# Exploitation of Rotational Symmetries to Solve the Swarm Initialization Problem

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- We have developed a sound, mathematical basis for solving the swarm initialization problem in a special case.
  - We will assume a circular swarm trajectory (eccentricity = 0, exact)
  - We define two operations which preserve swarm optimality:
    - Rotation of the swarm within some known space of valid rotations;
    - Transposition of any two satellites at any point in the orbit.
- **Quantization of the solution space** has enabled the use of discretized optimization methods chiefly, the *Munkres Algorithm*.
- A staged optimization approach solves the problem in successively more detailed passes.
- **Computational slowness** is the current major obstacle to implementing this method.
  - Careful algorithm selection;
  - Parallelization of the process;
  - Dividing stages by regularity of use;













In prior discussions, a swarm was defined to be a close-flying formation of satellites for which the following quantities could be defined:

## • Swarm envelope

- A **closed**, **convex set of points with no holes**, **gaps**, **or voids**, which contains all satellites in the swarm.
  - Mathematically speaking, the boundary of the envelope must be simply connected.
- The swarm envelope **may not contain points inaccessible to the satellites** (e.g., points at or below Earth's surface).
- For this discussion, we will use a **spherical envelope**.
- Must include a **reference point** that marks the envelope center.
  - This point **need not to be an element** of the swarm envelope.
  - For this discussion, we will use the center of the sphere.























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- Swarm trajectory
  - A **user-defined function of time** which specifies the position of the swarm envelope relative to any point with known coordinates.
    - This latter point can be taken to be the **center of the Earth**.
    - Position measured to the **envelope-relative reference point**.
  - If necessary, this definition may include a function to specify **the orientation of the swarm envelope**.
    - Unnecessary in the case of a spherical envelope.
  - While it is not a strict requirement in the general case, **trajectories should follow spacetime geodesics**.
  - For this discussion, we will assume that the swarm trajectory **can be approximated by a Keplerian orbit of zero eccentricity**.





















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- Swarm distribution
  - The **arrangement of satellites** within the sphere envelope as defined relative to the envelope-relative reference point.
  - The most abstract component of the satellite swarm, incorporating mission-specific parameters, an unknown number of degrees of freedom, and the swarm cost functional.
  - For this discussion, we will assume that the swarm distribution can be rotated about the envelope relative reference point.
- Swarm centroid
  - A function of the satellite positions which **determines the true center of the swarm**.
  - Distinct from the envelope-relative reference point, as the centroid is **determined by satellite positions, not envelope geometry**.
  - For this discussion, we will use the **arithmetic mean** of all satellite positions to define the swarm centroid.















For this discussion, we will default to our test case of **twelve satellites**:

- The swarm envelope is a **sphere of radius** *ρ* with reference at its center.
- The swarm trajectory is a **circular orbit of radius** *a*.
- The satellites in the swarm should be distributed at the **vertices of a regular icosahedron**.
  - The swarm may be **transformed by any rotation** about its center.
- The swarm centroid is defined using the arithmetic mean:

$$\vec{r}_0(t) = \frac{1}{12} \sum_{i=1}^{12} \vec{r}_i(t)$$









### **Test Swarm Geometry**





#### A rotating icosahedron is still an icosahedron.

Animation from Wikipedia.













## **Test Swarm Configuration**







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## **Formation Chains**











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## **Formation Chains**













- Note that *n* successive rotations by  $2\pi/n$  radians about a given axis will result in a net angle of  $2\pi$  radians.
- We may, however, rotate by *multiples* of  $2\pi/n$  radians.
  - Recognizing that a rotation by  $2\pi$  radians is equivalent to no rotation, it follows that rotations by multiples of n + 1 to 2n 1 are equivalent to rotations by **multiples of 1 to n 1.**
  - We may also consider the negative multiples (i.e., -(n 1) to -1).
  - Rotations by a multiple of 0 need only be considered about a single axis, since this corresponds to zero rotation.
- With two distinct rotations to check (the starting attitude and the change in attitude), we must check  $4m^2n(n-1) + 1$  formation chains.
- For 46 axes and 12 satellites, this corresponds to 1,117,249 chains.
- Some of these chains may be sufficiently close to others to qualify as "duplicate." **We may eliminate these duplicate chains**.
  - This process takes several days to a week or more on a single CPU, but only needs to be done once (the results may be saved and re-used).















- For 46 axes and 12 satellites, **with duplicate elimination**, we obtain **between 5,000 and 10,000** formation chains depending on tolerance.
- These formation chains are saved and used in the next step.
  - Note that, because these chains depend only on the selected value of the rotation, this process is **invariant with respect to the radii of the formation, or its orbit.**















- Once a group of formation chains are selected, we must then pull orbits from them by **selecting the first two formations out of the chain**.
- For this pair, we may connect **each satellite from the first formation to any position from the second.** 
  - Connecting the *i*th satellite to the *j*th position yields an orbit. Comparing this orbit to the formation chain yields a contribution to a cost function,  $J_{ij}$ .
  - For each *i* and *j*, we select the two formations with minimal  $J_{ij}$ .
- Note that, for each pair, we may propagate the assignment to another formation, so we must compare the values of  $J_{ij}$  for each formation in the chain.
- The optimal assignment may be determined using the Munkres Algorithm













## **Selecting Orbits**







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## **Selecting Orbits**



- The Munkres algorithm runs in  $O(n^4)$  time, so we are currently investigating ways to avoid running the Munkres algorithm in cases where it would clearly produce a suboptimal result.
  - For example, if the sum of the *n* smallest elements in the cost matrix, J is greater than the current best result, we may skip the algorithm.
- Once the best result has been found, that formation chain is selected, along with the assignment function determined by Munkres and the orbits it generates.
- Those orbits are passed on to the next stage of optimization, which uses continuous optimization to further refine the initial velocities and minimize the swarm cost functional.

















- Implement code to run the orbit selection protocol.
- Implement code to refining stage.
- Run this with known test cases, such as the LISA configuration.
- Compare performance for low-n vs. high-n cases.











