## 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;

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## Principal Definitions

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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## Principal Definitions



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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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## Principal Definitions



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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.

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

- The swarm envelope is a sphere of radius $\boldsymbol{\rho}$ with reference at its center.
- The swarm trajectory is a circular orbit of radius $\boldsymbol{a}$.
- 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:

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

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A rotating icosahedron is still an icosahedron. Animation from Wikipedia.



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- 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 $2 n-1$ are equivalent to rotations by multiples of 1 to $\boldsymbol{n - 1}$.
- 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 $\mathbf{4 m ^ { 2 }} \boldsymbol{n}(\boldsymbol{n}-\mathbf{1})+\mathbf{1}$ 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.

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## Selecting Orbits

- 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, $J_{i j}$.
- For each $i$ and $j$, we select the two formations with minimal $J_{i j}$.
- Note that, for each pair, we may propagate the assignment to another formation, so we must compare the values of $J_{i j}$ for each formation in the chain.
- The optimal assignment may be determined using the Munkres Algorithm

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## Selecting Orbits

- The Munkres algorithm runs in $O\left(n^{4}\right)$ 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.

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- 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.

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