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FUEL-OPTIMAL IMPULSIVE FIXED-TIME TRAJECTORIES IN THE LINEARIZED CIRCULAR RESTRICTED 3-BODY PROBLEM

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The problem of fixed-time fuel-optimal trajectories with high-thrust propulsion in the vicinity of a Lagrange point is tackled via the linear version of the primer vector theory. More precisely, the proximity to a Lagrange point i.e. any equilibrium point - stable or not - in the circular restricted three-body problem allows for a linearization of the dynamics. Furthermore, it is assumed that the spacecraft has ungimbaled thrusters, leading to a formulation of the cost function with the 1-norm for space coordinates, even though a generalization exists for steerable thrust and the 2-norm. In this context, the primer vector theory gives necessary and sufficient optimality conditions for admissible solutions to two-value boundary problems. Similarly to the case of rendezvous in the restricted two-body problem, the in-plane and out-of-plane trajectories being uncoupled, they can be treated independently. As a matter of fact, the out-of-plane dynamics is simple enough for the optimal control problem to be solved analytically via this indirect approach. As for the in-plane dynamics, the primer vector solution of the so-called primal problem is derived by solving a hierarchy of linear programs, as proposed recently for the aforementioned rendezvous. The optimal thrusting strategy is then numerically obtained from the necessary and sufficient conditions. Finally, in-plane and out-of-plane control laws are combined to form the complete 3-D fuel-optimal solution. Results are compared to the direct approach that consists in working on a discrete set of times in order to perform optimization in finite dimension. Examples are provided near various Lagrange points in the Sun-Earth and Earth-Moon systems, hinting at the extensive span of possible applications of this technique in station-keeping as well as mission analysis, for instance when connecting manifolds to achieve escape or capture.

I. INTRODUCTION

Lagrange points, a.k.a. libration points, are interesting in binary systems e.g. Earth-Moon, since they are motionless in the corotating frame, and thus offer continuous visibility over one or two of the main bodies. A favoured type of orbit around Lagrange points is the so-called halo one, due to its periodicity. Depending on the binary system, such orbits can be dynamically unstable, creating a need for station-keeping. Even in a stable configuration, orbital perturbations such as radiation pressure can be enough to require trajectory corrections. Hence the minimization of fuel consumption for trajectories in the vicinity of libration points is a topic of high interest. This problem can be formulated as an optimal control problem and tack-
led by a direct or indirect approach. Generally speaking, the direct approach consists in discretizing the continuous control variables into a finite number of parameters that can be fed to a nonlinear programming solver\(^9\). On the other hand, the indirect approach builds on the nature of the problem and aims at guaranteeing the optimality of the solution by looking at the so-called co-state variables, a somewhat generalisation of the Lagrange multipliers\(^4\). For space trajectories with high-thrust propulsion and parametrized with Cartesian coordinates, the indirect approach is best known as the primer vector theory, due to the work of Lawden in the early 1960’s\(^{16}\). In short, it states optimality conditions for a trajectory in terms of a vector (primer vector) whose magnitude has to be always smaller or equal to one for optimality certification. The dates or locations where it reaches unit norm are the ones where the optimal maneuvers should occur. While these optimality conditions are only necessary with nonlinear dynamics, they become sufficient in a linear setting, as proven rigorously by Neustadt\(^{17}\). The general, non-linear theory has already been applied to high-thrust trajectories in the restricted 3-body problem, see for instance the work of Hiday-Johnston \& Howell\(^8\) or Davis et al\(^9\). Despite their more general framework, the optimality conditions in this formulation lack the sufficiency of the linearized dynamics that are yet available in the vicinity of a Lagrange point. Furthermore, the optimal solution to this control problem under the linear approximation, along with its co-state, could be a good initial guess for the nonlinear case to be tackled indirectly. In this work, a modern numerical approach, based on convergent discretization methods for the solution of semi-infinite convex problem (SICP)\(^{20}\) and heuristically improved from the one used for the restricted 2-body problem in\(^1\), is introduced for trajectories near any of the libration points in the circular restricted 3-body problem (CR3BP). In order to speed up the procedure, it is coupled with recent analytical results in the linear theory of the primer vector\(^2\).

Notations: The velocity increment at \(\nu_i\) will be denoted by \(\Delta V(\nu_i)\). \(\{b_i\}_{i=1}^{N}\) is a sequence of variables \(b_i\), \(i = 1, \cdots, N\), and \(\text{sgn}(z)\) is the sign function of the variable \(z\). The prime denotes differentiation with respect to the true anomaly \(\nu\). \(O_{p \times m}\) and \(I_m\) denote respectively the null matrix of dimensions \(p \times m\) and the identity matrix of dimension \(m\). Let \(r \in \mathbb{N}_+\) and \((p, q) \in \mathbb{N} \times \mathbb{N}\) such that: \(1 \leq p \leq \infty\) and \(\frac{1}{q} + \frac{1}{q'} = 1\). Classically, \(C([\nu_0, \nu_f], \mathbb{R}^r)\) is the Banach space of continuous functions \(f : [\nu_0, \nu_f] \to \mathbb{R}^r\) equipped with the norm \(\|f\|_q = \sup_{\nu_0 \leq \nu \leq \nu_f} \|f(\nu)\|_q\). Denote by \(L_{1,p}([\nu_0, \nu_f], \mathbb{R}^r)\) the normed linear space of Lebesgue integrable functions from \([\nu_0, \nu_f]\) to \(\mathbb{R}^r\) with the norm \(\|u\|_{1,p} = \int_{\nu_0}^{\nu_f} \|u(\nu)\|_p d\nu\). Let \(\text{BV}([\nu_0, \nu_f], \mathbb{R}^r)\) be the space of functions of bounded variation over the interval \([\nu_0, \nu_f]\) with: \(\|g\|_{1,p} = \sup_{P_k} \sum_{i=1}^{\kappa} \|g(\nu_i) - g(\nu_{i-1})\|_p\), where the supremum is taken over all finite partitions \(P_k = (\nu_i)_{i=1}^\kappa\) of \([\nu_0, \nu_f]\). For a symmetric real matrix \(S \in \mathbb{R}^{n \times n}\), the notation \(S \preceq 0\) \((S \succeq 0\) stands for the negative (positive) semi-definiteness of \(S\). Finally, \(\chi_A\) is the indicator function of the set \(A\).

II. LINEARIZED DYNAMICS OF THE CR3BP

The circular restricted 3-body problem deals with the motion of a negligible mass \(m\) (a spacecraft for instance) affected by Keplerian gravity field of two primary massive bodies \(m_2 < m_1\) orbiting in a circular orbit around their barycentre \(B\). This relative motion is classically studied in a non-inertial synodic frame \((\vec{i}, \vec{j}, \vec{k})\) rotating with the two major point-masses \(m_1\) and \(m_2\) which origin is located at \(B\). The \(x\)-axis is directed from the primary \(m_1\) to the primary \(m_2\), the \(z\)-axis is parallel to the direction of the orbital angular velocity vector of the primary system with respect to an inertial frame and the \(y\)-axis completes the right-handed system as depicted on Figure 1. The mean motion \(\omega\) of the primary system is constant and given by:

\[
\omega^2 R^3 = G(m_1 + m_2),
\]

where \(G\) is the gravitational constant and \(R\) the distance between \(m_1\) and \(m_2\).

---

\(\Delta V(\nu_i)\) \(O_{p \times m}\) \(I_m\) \(\text{sgn}(z)\) \(C([\nu_0, \nu_f], \mathbb{R}^r)\) \(L_{1,p}([\nu_0, \nu_f], \mathbb{R}^r)\) \(\text{BV}([\nu_0, \nu_f], \mathbb{R}^r)\) \(\|g\|_{1,p}\) \(\chi_A\) \(\omega\) \(G\) \(R\) \(\|u\|_{1,p}\) \(\|g\|_{1,p}\) \(\sup\) \(\sum\) \(\kappa\) \(\nu_0\) \(\nu_f\) \(\mathbb{R}^r\) \(\mathbb{R}^{n \times n}\) \(S \preceq 0\) \(S \succeq 0\) \(\chi_A\) \(\omega\) \(G\) \(R\) \(\|u\|_{1,p}\) \(\|g\|_{1,p}\) \(\sup\) \(\sum\) \(\kappa\) \(\nu_0\) \(\nu_f\) \(\mathbb{R}^r\) \(\mathbb{R}^{n \times n}\) \(S \preceq 0\) \(S \succeq 0\) \(\chi_A\) \(\omega\) \(G\) \(R\) \(\|u\|_{1,p}\) \(\|g\|_{1,p}\) \(\sup\) \(\sum\) \(\kappa\) \(\nu_0\) \(\nu_f\) \(\mathbb{R}^r\) \(\mathbb{R}^{n \times n}\) \(S \preceq 0\) \(S \succeq 0\) \(\chi_A\) \(\omega\) \(G\) \(R\)
as the mass ratio of the primary system (also defined as the non-dimensional distance of \( m_2 \) to \( B \)), it is readily obtained that:

\[
\begin{align*}
    x_1 &= -\frac{Rm_2}{m_1 + m_2} = -R\mu, \\
    x_2 &= \frac{Rm_1}{m_1 + m_2} = R(1 - \mu).
\end{align*}
\]

Based on a Lagrangian formulation of the CR3TBP and changing the independent variable from the time \( t \) to the true anomaly \( \nu \) of the primary system, the non-linear equations of motion of the third body \( m \) expressed in cartesian coordinates \((x, y, z)\) in the rotating frame are given by\(^7\) by:

\[
\begin{align*}
    x'' &= \frac{R^3(1 - \mu)(x + R\mu)}{\|\vec{p}_1\|^3} - \frac{R^3\mu(x - R(1 - \mu))}{\|\vec{p}_2\|^3} + 2y' + x, \\
    y'' &= \frac{R^3y}{\|\vec{p}_1\|^3} - \frac{\mu R^3y}{\|\vec{p}_2\|^3} - 2x' + y, \\
    z'' &= \frac{R^3z}{\|\vec{p}_1\|^3} - \frac{\mu R^3z}{\|\vec{p}_2\|^3},
\end{align*}
\]

where \( \|\vec{p}_1\| \) and \( \|\vec{p}_2\| \) are the distances from the third body to respectively \( m_1 \) and \( m_2 \), \((\cdot)'\) is the derivative with respect to the true anomaly \( \nu \) and \( \dot{\nu} = \omega \). A normalized form \((R = 1 \text{ and } \omega = 1)\) of these equations may be found in a more recent reference\(^8\).

### II.ii Linearized equations around the libration points

It is well known that the dynamical system represented by Equations [5]-[7] exhibits five equilibrium points, the so-called libration points, which are stationary points with respect to the rotating system. The negligible mass \( m \) placed at these points with zero initial velocity will remain there if there is no external perturbation. All five equilibrium points lie in the \( x - y \) plane and two families of libration points are distinguished. The **collinear points** \( L_1-L_3 \) for which \( y = 0 \) and \( z = 0 \), satisfy the following equation:

\[
x = \frac{(1 - \mu)R^3}{(x + R\mu)^2} + \frac{\mu R^3}{(x - R(1 - \mu))^2}.
\]

The **triangular** or **equilateral points** \( L_4, \ L_5 \) are defined by \( z = 0 \) and:

\[
\begin{align*}
    x &= \frac{R}{2} - \mu R, \\
    y &= \pm \frac{\sqrt{3}}{2} R.
\end{align*}
\]

A typical geometrical configuration of all libration points is given in Figure 1.

Defining the small perturbation of a nearby trajectory around an equilibrium point \( L_i \), where \( i \in \{1, \ldots, 5\} \), the linearized autonomous equations of relative motion of \( m \) around the Libration point \( L_i \) are then given in state-space form as:

\[
X'(\nu') = AX(\nu),
\]

where \( X^T = [\delta x \ \delta y \ \delta z \ \delta x' \ \delta y' \ \delta z'] \) and

\[
A = \begin{bmatrix}
    0 & 0 & 1 & 0 & 0 & 0 \\
    0 & 0 & 0 & 0 & 0 & 1 \\
    0 & 0 & 0 & 0 & 0 & 0 \\
    1 - U_{xx} & -U_{xy} & 0 & 0 & 2 & 0 \\
    -U_{xy} & 1 - U_{yy} & 0 & -2 & 0 & 0 \\
    0 & 0 & -U_{zz} & 0 & 0 & 0
\end{bmatrix},
\]

\[
U_{xx} = \frac{R^3(1 - \mu)}{r_{1L_i}^3} + \frac{R^3\mu}{r_{2L_i}^3},
\]

\[
U_{xy} = \frac{3(1 - \mu)R^3(x_{L_i} - R(1 - \mu))}{r_{1L_i}^5} + \frac{3\mu R^3(x_{L_i} + R\mu)}{r_{2L_i}^5},
\]

\[
U_{yy} = \frac{3(1 - \mu)R^3y_{L_i}^2}{r_{1L_i}^5} + \frac{3\mu R^3y_{L_i}^2}{r_{2L_i}^5},
\]

and

\[
r_{1L_i} = \sqrt{(x_{L_i} + R\mu)^2 + y_{L_i}^2},
\]

\[
r_{2L_i} = \sqrt{(x_{L_i} - R(1 - \mu))^2 + y_{L_i}^2}.
\]

The dynamical matrix \( A \) is a time-invariant matrix and its associated transition matrix \( \Phi(\nu, \nu_0) \) for Equation [11] is given by \( \exp(A(\nu - \nu_0)) \). Note that the \( x - y \) in-plane and out-of-plane motions are completely decoupled. Since \( U_{zz} > 0 \), the out-of-plane motion is an harmonic oscillator with pulsation \( \sqrt{U_{zz}} \) while the in-plane trajectories may be bounded or not depending on \( \mu \) and the initial conditions.

### III. A PRIMER VECTOR APPROACH

The primer vector theory is the historical name given to an indirect approach to the design of fuel-optimal
space trajectories. The seminal results dates back to the fifties and the beginning of the sixties with the works of Lawden\textsuperscript{13,14} which have been rigorously confirmed by Neustad in linear\textsuperscript{17} and nonlinear\textsuperscript{18} settings right after. In this paper, since we are interested by designing fuel transfer trajectories around the libration points, only its linear version is used and described. The main lines of the review of the theoretical background of the proposed results are based on the reference\textsuperscript{9}.

III.i Problem statement and its moment counterpart

Designing minimum-fuel transfer trajectories around the libration points in the CR3BP setting naturally boils down to solving the following linear optimal control problem:

**Problem 3.1.** (Optimal control problem)

Find \( \bar{u} \in L_{1,p}([\nu_0, \nu_f], \mathbb{R}^3) \) solution of the optimal control problem:

\[
\inf_{\psi} \|\psi\|_{L^p} = \inf_{\psi} \int_{\nu_0}^{\nu_f} \|\psi(\nu)\|_p d\nu \\
\text{s.t. } X'(\nu) = AX(\nu) + Bu(\nu), \\
\hspace{1cm} X(\nu_0) = X_0, X(\nu_f) = X_f \in \mathbb{R}^n, \\
\hspace{1cm} \forall \nu \in [\nu_0, \nu_f], \nu_0, \nu_f \text{ fixed,}
\]

where \( B^T = [O_{(n-r)\times r} \hspace{1cm} I_r] \). The state vector dimension and the number of inputs in [10] are denoted \( n \) and \( r \), respectively with \( n = 2, r = 1 \) for the out-of-plane case and \( n = 4, r = 1 \) for the in-plane case. The value of the integer \( p \) actually depends on the geometric configuration of the thrusters. For a steerable thruster, \( p = 2 \), while for \( 6 \) unimbabled identical thrusters, \( p = 1 \).

Following the approach from\textsuperscript{17}. Problem 3.1 is now transformed into an equivalent problem of moment by integrating equation [18]. Let us define a fundamental matrix \( \varphi(\nu) \) for Equation [10] (i.e. \( \Phi(\nu, \nu_0) = \varphi(\nu)\varphi^{-1}(\nu_0) \)) and the matrix \( Y(\nu) = \varphi^{-1}(\nu)B = \begin{bmatrix} y_1(\nu) \cdots y_n(\nu) \end{bmatrix}^T \in \mathbb{R}^{n \times r} \), then:

\[
\begin{aligned}
c &= \varphi^{-1}(\nu_f)X(\nu_f) - \varphi^{-1}(\nu_0)X_0 \\
&= \int_{\nu_0}^{\nu_f} \varphi^{-1}(\sigma)B(\sigma)u(\sigma)d\sigma \\
&= \int_{\nu_0}^{\nu_f} Y(\sigma)u(\sigma)d\sigma.
\end{aligned}
\]

It is important to notice for the remainder of the analysis that for the specific matrices \( Y(\nu) \) encountered in the studied problem, \( y_1(\nu) \cdots y_n(\nu) \) are linearly independent elements of \( C([\nu_0, \nu_f], \mathbb{R}^r) \). This will be assumed in the rest of the paper. It follows from [19] that Problem 3.1 can be equivalently written as:

**Problem 3.2.** (Minimum norm moment problem) Find \( \bar{u}(t) \in L_{1,p}([\nu_0, \nu_f], \mathbb{R}^r) \) solution of the minimum norm moment problem:

\[
\begin{aligned}
\inf_{\psi} \|\psi\|_{L^p} &= \inf_{\psi} \int_{\nu_0}^{\nu_f} \|\psi(\nu)\|_p d\nu \\
\text{s.t. } &\int_{\nu_0}^{\nu_f} Y(\sigma)u(\sigma)d\sigma = c, \nu_0, \nu_f \text{ fixed.}
\end{aligned}
\]

It is well-known that Problem 3.2 may not reach its optimal solution due to concentration effects (see the reference\textsuperscript{21}) . It is then necessary to resort to a relaxation scheme by embedding the space \( L_{1,p}([\nu_0, \nu_f], \mathbb{R}^r) \) in the dual space \( C^*([\nu_0, \nu_f], \mathbb{R}^r) \) of the Banach space \( C([\nu_0, \nu_f], \mathbb{R}^r) \).

III.ii A relaxed problem and its SICP formulation

Generalized solutions of the original Problem 3.2 may be obtained as the solutions of the following relaxed moment problem.

**Problem 3.3.** (Relaxed moment problem)

Determine \( \check{g} \in BV([\nu_0, \nu_f], \mathbb{R}^r) \) solution of the following problem:

\[
\begin{aligned}
\inf_{\check{g}} \|\check{g}\|_{L^p} &= \inf_{\check{g}} \max_{\nu \in [\nu_0, \nu_f]} \|g(\nu) - g(\nu_{\nu_0})\|_p \\
\text{s.t. } &\int_{\nu_0}^{\nu_f} Y(\nu)d\nu = c.
\end{aligned}
\]

\( P_\kappa = \{\nu_0 = \nu_1 < \nu_2, \cdots, \nu_{\kappa} = \nu_f\} \) is any finite partition of \( [\nu_0, \nu_f] \). It is shown in\textsuperscript{17} that the infimum of Problem 3.3 is reached and that it is equal to the infimum of Problem 3.2, denoted by \( \check{\eta} \) in what follows. The association between the space \( BV([\nu_0, \nu_f], \mathbb{R}^r) \) and the dual \( C^*([\nu_0, \nu_f], \mathbb{R}^r) \) of the space \( C([\nu_0, \nu_f], \mathbb{R}^r) \) defined in the Riesz Representation Theorem and the use of a duality principle based on the extension form of the Hahn-Banach theorem\textsuperscript{16} are the basic elements for the next result, originally given in\textsuperscript{17} in its complete form and partially in\textsuperscript{15} for particular optimization problems. Here, we follow the lines developed in the textbook of [16, Chapter 5].

**Theorem 1.** Let \( y_i(\nu) \in C([\nu_0, \nu_f], \mathbb{R}^r), \forall i = 1, \cdots, n, \) \( q = \frac{p}{p-1} \) the conjugate index and suppose that

\[
D = \left\{ g \in BV([\nu_0, \nu_f], \mathbb{R}^r) : \int_{\nu_0}^{\nu_f} Y(\nu)d\nu = c \right\},
\]

is a non empty set then

\[
\check{\eta} = \min_{\check{g} \in D} \|\check{g}\|_{L^p} = \max_{\|Y^T(\nu)\lambda\|_q \leq 1} \|C^T\lambda\|.
\]
In addition, let \( \bar{y} \) and \( \bar{\lambda} \) be optimal solutions of [23],
\[
\bar{\lambda} = \arg \left[ \max_{\lambda} \left| y^T(\nu) \right| \right] \text{ and let } \bar{y}(\nu) = \sum_{i=1}^{n} \lambda_i y_i(\nu) = Y^T(\nu) \bar{\lambda} \in \mathbb{R}^r.
\]
Then the optimal \( y \) is such that:
\[
\int_{q_0}^{q_f} \bar{X}(\nu)\bar{y}(\nu) d\nu = \sup_{\nu_0 \leq \nu \leq \nu_f} \left\| \bar{y}(\cdot) \right\|_q \| \bar{y} \|_{tv,p}.
\]  

The two problems defined in Eq. [23] may be considered as dual through the equality of the optimal values of their respective objectives and the relation between their solutions thanks to the condition in Eq. [24]. This results in a significant simplification: The infinite-dimensional optimization Problem 3.3 has been converted to a search of an optimal vector \( \bar{\lambda} \) in a finite-dimensional vector space submitted to a continuum of constraints, yielding a semi-infinite convex problem (SICP):

**Problem 3.4.** (SICP problem) Find \( \bar{\lambda} \in \mathbb{R}^n \) solution of
\[
\bar{\mu} = \min_{\lambda \in \mathbb{R}^n} -\bar{c}^T \lambda \quad \left| Y^T(\nu) \lambda \right|_{tv} \leq 1.
\]

Note that \( \bar{\mu} = -\bar{\eta} \). Once its solution is obtained, the relation [24] between the function \( \bar{y}(\cdot) \) element of the Banach space \( C([\nu_0, \nu_f], \mathbb{R}^r) \) and the optimal vector \( \bar{\mu} \) is particularly important to get back to the optimal bounded variation solution of the relaxed Problem 3.3 as shown in the following result from [17].

**Theorem 2.** Let \( y_i(\cdot) \in C([\nu_0, \nu_f], \mathbb{R}^r) \), \( i = 1, \ldots, n \) and \( \bar{\lambda} \in \mathbb{R}^n \) be an optimal solution of Problem (25). Define the sets \( \Gamma = \{ \bar{\nu} \in [\nu_0, \nu_f] : \| \bar{y}(\bar{\nu}) \|_q = 1 \} \) and \( \bar{\Gamma} = \bigcup_{i=1}^{n} \Gamma \). The compactness of \( \bar{\Gamma} \) for \( p = 1 \). There is an optimal solution \( \bar{\gamma}(\cdot) \in BV ([\nu_0, \nu_f], \mathbb{R}^r) \), which is a step function with at most \( n \) points of discontinuity \( \bar{\nu}_j \in \Gamma, j = 1, \ldots, N \leq n \). Its jumps are given by:
\[
g_\delta(\bar{\nu}_j) - g_\delta(\bar{\nu}_j^-) = \alpha_\delta \text{sgn}(\bar{y}_s(\bar{\nu}_j)) \chi_{\Gamma_j}, \quad \alpha_\delta > 0, \quad \text{when } p = 1,
\]
or
\[
g_\delta(\bar{\nu}_j) - g_\delta(\bar{\nu}_j^-) = \alpha_\delta |\bar{y}_s(\bar{\nu}_j)|^{q-1} \text{sgn}(\bar{y}_s(\bar{\nu}_j)),
\]
for \( 1 < p < \infty \),
\[
\text{when } 1 < p < \infty,
\]
where \( \beta_i(\bar{\nu}_j) \) are given by:
\[
\beta_i(\bar{\nu}_j) = \begin{cases} 
\sum_{s=1}^{r} y_{i,s}(\bar{\nu}_j) sgn(\bar{y}_s(\bar{\nu}_j)), & \text{when } p = 1, \\
\sum_{s=1}^{r} y_{i,s}(\bar{\nu}_j) |\bar{y}_s(\bar{\nu}_j)|^{q-1} sgn(\bar{y}_s(\bar{\nu}_j)), & \text{when } 1 < p < \infty.
\end{cases}
\]
for all \( j = 1, \ldots, N \).

This theorem states important results that have been known for a while in the aerospace community but whose value has not been completely exploited to derive efficient numerical algorithms for impulsive maneuvers design. First, it says that the optimal controlled trajectory for the minimum-fuel linearized optimal control problem is purely impulsive and that the number of impulses is upper-limited by \( n \) which is the dimension of the fixed final conditions of the optimal control problem.

**Remark 1.** It is also shown in [17] that a sequence of functions \( u_{i,\epsilon}(\cdot) \in L_{1,p}([\nu_0, \nu_f], \mathbb{R}^r) \) converges to a linear combination of \( \delta(\cdot) \) functions corresponding to the function \( g(\cdot) \) with equal norms. Let \( \Delta V(\bar{\nu}_j) = \bar{g}(\bar{\nu}_j) - g(\bar{\nu}_j^-) \), then roughly speaking, this may be described by:
\[
u_{\epsilon}(\nu) \to \sum_{j=1}^{N} \Delta V(\bar{\nu}_j) \delta(\bar{\nu}_j - \nu), \quad \epsilon \to 0.
\]

Indeed, the initial optimal control problem amounts to find the sequences of optimal impulse locations \( \{ \bar{\nu}_i \}_{i=1,\ldots,N} \) and optimal impulse vectors \( \{ \Delta V(\bar{\nu}_i) \}_{i=1,\ldots,N} \) verifying the boundary equation:
\[
\epsilon = \sum_{i=1}^{N} Y(\bar{\nu}_i) \Delta V(\bar{\nu}_i).
\]

The vector \( p_\delta(\nu) = Y^T(\nu) \lambda \) involved in [25] is nothing but the primer vector initially defined in the seminal work of Lawden [14]. In this reference, the primer vector \( p_\delta(\nu) \) is defined as the velocity adjoint vector arising from applying the Pontryagin Maximum Principle to optimal trajectory problems or Lagrangian duality as in [5] where the vector \( \lambda \) is the optimal Lagrange multiplier. For an optimal impulsive trajectory, the primer vector \( p_\delta(\nu) \) must satisfy the well-known Lawden’s necessary and sufficient optimality conditions recalled in [4]. In summary, finding the optimal solution to the primal problem is an indirect way of solving the original trajectory problem. As a result, efficient algorithms to solve the former are extremely valuable. The next section proposes a new procedure based on a discretization algorithm for the solution of the semi-infinite programming Problem 3.4.
IV. A MODERN NUMERICAL APPROACH

The resolution of Problem 3.4 when dealing with trajectories around Lagrange points is analogous to the solution of the circular linearized rendezvous 2-body problem\(^1\). Actually, the out-of-plane dynamics being a harmonic oscillator as well, the optimal solution for \(n = 1\) is analytical and identical to the one given by Prussing\(^2\). As for \(n \geq 2\), it can be solved numerically as in Arzelier et al\(^1\). Basically, the primer vector is obtained iteratively by solving a sequence of convex problems in \(\mathbb{R}^2\). An improved heuristic version is proposed in this paper. Moreover, by relying on properties of the 1 and \(\infty\) norms, concatenating the projected in-plane and out-of-plane optimal solution yields the optimal solution for \(n = 3\) when \(p = 1\). Taking advantage of this property makes for a faster resolution under the complete dynamics. This section details the main results at the basis of the algorithm used to compute the minimum-fuel sequence of impulsive maneuvers.

The general, numerical approach proposed in\(^1\) and based on discretization methods or exchange methods reminds the numerical solution of the Chebyshev approximation problem\(^20\). It intends as solving a series of convex problems 3.4 for which the feasible set described by infinitely many constraints is replaced by a discrete approximation involving a finite set of constraints defined on a given grid \(\Theta_i\) of locations in the interval \([\nu_0, \nu_f]\). For \(p = 1\), the discretized version of 3.4 boils down to a linear program:

**Problem 4.1.** Discretized SICP (\(p = 1\))

\[
\max_{\lambda} \ c^T \cdot \lambda \\
\text{s.t.} \quad -1 \leq |Y^T \cdot \lambda(\nu)| \leq 1, \ \forall \ \nu \in \Theta_i, \ \text{for} \ j = 1, \ldots, n,
\]

while for \(p = 2\), it falls into the scope of Semi-Definite Programming:

**Problem 4.2.** Discretized SICP (\(p = 2\))

\[
\max_{\lambda} \ c^T \cdot \lambda \\
\text{s.t.} \quad \left[ \begin{array}{c} 1 \\ Y^T \cdot \lambda(\nu) \\ 1_n \end{array} \right] \geq 0 \ \forall \ \nu \in \Theta_i.
\]

The key point in the method is that, by properly building a sequence of sets \(\Theta_i\), the solutions \(\lambda^{(i)}\) of the discretized problems 4.1 or 4.2 will converge to the solution of 3.4. Its efficiency strongly relies on the fact that Problems 4.1 and 4.2 can be solved efficiently by dedicated methods like simplex or interior-point methods\(^3\).

It has been shown\(^1\) that the sequence of discretized sets \(\Theta_{i+1} = \Theta_i \cup \{ \text{argmax}_{\nu} \| Y^T \cdot \lambda^{(i)}(\nu) \|_q \} \) verifies the necessary assumptions of\(^5\) and converges to the optimal solution of 3.4. In this paper, following incentives from exchange methods\(^2\) and ideas from\(^2\), a heuristic improvement is made by also removing all incentives from exchange methods. In other words, \(\Theta_{i+1} = \{ \nu \in \Theta_i, \| Y^T \cdot \lambda(\nu)\|_q = 1 \} \cup \{ \text{argmax}_{\nu} \| Y^T \cdot \lambda^{(i)}(\nu) \|_q \} \). Also used in\(^10\), this modification leads to a reduction of the number of constraints involved in Problems 4.1 or 4.2. This is summarized in the algorithm below written in pseudo-code.

**Input:** interval \(\Theta = [\nu_0, \nu_f]\), matrix \(Y(\nu)\), initial condition \(c\), accuracy \(\varepsilon\)

**Output:** \(\mu^{(i)}\) and \(\lambda^{(i)}\) numerical solution of Pb. 3.4

**Init:**

\(i \leftarrow 0;\)

\(\text{Convergence} \leftarrow \text{False};\)

\(\Theta_0 \leftarrow \{ \theta_0, \theta_1 \} \subset \Theta \text{ s.t. } \theta_0 - \theta_1 \neq k\pi;\)

**while** Convergence=False **do**

**Find** \(\lambda^{(i)}\) solution of discretized problem:

\[
\mu^{(i)} = \inf_{\lambda \in \mathbb{R}^n} -c^T \lambda \\
\text{s.t.} \quad \| Y^T(\theta_k) \lambda \|_q \leq 1 \quad \text{for all } \theta_k \in \Theta_i
\]

\(\text{if } \text{Convergence=}\text{False then}\)

\(i \leftarrow i + 1;\)

\(\Theta_i \leftarrow \Theta_{i-1} \cup \left\{ \arg \left( \max_{\theta \in \Theta} \| Y^T(\theta) \lambda^{(i)} \|_q \right) \right\};\)

\(\Theta_i \leftarrow \Theta_i - \left\{ \theta_k \| Y(\theta_k) \lambda^{(i)} \|_q < 1 - \varepsilon \right\};\)

**end**

**return** \(\mu^{(i)}, \lambda^{(i)}\).

**Algorithm 1:** Numerical procedure for solving Problem 3.4

Once Algorithm 1 returns the near-optimal value of \(\lambda\), which gives the near-optimal primer vector, Theorem 2 is used to reconstruct the instantaneous velocity increments.

V. RESULTS AND DISCUSSION

The algorithm described in the previous Section has been implemented in Python, making use of the scipy.linprog routine for Linear Programming and cvx solver for Semi-Definite Programming. The practical results presented thereafter concern the \(L_2\) libration point of the Earth-Moon system which is of particular interest as the tentative location of a fuel-depot to be used by interplanetary spacecraft. The method is nonetheless applicable to any Lagrange point in any binary system. A set of initial and final values for the transfer (in- and out-
of-plane), reported in Tables 1, was generated randomly, but could be chosen for specific purposes such as station-keeping or transfer between Halo orbits. Comparisons are twofold: between 1 and 2-norm costs on the one hand and between indirect and direct approaches on the other hand.

For an a priori fixed number of impulsive maneuvers at given locations, a convex problem (LP or SDP) is formulated and solved numerically by discretizing the possible locations of burns. Its solution is therefore suboptimal, depending strongly upon the number of impulses\textsuperscript{12,15}. Introducing slack variables, 1- and 2-norm minimization respectively boils down to a linear and semi-definite program. Due to the greater computational cost of the latter, the grid is chosen to be sparser here e.g. 1000 points against 10000. On the other hand, the indirect approach uses the thinner version when checking the magnitude of the primer vector.

Table 1: Initial and final conditions

<table>
<thead>
<tr>
<th>$\nu_0$ (rad)</th>
<th>$\delta x_0$ (m)</th>
<th>$\delta y_0$ (m)</th>
<th>$\delta z_0$ (m)</th>
</tr>
</thead>
<tbody>
<tr>
<td>3.322</td>
<td>6449.40</td>
<td>65117.03</td>
<td>22814.91</td>
</tr>
<tr>
<td>$\delta \dot{x}_0$ (m/s)</td>
<td>$\delta \dot{y}_0$ (m/s)</td>
<td>$\delta \dot{z}_0$ (m/s)</td>
<td></td>
</tr>
<tr>
<td>-0.0312</td>
<td>0.0392</td>
<td>0.2114</td>
<td></td>
</tr>
</tbody>
</table>

<table>
<thead>
<tr>
<th>$\nu_f$ (rad)</th>
<th>$\delta x_f$ (m)</th>
<th>$\delta y_f$ (m)</th>
<th>$\delta z_f$ (m)</th>
</tr>
</thead>
<tbody>
<tr>
<td>4.737</td>
<td>59066.09</td>
<td>67728.64</td>
<td>84015.47</td>
</tr>
<tr>
<td>$\delta \dot{x}_f$ (m/s)</td>
<td>$\delta \dot{y}_f$ (m/s)</td>
<td>$\delta \dot{z}_f$ (m/s)</td>
<td></td>
</tr>
<tr>
<td>-0.1087</td>
<td>0.1616</td>
<td>-0.1730</td>
<td></td>
</tr>
</tbody>
</table>

Figure 3 depicts the primer vector history for $p = 1$ whereas Figure 2 is for the fuel-consumption, the relative positions and velocities obtained via the indirect approach. The corresponding trajectory is showed in Figure 4. The same plots for $p = 2$ are respectively on Figures 6, 5 and 7. Tables 2 and 3 contain the control laws respectively for $p = 1$ and 2. The optimality of the solutions is certified by the norm of the primer vector being always less or equal to one, the latter occurring exclusively at times of burns.

Table 2: Location and components of optimal impulses for $p = 1$, $N^* = 4$

<table>
<thead>
<tr>
<th>$\nu$ (rad)</th>
<th>$\Delta V_x$ (m/s)</th>
<th>$\Delta V_y$ (m/s)</th>
<th>$\Delta V_z$ (m/s)</th>
<th>$|\Delta V|_1$ (m/s)</th>
</tr>
</thead>
<tbody>
<tr>
<td>3.322</td>
<td>0.0126</td>
<td>0</td>
<td>0</td>
<td>0.0126</td>
</tr>
<tr>
<td>3.987</td>
<td>0</td>
<td>0.05530</td>
<td>0.3617</td>
<td>0.05530</td>
</tr>
<tr>
<td>4.030</td>
<td>0</td>
<td>0</td>
<td>0</td>
<td>0.01570</td>
</tr>
<tr>
<td>4.737</td>
<td>0</td>
<td>0.1570</td>
<td>0</td>
<td>0.5530</td>
</tr>
</tbody>
</table>

Table 3: Location and components of optimal impulses for $p = 2$, $N^* = 2$

<table>
<thead>
<tr>
<th>$\nu$ (rad)</th>
<th>$\Delta V_x$ (m/s)</th>
<th>$\Delta V_y$ (m/s)</th>
<th>$\Delta V_z$ (m/s)</th>
<th>$|\Delta V|_2$ (m/s)</th>
</tr>
</thead>
<tbody>
<tr>
<td>3.928</td>
<td>-0.0181</td>
<td>0.5173</td>
<td>0.1541</td>
<td>0.5401</td>
</tr>
<tr>
<td>4.737</td>
<td>-0.5677</td>
<td>0.4595</td>
<td>0.0165</td>
<td>0.6850</td>
</tr>
</tbody>
</table>

The minimal 1- and 2-norm trajectories obtained via the indirect approach are visually very similar. The main difference comes from the fact that the minimal 1-norm
control has more burns i.e. four rather than two. This can be explained by the fact that it is more efficient with $p = 2$ to minimize the number of maneuvers. Here one can see that somehow the two middle impulses of $p = 1$ are merged into a single one for $p = 2$. This also translates into an optimal solution for the 2-norm with no zero component, unlike for the 1-norm. Moreover, the mid-course maneuvers exhibited by both minimal control laws is a clear setback for the analytic yet generally sub-optimal 2-impulse method that consists in designing an initial maneuver to achieve transfer in position at the final location where the second and last impulse matches the desired velocity. The fuel-optimal solution obtained for $p = 2$ in particular shows the advantage of an initial coasting arc.

Table 4: Comparison between direct and indirect approaches for $p = 1$

<table>
<thead>
<tr>
<th>Approach</th>
<th>Cost (m/s)</th>
<th>Computation time (s)</th>
</tr>
</thead>
<tbody>
<tr>
<td>Indirect</td>
<td>1.6384</td>
<td>0.8447</td>
</tr>
<tr>
<td>Direct</td>
<td>1.6384</td>
<td>1.3974</td>
</tr>
</tbody>
</table>

Table 5: Comparison between direct and indirect approaches for $p = 2$

<table>
<thead>
<tr>
<th>Approach</th>
<th>Cost (m/s)</th>
<th>Computation time (s)</th>
</tr>
</thead>
<tbody>
<tr>
<td>Indirect</td>
<td>1.2251</td>
<td>1.3731</td>
</tr>
<tr>
<td>Direct</td>
<td>1.2251</td>
<td>160.5454</td>
</tr>
</tbody>
</table>

Objective functions and timings are reported for direct and indirect approaches in Table 4 for the 1-norm and Table 5 for the 2-norm. For the two cases, both approaches return an identical fuel-consumption. However, due to need for a fine locations grid for the possible burns, the direct method requires in general a larger computing effort.

VI. CONCLUSIONS

In this paper, a modern approach to the linear primer vector theory has been proposed for fuel-optimal trajectories near any Lagrange point, extending the previous work of the authors with the linearized restricted elliptical 2-body problem. It combines recent results that are both analytical and numerical to provide the user with a fast, efficient way of computing the optimal burns. Output of a sequence of linear or semi-definite programs (depending on the geometrical configuration of the thrusters), the minimal consumption can be evaluated without the actual need of calculating the maneuvers. The latter, which cannot be achieved by a direct approach, turns this method into a powerful tool for mission analysis. Still, the convergence of the used algorithm remains an open problem to be studied in coming developments. Future work includes also considering the elliptical case i.e. the non-circular generalization of the Lagrange points. Perspectives are to initialize an indirect solving approach to the control problem under the non-linear dynamics with the solution of the linearized version.

REFERENCES


