(P_hist[k, :, :]) for k in
↳ range(N_steps)], axis=0)) # (N_steps, 6)

return {
    "x_hat_hist": x_hat_hist,
    "e_hist": e_hist,
    "P_hist": P_hist,
    "sigma_hist": sigma_hist,
    "y_pred_hist": y_pred_hist,
    "innov_hist": innov_hist,
    "S_hist": S_hist,
}

# =====
# 4(a) - Single Run
# =====
# Start with truth, will tune later
Q_KF = Q_true.copy()
R_KF = R_true.copy()

# Get Truth
x_true, y_meas = simulate_truth_trajectory(
    Q=Q_true,
    R=R_true,
    delta_x0_mean=delta_x0_mean,
    P0_plus=P0_plus,
)

```

```

        x_traj_nom=x_traj_nom,
        u_nom=u_nom,
        L_nom=L_nom,
        dt=dt,
        rng=rng
    )

# Run the LKF
result = run_LKF_on_truth(
    Q_KF=Q_KF, R_KF=R_KF,
    delta_x0_mean=delta_x0_mean,
    P0_plus=P0_plus,
    x_traj_nom=x_traj_nom,
    y_traj_nom=y_traj_nom,
    F_list=F_list,
    H_list=H_list,
    x_true=x_true,
    y_meas=y_meas
)

```

x_hat_hist = result["x_hat_hist"]
y_pred_hist = result["y_pred_hist"]
sigma_hist = result["sigma_hist"]
e_hist = result["e_hist"]

```

def plot_single_run_results(
    tvec,
    x_true,
    y_meas,
    x_hat_hist,
    y_pred_hist,
    e_hist,
    sigma_hist,
    state_labels,
    meas_labels,
    filter_name="KF",
    save_prefix=None
):
    """
    Shared plotting for Parts 4(a) and 5(a).

```

Produces three figures:

1. *Truth vs estimate (states)*
2. *Noisy measurement vs predicted measurement*
3. *State estimation errors with +/- 2 sigma bounds*

```

If save_prefix is provided, files are saved as:
    f"{save_prefix}_states.png"
    f"{save_prefix}_measurements.png"
    f"{save_prefix}_errors.png"
"""

n = x_true.shape[1]
p = y_meas.shape[1]

state_labels = [
    r"\xi_g [m]",
    r"\eta_g [m]",
    r"\theta_g [rad]",
    r"\xi_a [m]",
    r"\eta_a [m]",
    r"\theta_a [rad]",
]
# Writing this manually since the format given in .mat file doesn't work
# with python
meas_labels = [
    r"\gamma_{ag} [rad]",
    r"\rho_{ga} [m]",
    r"\gamma_{ga} [rad]",
    r"\xi_a [m]",
    r"\eta_a [m]",
]
# State Trajectories
fig1, axes1 = plt.subplots(3, 2, figsize=(12, 9))
axes1 = axes1.ravel()

for i in range(n):
    ax = axes1[i]
    ax.plot(tvec, x_true[:, i], label="Nonlinear truth", linewidth=1.8)
    ax.plot(tvec, x_hat_hist[:, i], "--", label=f"{filter_name} estimate", linewidth=1.8)
    ax.set_ylabel(state_labels[i])
    ax.set_xlabel("Time [s]")
    ax.grid(True)
    if i == 0:
        ax.legend(loc="best")

fig1.suptitle(f"{filter_name}: Truth vs Estimated States", fontsize=14)
fig1.tight_layout(rect=[0, 0.03, 1, 0.95])

```

```

if save_prefix is not None:
    fig1.savefig(f"{save_prefix}_states.png")

# Measurements
fig2, axes2 = plt.subplots(3, 2, figsize=(12, 9))
axes2 = axes2.ravel()

for i in range(p):
    ax = axes2[i]
    ax.plot(tvec, y_meas[:, i], label="Noisy measurement", linewidth=1.8)
    ax.plot(tvec, y_pred_hist[:, i], "--", label=f"{filter_name} predicted_{meas}", linewidth=1.8)
    ax.set_ylabel(meas_labels[i])
    ax.set_xlabel("Time [s]")
    ax.grid(True)
    if i == 0:
        ax.legend(loc="best")

axes2[-1].axis("off")

fig2.suptitle(f"{filter_name}: Measurements vs Predicted Measurements", fontsize=14)
fig2.tight_layout(rect=[0, 0.03, 1, 0.95])

if save_prefix is not None:
    fig2.savefig(f"{save_prefix}_measurements.png")

# State Estimation Error +2 sigma
fig3, axes3 = plt.subplots(3, 2, figsize=(12, 9))
axes3 = axes3.ravel()

for i in range(n):
    ax = axes3[i]
    ax.plot(tvec, e_hist[:, i], label="Estimation error", linewidth=1.5)
    ax.plot(tvec, 2.0 * sigma_hist[:, i], "k--", label=r"$+2\sigma$",
            linewidth=1.0)
    ax.plot(tvec, -2.0 * sigma_hist[:, i], "k--", label=r"$-2\sigma$",
            linewidth=1.0)
    ax.set_ylabel(state_labels[i])
    ax.set_xlabel("Time [s]")
    ax.grid(True)
    if i == 0:
        ax.legend(loc="best")

fig3.suptitle(rf"{filter_name}: Estimation Errors with $\pm2\sigma$ Bounds (Single Run)", fontsize=14)
fig3.tight_layout(rect=[0, 0.03, 1, 0.95])

```

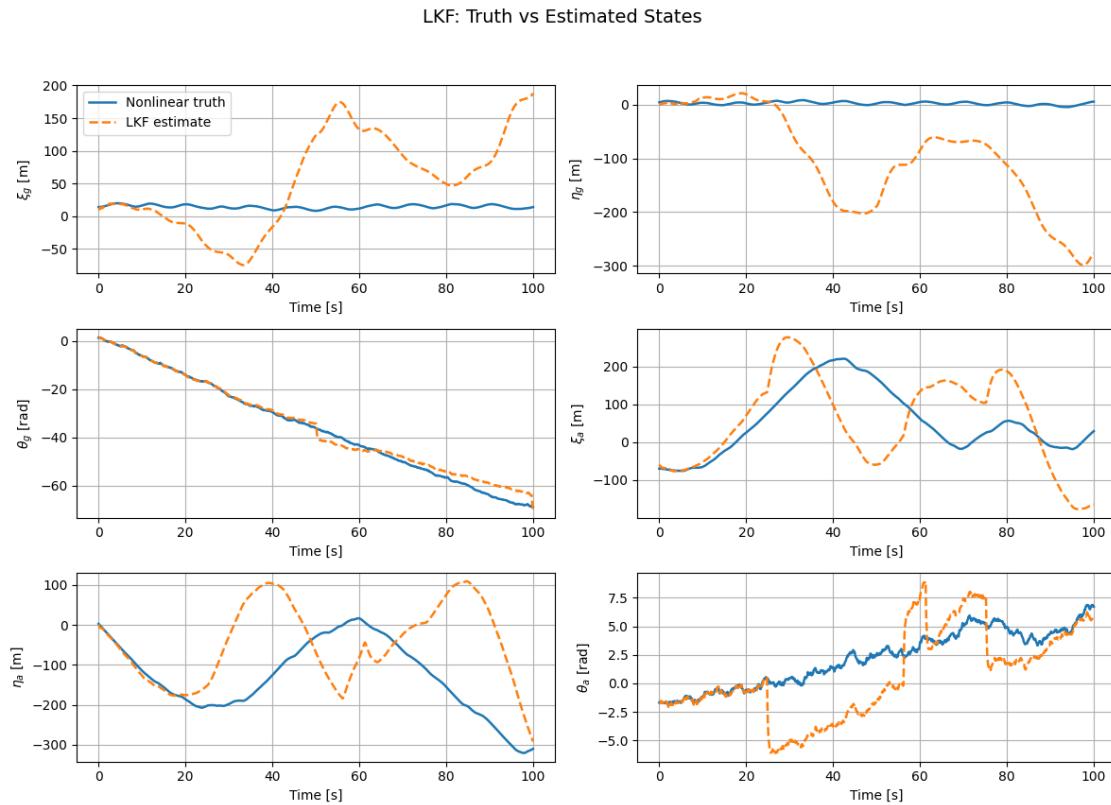
```

if save_prefix is not None:
    fig3.savefig(f"{save_prefix}_errors.png")

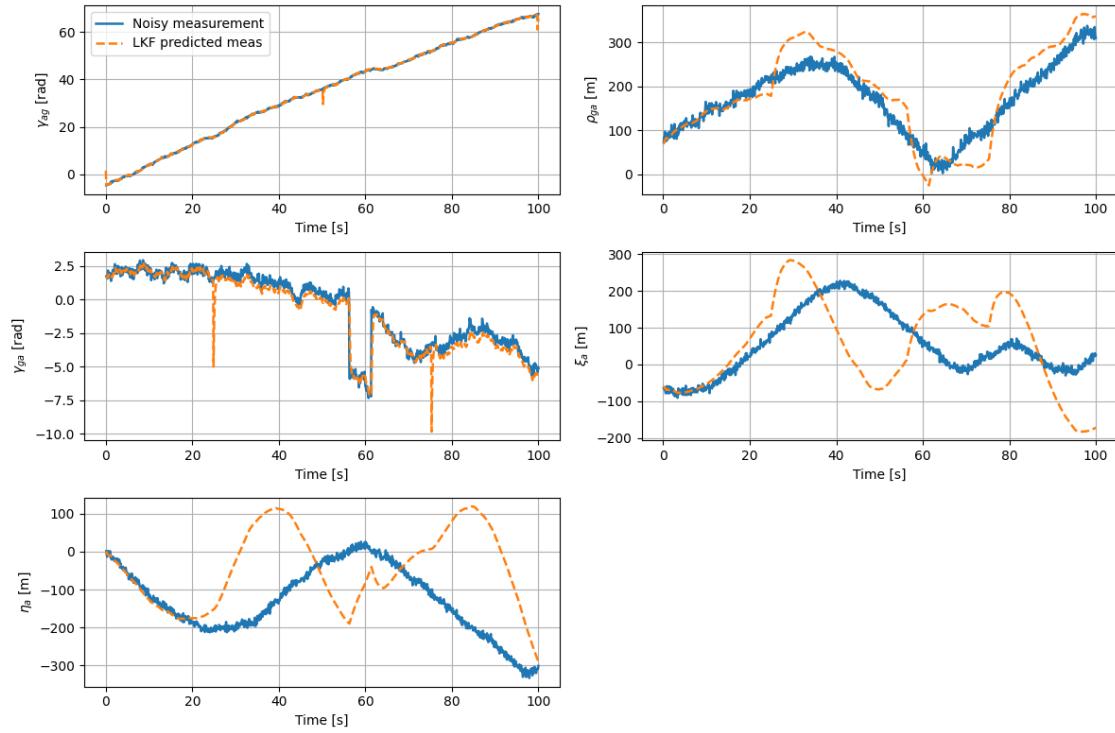
plt.show()

plot_single_run_results(
    tvec=tvec,
    x_true=x_true,
    y_meas=y_meas,
    x_hat_hist=x_hat_hist,
    y_pred_hist=y_pred_hist,
    e_hist=e_hist,
    sigma_hist=sigma_hist,
    state_labels=state_labels,
    meas_labels=meas_labels,
    filter_name="LKF",
    save_prefix="4a_LKF"
)

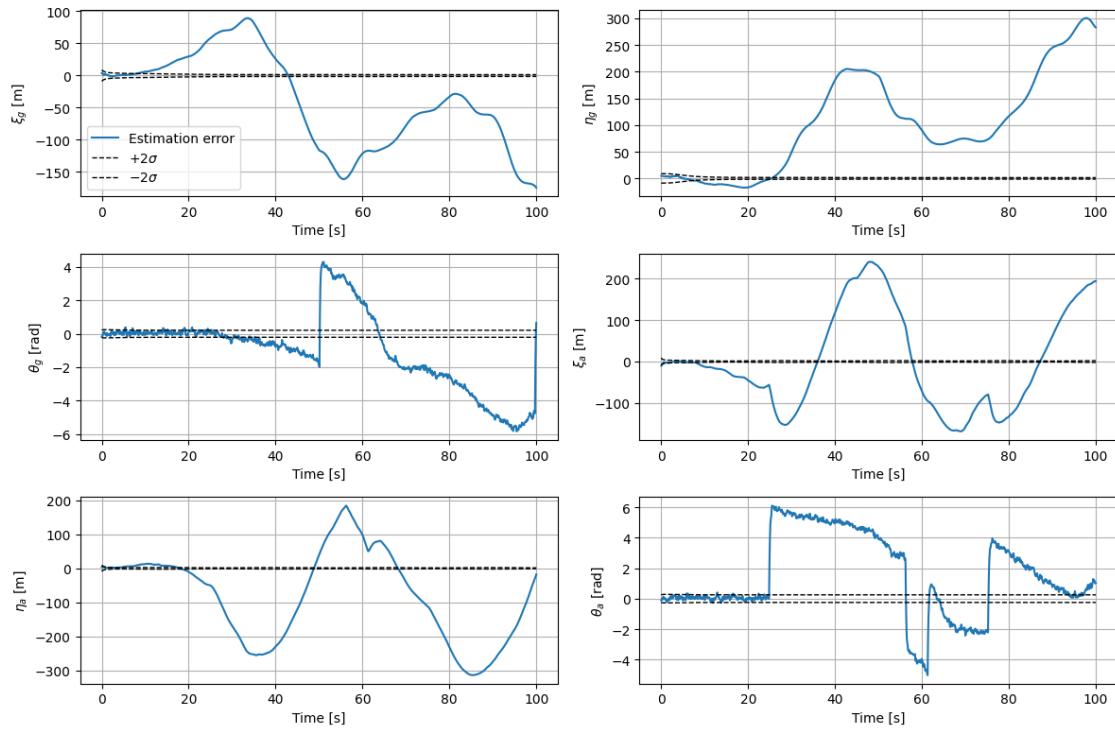
```



LKF: Measurements vs Predicted Measurements



LKF: Estimation Errors with $\pm 2\sigma$ Bounds (Single Run)



1 4 part (b) and (c) - LKF NEES and NIS

```
[5]: from scipy.stats import chi2

def run_mc_TMT(
    filter_runner,      # run_LKF_on_truth or run_EKF_on_truth
    Q_true, R_true,
    Q_KF, R_KF,
    delta_x0_mean, P0_plus,
    x_traj_nom, y_traj_nom,
    F_list, H_list,    # only used by LKF; EKF can ignore via wrapper
    u_nom, L_nom, dt,
    N_mc, alpha, rng
):
    """
    Run Monte Carlo Truth Model Tests
    Returns NEES and NIS data.
    filter_runner can either be the LKF or EKF function
    """

```

```

N_steps, n = x_traj_nom.shape
p = R_true.shape[0]

NEES = np.zeros((N_mc, N_steps))
NIS  = np.zeros((N_mc, N_steps))

for i in range(N_mc):
    # Get truth
    x_true, y_meas = simulate_truth_trajectory(
        Q=Q_true,
        R=R_true,
        delta_x0_mean=delta_x0_mean,
        P0_plus=P0_plus,
        x_traj_nom=x_traj_nom,
        u_nom=u_nom,
        L_nom=L_nom,
        dt=dt,
        rng=rng
    )

    # Run chosen filter on this truth
    if filter_runner is run_LKF_on_truth:
        result = filter_runner(
            Q_KF=Q_KF, R_KF=R_KF,
            delta_x0_mean=delta_x0_mean,
            P0_plus=P0_plus,
            x_traj_nom=x_traj_nom,
            y_traj_nom=y_traj_nom,
            F_list=F_list,
            H_list=H_list,
            x_true=x_true,
            y_meas=y_meas
        )
    else:
        result = filter_runner(
            Q_KF=Q_KF, R_KF=R_KF,
            delta_x0_mean=delta_x0_mean,
            P0_plus=P0_plus,
            x_traj_nom=x_traj_nom,
            u_nom=u_nom,
            L_nom=L_nom,
            x_true=x_true,
            y_meas=y_meas,
            dt=dt
        )

    e_hist      = result["e_hist"]

```

```

P_hist      = result["P_hist"]
innov_hist = result["innov_hist"]
S_hist      = result["S_hist"]

# NEES
for k in range(N_steps):
    e_k = e_hist[k, :]
    P_k = P_hist[k, :, :]
    NEES[i, k] = e_k @ np.linalg.solve(P_k, e_k)

# NIS (skip k=0)
for k in range(1, N_steps):
    innov_k = innov_hist[k, :]
    S_k = S_hist[k, :, :]
    NIS[i, k] = innov_k @ np.linalg.solve(S_k, innov_k)

return NEES, NIS

def plot_consistency_stat(
    tvec_plot,
    samples,           # shape (N_mc, N_steps_plot)
    dof,              # n for NEES, p for NIS
    alpha,
    quantity_name,   # e.g. "NEES (LKF)" or "NIS (EKF)"
    ylabel,           # e.g. "NEES" or "NIS"
    legend_loc="upper left"
):
    """
    Plot consistency statistic (NEES or NIS) vs time with chi-square bounds.

    tvec_plot : (N_steps_plot,) time array
    samples   : (N_mc, N_steps_plot) array of per-run stats
    dof       : degrees of freedom (n for NEES, p for NIS)
    alpha     : significance level (e.g. 0.05 for 95% CI)
    """
    N_mc, N_steps_plot = samples.shape

    # Sample mean over Monte Carlo runs
    stat_mean = np.mean(samples, axis=0)

    # Chi-square bounds for the mean
    df_total = N_mc * dof
    lower_bound = chi2.ppf(alpha / 2.0, df=df_total) / N_mc
    upper_bound = chi2.ppf(1.0 - alpha / 2.0, df=df_total) / N_mc

    print(f"{quantity_name} mean bounds at {100*(1-alpha):.1f}% CI:")

```

```

print(f"  lower: {lower_bound:.3f}, upper: {upper_bound:.3f}")
print(f"  expected mean = dof = {dof}")

# ----- Plot -----
fig, ax = plt.subplots(figsize=(12, 5))

# Plot a few transparent trajectories to show spread
for i in range(min(N_mc, 20)):
    ax.plot(tvec_plot, samples[i, :], color="0.8", linewidth=0.6, alpha=0.6)

# Plot mean across MC runs
ax.plot(tvec_plot, stat_mean, label=f"Mean {quantity_name}", linewidth=2)

# Expected value and bounds
ax.hlines(dof, tvec_plot[0], tvec_plot[-1],
           colors="g", linestyles="--",
           label=f"Expected mean = {dof}")
ax.hlines(lower_bound, tvec_plot[0], tvec_plot[-1],
           colors="r", linestyles="--",
           label=f"Lower bound ({100*(1-alpha):.1f}% CI)")
ax.hlines(upper_bound, tvec_plot[0], tvec_plot[-1],
           colors="r", linestyles="--",
           label=f"Upper bound ({100*(1-alpha):.1f}% CI)")

ax.set_xlabel("Time [s]")
ax.set_ylabel(ylabel)
ax.set_title(f"{quantity_name} Consistency Test")
ax.grid(True)
ax.legend(loc=legend_loc)

plt.tight_layout()
plt.show()

# =====
# Run LKF Simulation
# =====
# Monte Carlo parameters
N_mc = 200 # number of Monte Carlo runs
alpha = 0.05 # significance level for 95% confidence interval

# Choose noise matrices - can be tuned
Q_KF = Q_true.copy()
R_KF = R_true.copy()

# Reproducible results while still having each trial be different
rng = np.random.default_rng(123)

```

```

NEES_LKF, NIS_LKF = run_mc_TMT(
    filter_runner=run_LKF_on_truth,
    Q_true=Q_true, R_true=R_true,
    Q_KF=Q_KF, R_KF=R_KF,
    delta_x0_mean=delta_x0_mean, P0_plus=P0_plus,
    x_traj_nom=x_traj_nom, y_traj_nom=y_traj_nom,
    F_list=F_list, H_list=H_list,
    u_nom=u_nom, L_nom=L_nom, dt=dt,
    N_mc=N_mc, alpha=alpha, rng=rng
)

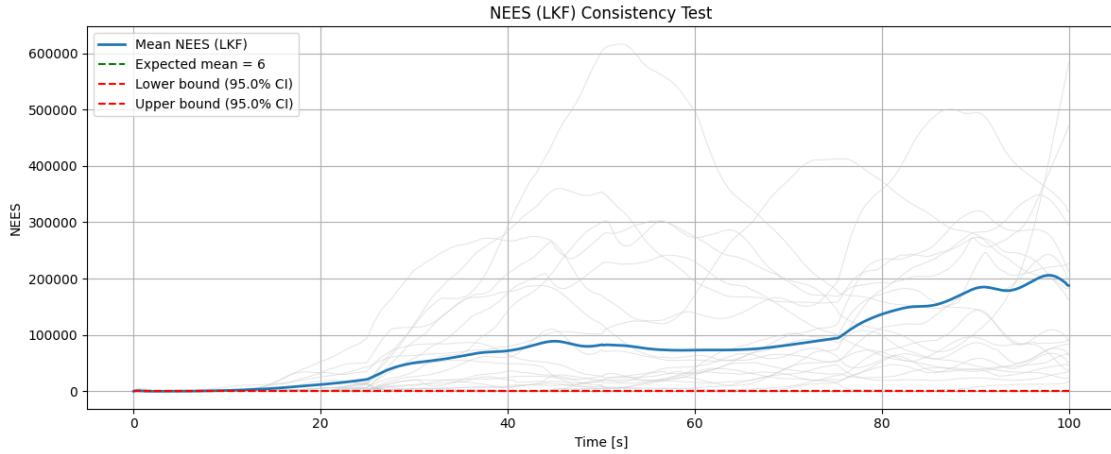
# =====
# Plot LKF Results
# =====

# NEES
plot_consistency_stat(
    tvec_plot=tvec,
    samples=NEES_LKF,
    dof=n,
    alpha=alpha,
    quantity_name="NEES (LKF)",
    ylabel="NEES",
    legend_loc="upper left",
)

# NIS - skip 0
plot_consistency_stat(
    tvec_plot=tvec[1:],
    samples=NIS_LKF[:, 1:],
    dof=p,
    alpha=alpha,
    quantity_name="NIS (LKF)",
    ylabel="NIS",
    legend_loc="upper right",
)

```

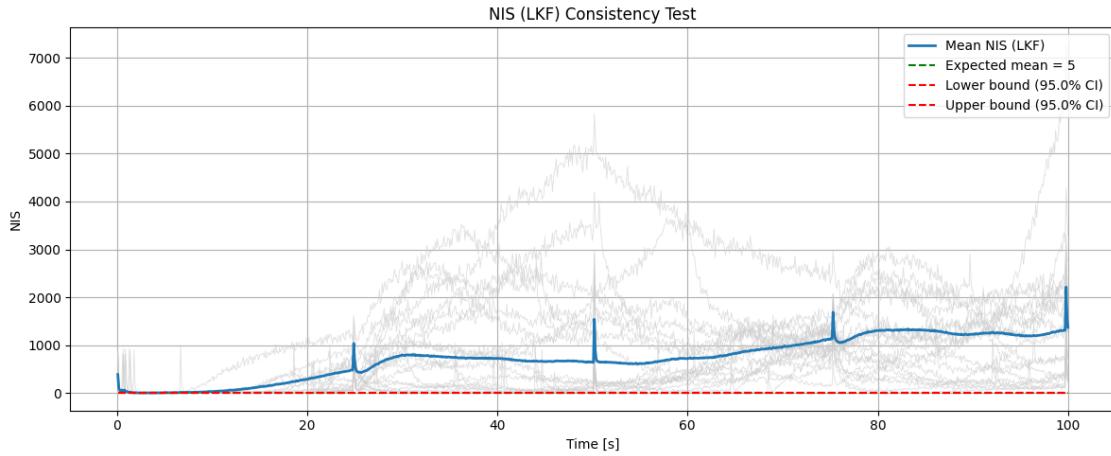
NEES (LKF) mean bounds at 95.0% CI:
lower: 5.529, upper: 6.489
expected mean = dof = 6



NIS (LKF) mean bounds at 95.0% CI:

lower: 4.571, upper: 5.448

expected mean = dof = 5



Issues:

- Nonlinear truth vs linearized filter about a noise-free nominal trajectory
- Process noise strong enough that x_{true} drifts far from $x_{\text{traj_nom}}$

2 Question 5 - EKF

```
[6]: def run_EKF_on_truth(
    Q_KF, R_KF,
    delta_x0_mean, P0_plus,
    x_traj_nom,
    u_nom, L_nom,
    x_true, y_meas,
```

```

dt
):
"""
Run ONE EKF pass on a given truth trajectory and measurements.
Same return structure as run_LKF_on_truth.
"""

N_steps, n = x_traj_nom.shape
p = R_KF.shape[0]

# Initial filter state: nominal + mean perturbation
x_hat_plus = x_traj_nom[0, :] + delta_x0_mean
P_plus = P0_plus.copy()

# Storage
x_hat_hist = np.zeros((N_steps, n))
P_hist = np.zeros((N_steps, n, n))
y_pred_hist = np.zeros((N_steps, p))
innov_hist = np.zeros((N_steps, p))
S_hist = np.zeros((N_steps, p, p))

# Initial conditions
x_hat_hist[0, :] = x_hat_plus
P_hist[0, :, :] = P_plus
y_pred_hist[0, :] = np.nan
innov_hist[0, :] = np.zeros(p)
S_hist[0, :, :] = np.eye(p)

for k in range(N_steps):
    # Time update (predict)
    x_hat_minus = x_hat_plus + dt * f_ct(x_hat_plus, u_nom, L_nom)
    # Get matrices for current iteration
    F_k = np.eye(n) + dt * np.array(A_eval(x_hat_plus, u_nom, L_nom), ↴
                                     dtype=float)
    H_k = C_k = np.array(C_eval(x_hat_minus), dtype=float)
    # Continue prediction step
    P_minus = F_k @ P_plus @ F_k.T + Q_KF
    S_k = H_k @ P_minus @ H_k.T + R_KF
    K_k = P_minus @ H_k.T @ np.linalg.inv(S_k)

    # Get innovation term
    y_meas_k = y_meas[k, :]
    y_pred_k = h_meas(x_hat_minus)      # EKF uses estimate instead of nominal
    innov_k = np.zeros(p) if k == 0 else y_meas_k - y_pred_k      # t=0 data ↴
    ↴is NaNs

    # Measurement update
    x_hat_plus = x_hat_minus + K_k @ innov_k

```

```

P_plus = (np.eye(n) - K_k @ H_k) @ P_minus

# Log everything
x_hat_hist[k, :] = x_hat_plus
P_hist[k, :, :] = P_plus
y_pred_hist[k, :] = y_pred_k
innov_hist[k, :] = innov_k
S_hist[k, :, :] = S_k

# Errors and sigmas
e_hist = x_true - x_hat_hist
sigma_hist = np.sqrt(np.stack([np.diag(P_hist[k, :, :]) for k in
                                range(N_steps)]), axis=0))

return {
    "x_hat_hist": x_hat_hist,
    "e_hist": e_hist,
    "P_hist": P_hist,
    "sigma_hist": sigma_hist,
    "y_pred_hist": y_pred_hist,
    "innov_hist": innov_hist,
    "S_hist": S_hist,
}
}

# =====
# Part 5 (a) - Single-run TMT
# =====

# You can reuse the same Q_KF, R_KF, delta_x0_mean, P0_plus as LKF
Q_KF_EKF = Q_true.copy()
R_KF_EKF = R_true.copy()

# New RNG for this section (optional, just to decouple from LKF run)
rng_ekf = np.random.default_rng(321)

# Generate one nonlinear truth trajectory + noisy measurements
x_true_ekf, y_meas_ekf = simulate_truth_trajectory(
    Q=Q_true,
    R=R_true,
    delta_x0_mean=delta_x0_mean,
    P0_plus=P0_plus,
    x_traj_nom=x_traj_nom,
    u_nom=u_nom,
    L_nom=L_nom,
    dt=dt,
    rng=rng_ekf
)

```

```

)
# Run EKF on this truth
result_EKF = run_EKF_on_truth(
    Q_KF=Q_KF_EKF, R_KF=R_KF_EKF,
    delta_x0_mean=delta_x0_mean,
    P0_plus=P0_plus,
    x_traj_nom=x_traj_nom,
    u_nom=u_nom,
    L_nom=L_nom,
    x_true=x_true_ekf,
    y_meas=y_meas_ekf,
    dt=dt
)

x_hat_hist_ekf  = result_EKF["x_hat_hist"]
y_pred_hist_ekf = result_EKF["y_pred_hist"]
sigma_hist_ekf  = result_EKF["sigma_hist"]
e_hist_ekf       = result_EKF["e_hist"]      # (N_steps, n)

plot_single_run_results(
    tvec=tvec,
    x_true=x_true_ekf,
    y_meas=y_meas_ekf,
    x_hat_hist=result_EKF["x_hat_hist"],
    y_pred_hist=result_EKF["y_pred_hist"],
    e_hist=result_EKF["e_hist"],
    sigma_hist=result_EKF["sigma_hist"],
    state_labels=state_labels,
    meas_labels=meas_labels,
    filter_name="EKF",
    save_prefix="5a_EKF"
)

# =====
# 5 (b) and (c) - NEES and NIS
# =====
NEES_EKF, NIS_EKF = run_mc_TMT(
    filter_runner=run_EKF_on_truth,
    Q_true=Q_true, R_true=R_true,
    Q_KF=Q_KF, R_KF=R_KF,
    delta_x0_mean=delta_x0_mean, P0_plus=P0_plus,
    x_traj_nom=x_traj_nom, y_traj_nom=y_traj_nom,
    F_list=F_list, H_list=H_list,
    u_nom=u_nom, L_nom=L_nom, dt=dt,
)

```

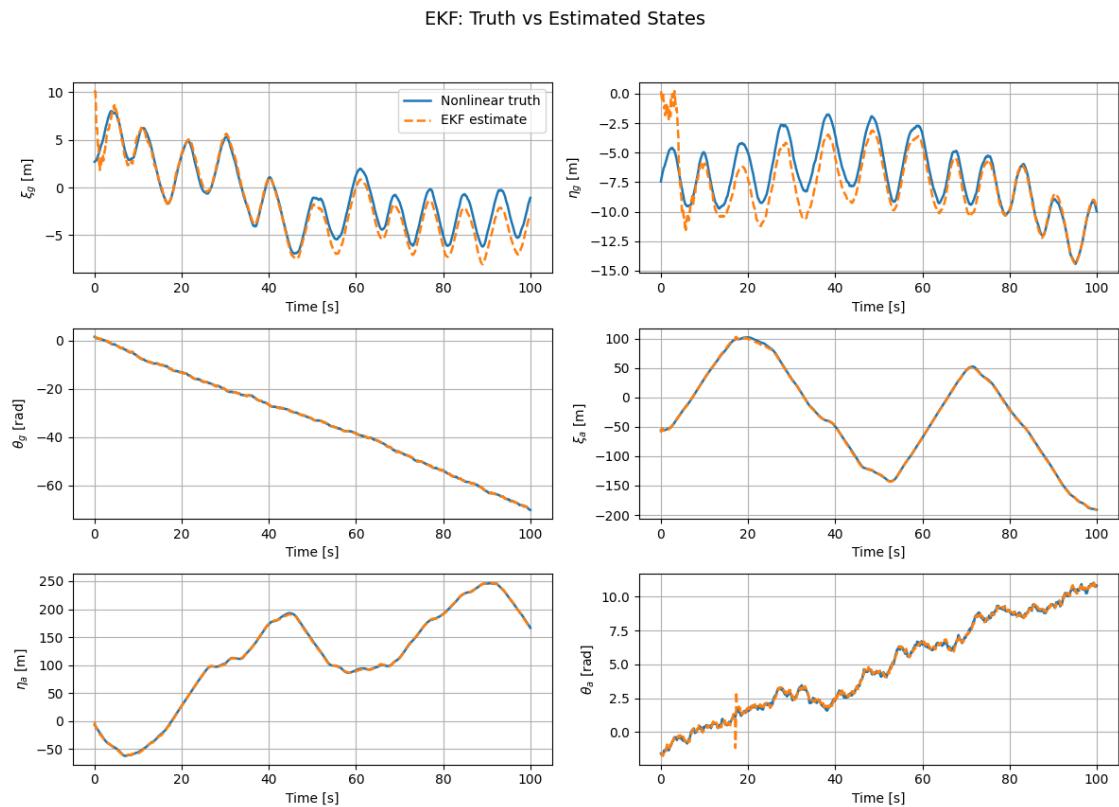
```

N_mc=N_mc, alpha=alpha, rng=rng
)

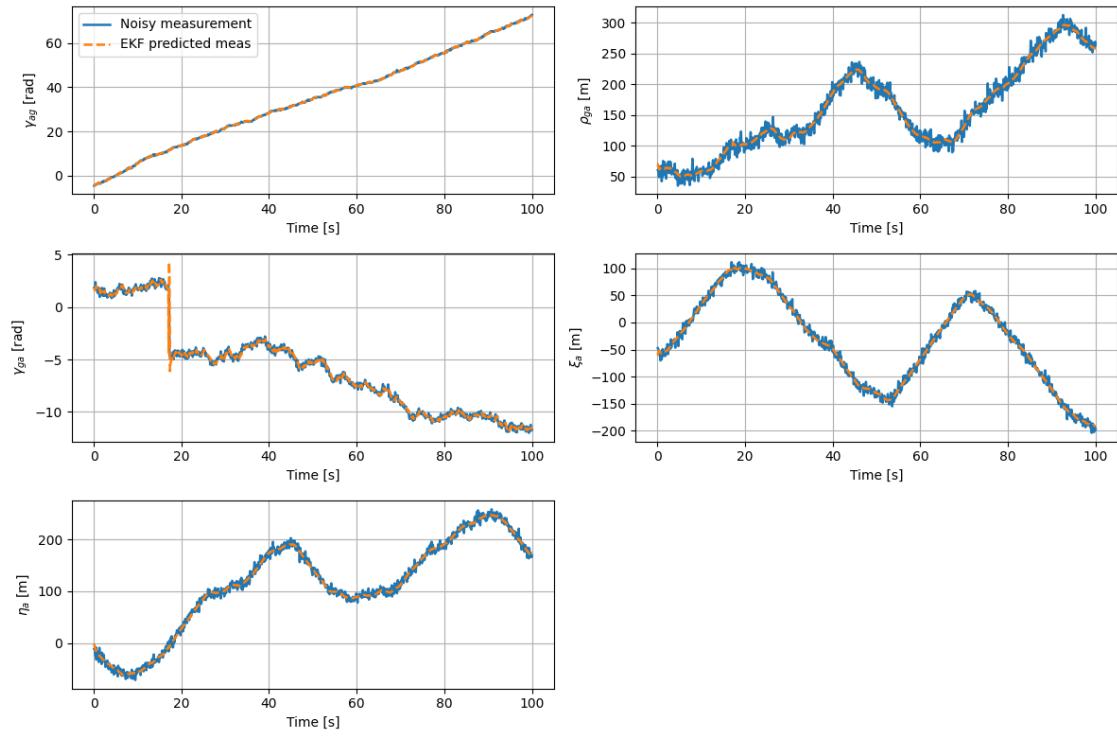
plot_consistency_stat(
    tvec_plot=tvec,
    samples=NEES_EKF,
    dof=n,
    alpha=alpha,
    quantity_name="NEES (EKF)",
    ylabel="NEES",
    legend_loc="upper left",
)

plot_consistency_stat(
    tvec_plot=tvec[1:],
    samples=NIS_EKF[:, 1:],
    dof=p,
    alpha=alpha,
    quantity_name="NIS (EKF)",
    ylabel="NIS",
    legend_loc="upper right",
)

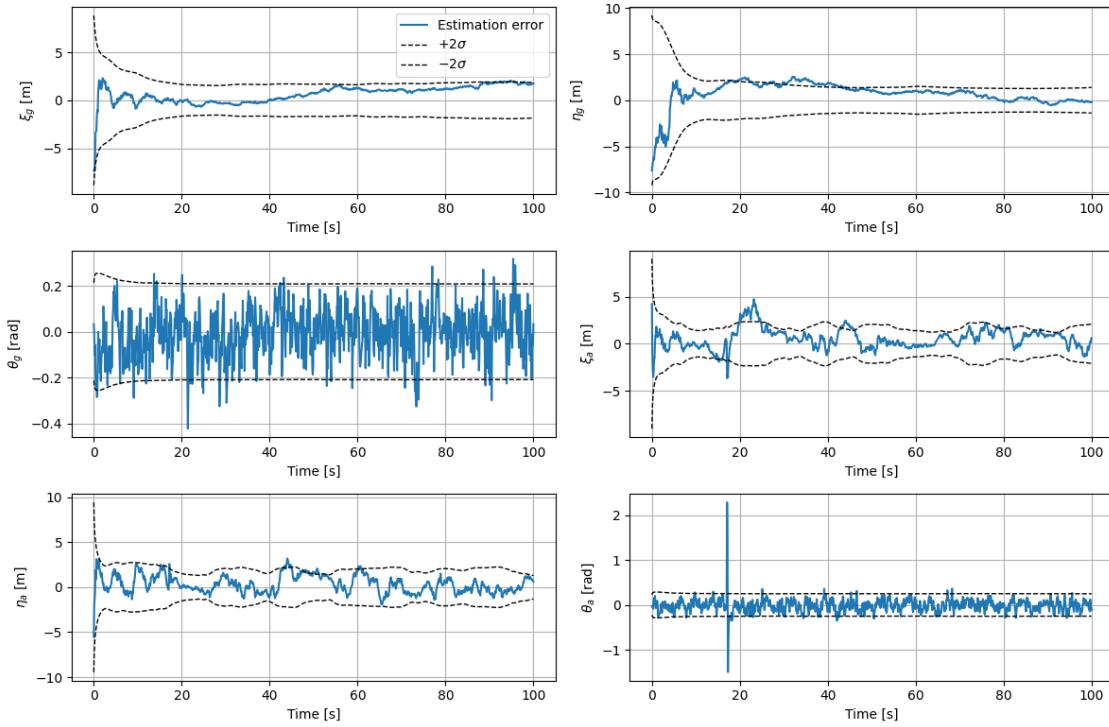
```



EKF: Measurements vs Predicted Measurements



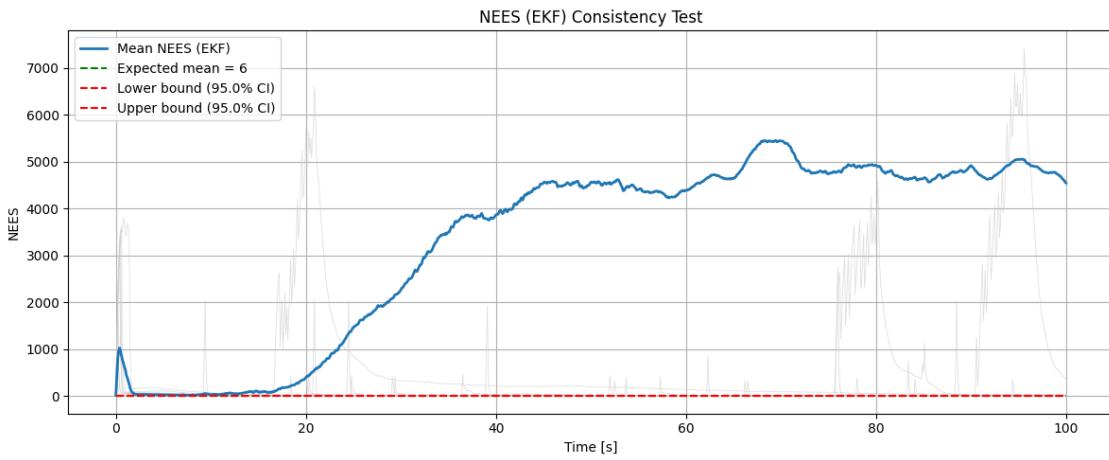
EKF: Estimation Errors with $\pm 2\sigma$ Bounds (Single Run)



NEES (EKF) mean bounds at 95.0% CI:

lower: 5.529, upper: 6.489

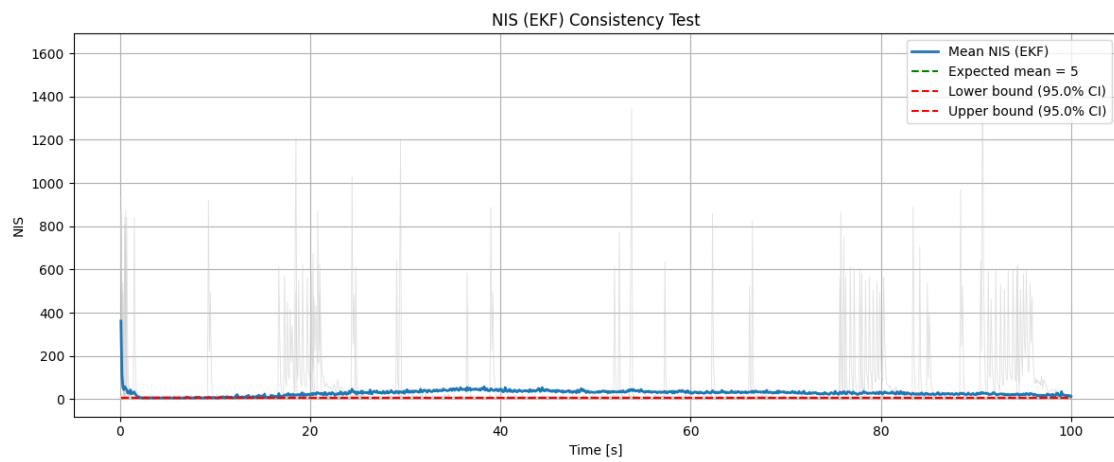
expected mean = dof = 6



NIS (EKF) mean bounds at 95.0% CI:

lower: 4.571, upper: 5.448

expected mean = dof = 5



[]: