Pole optimization for exponential integration

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Contents

1 Introduction

This example is concered with the computation of a family of common-denominator rational approximants for the two-parameter function $\exp(-tz)$ using RKFIT [2, 3]. This corresponds to the example from [3, Section 6.2]. Let us consider the problem of solving a linear constant-coefficient initial-value problem

$$\boldsymbol{u}'(t) + L\boldsymbol{u}(t) = \boldsymbol{0}, \quad \boldsymbol{u}(0) = \boldsymbol{u}_0,$$

at several time points $t_1, \ldots, t_\ell$. The exact solutions $\boldsymbol{u}(t_j)$ are given in terms of the matrix exponential as $\boldsymbol{u}(t_j) = \exp(-t_j L)\boldsymbol{u}_0$. A popular approach for approximating $\boldsymbol{u}(t_j)$ is to use rational functions $r^{[j]}$ of the form

$$r^{[j]}(z) = \frac{\sigma_1^{[j]}}{\xi_1 - z} + \frac{\sigma_2^{[j]}}{\xi_2 - z} + \cdots + \frac{\sigma_m^{[j]}}{\xi_m - z},$$

constructed so that $r^{[j]}(L)\boldsymbol{u}_0 \approx \boldsymbol{u}(t_j)$. Note that the poles of $r^{[j]}$ do not depend on t_j and we have

$$r^{[j]}(L)\boldsymbol{u}_0 = \sum_{i=1}^{m} \sigma_i^{[j]} (\xi_i I - L)^{-1} \boldsymbol{u}_0,$$

the evaluation of which amounts to the solution of m linear systems. Such common-pole approximants have great computational advantage, in particular, in combination with direct solvers (as the LU factorizations of $\xi_i I - L$ can be reused for each t_j) and when the linear systems are assigned to parallel processors.

2 Surrogate approach

In order to use RKFIT for finding "good" poles $\xi_1, \ldots, \xi_m$ of the rational functions $r^{[j]}$, we propose a surrogate approach similar to that in [4]. Let $A = \mathrm{diag}(\lambda_1, \ldots, \lambda_N)$ be a diagonal matrix whose eigenvalues are a "sufficiently dense" discretization of the positive semiaxis $\lambda \geq 0$. In this example we take $N = 500$ logspaced eigenvalues in the interval $[10^{-6}, 10^6]$. Further, we define $\ell = 41$ logspaced time points t_j in the interval $[10^{-1}, 10^1]$, and the matrices $F^{[j]} = \exp(-t_j A)$. We also define $\boldsymbol{b} = [1, \ldots, 1]^T$ to assign equal weight to each eigenvalue of A.

```
N  = 500;
ee = [0 , logspace(-6, 6, N-1)];
A  = spdiags(ee(:), 0, N, N);
b  = ones(N, 1);
t  = logspace(-1, 1, 41);
for j = 1:length(t)
  F{j} = spdiags(exp(-t(j)*ee(:)), 0, N, N);
end
```

We then run the RKFIT algorithm for finding a family of rational functions $r^{[j]}$ of type $(m-1, m)$ with $m = 12$ so that $\|F^{[j]}\boldsymbol{b} - r^{[j]}(A)\boldsymbol{b}\|_2$ is minimized for all $j = 1, \ldots, \ell$.

```
m = 12; k = -1;                % type (11, 12)
xi = inf(1,m);                 % initial poles at infinity
param.k = k;                   % subdiagonal approximant
param.maxit = 6;               % at most 6 RKFIT iterations
param.tol   = 0;               % exactly 6 iterations
param.real  = 1;               % data is real-valued
[xi, ratfun, misfit, out] = rkfit(F, A, b, xi, param);
```

3 The RKFIT outputs

The first output argument of RKFIT is a vector `xi` collecting the poles $\xi_1, \ldots, \xi_m$ of the rational Krylov space. The second output `ratfun` is a cell array each cell of which is a `rkfun`, a datatype representing a rational function. All rational functions in this cell array share the same denominator with roots $\xi_1, \ldots, \xi_m$. The next output parameter is a vector containing the computed relative misfit after each RKFIT iteration. The relative misfit is defined as (cf. eq. (1.5) in [3])

$$\mathrm{misfit} = \sqrt{\frac{\sum_{j=1}^{\ell} \|F^{[j]}\boldsymbol{b} - r^{[j]}(A)\boldsymbol{b}\|_F^2}{\sum_{j=1}^{\ell} \|F^{[j]}\boldsymbol{b}\|_F^2}}.$$

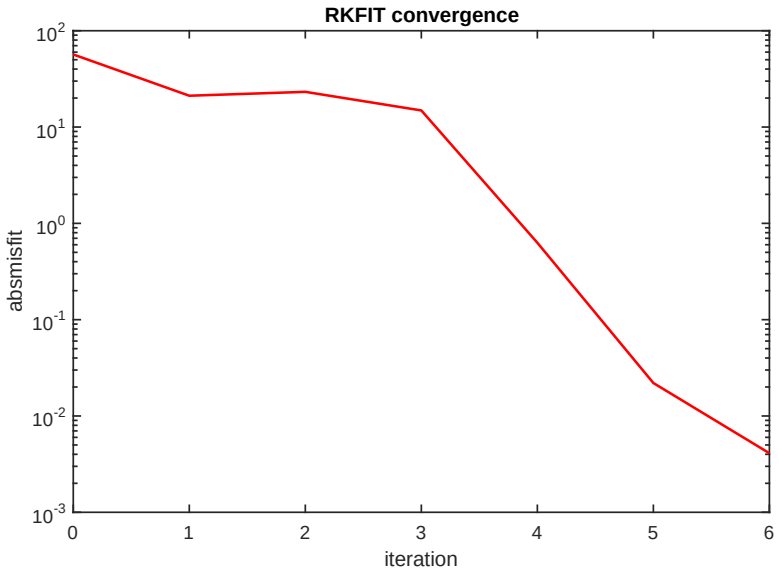
We can easily verify that the last entry of `misfit` indeed corresponds to this formula:

```
num = 0; den = 0;
for j = 1:length(ratfun)
  num = num + norm(F{j}*b - ratfun{j}(A,b), 'fro')^2;
  den = den + norm(F{j}*b, 'fro')^2;
end
disp([misfit(end) sqrt(num/den)])
```

```
   4.2516e-05   4.2516e-05
```

Here is a plot of the `misfit` vector, giving an idea of the RKFIT convergence:

```
figure
semilogy(0:6, [out.misfit_initial, misfit]*sqrt(den), 'r-');
xlabel('iteration');
ylabel('absmisfit')
title('RKFIT convergence')
```



4 Verifing the accuracy

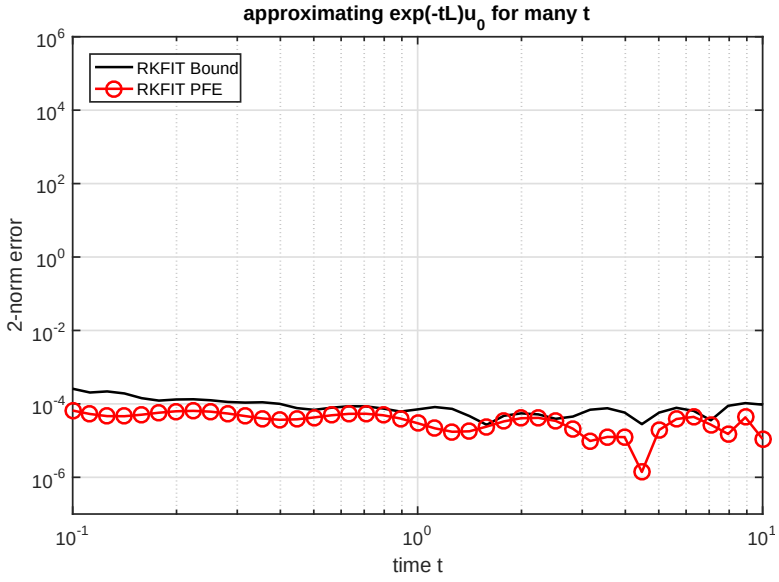
To evaluate the quality of the common-denominator rational approximants for all $\ell = 41$ time points t_j, we perform an experiment similar to that in [5, Figure 6.1] by approximating $\boldsymbol{u}(t_j) = \exp(-t_j L)\boldsymbol{u}_0$ and comparing the result to MATLAB's `expm`. Here, L is a 841×841 finite-difference discretization of the scaled 2D Laplace operator -0.02Δ on the domain $[-1, 1]^2$ with homogeneous Dirichlet boundary condition, and $\boldsymbol{u}_0$ corresponds to the discretization of $u_0(x, y) = (1 - x^2)(1 - y^2)e^x$ on that domain.

```
% Parts of the following code have been taken from [5].
J = 30; h = 2/J; s = (-1+h:h:1-h)';  % in [3,5] J = 50 is used
[xx,yy] = meshgrid(s,s);             % 2D grid
x = xx(:); y = yy(:);                % 2D grid stretched to 1D
L = 0.02*gallery('poisson',J-1)/h^2; % 2D Laplacian
v = (1-x.^2).*(1-y.^2).*exp(x);      % initial condition
v = v/norm(v);
```

```
for j = 1:length(t)
  exac(:,j) = expm(-t(j)*L)*v;
  rat = ratfun{j}(L,v);
  err_rat(j) = norm(rat - exac(:,j));
  bnd(j) = norm(ratfun{j}(A,b) - F{j}*b,inf);
end
```

We now plot the error $\|\boldsymbol{u}(t_j) - r^{[j]}(L)\boldsymbol{u}_0\|_2$ for each time point t_j (curve with red circles), together with the approximate upper error bound $\|\exp(-t_j A)\boldsymbol{b} - r^{[j]}(A)\boldsymbol{b}\|_\infty$ (black curve), which can be easily computed by direct evaluation. We find that the error is indeed approximately uniform and smaller than 1.1×10^{-4} over the time interval $[10^{-1}, 10^{1}]$.

```
figure
loglog(t, bnd, 'k-')
hold on
loglog(t, err_rat, 'r-o')
xlabel('time t'); ylabel('2-norm error')
legend('RKFIT Bound', 'RKFIT PFE', 'Location', 'NorthWest')
title('approximating exp(-tL)u_0 for many t')
grid on
axis([0.1, 10, 1e-7, 1e6])
```



5 Conversion to partial fraction form

When evaluating the rational functions $r^{[j]}$ on a parallel computer, it is convenient to have their partial fraction expansions at hand. The `rkfun` class provides a method called `residue` for this purpose. This method supports the use of MATLAB's variable precision (VPA) capabilities, or the Advanpix Multiple Precision (MP) toolbox [1].

For example, here are the residues and poles of the first rational function $r^{[1]}$ corresponding to $\exp(-t_1 A)\boldsymbol{b}$:

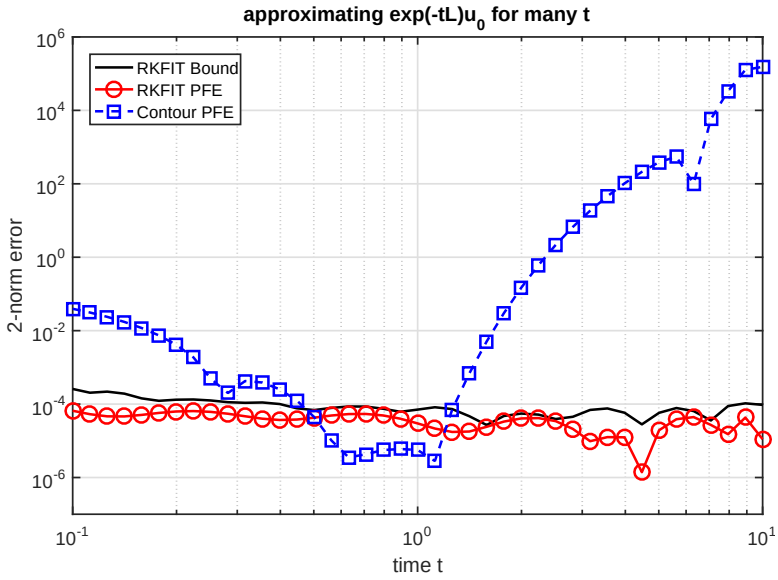
```
try mp(1); catch e, try mp=@(x)vpa(x); mp(1); catch e, mp=@(x)x;
warning('MP & VPA are unavailable. Using double.'); end, end

[resid, xi, absterm, cnd, pf] = residue(mp(ratfun{1}));
disp(double([resid , xi]))
```

```
   1.0366e-01 - 5.2768e-03i  -3.6241e-01 + 1.9079e-01i
   1.0366e-01 + 5.2768e-03i  -3.6241e-01 - 1.9079e-01i
   1.1852e-01 + 6.1657e-04i  -2.3171e-01 + 6.3413e-01i
   1.1852e-01 - 6.1657e-04i  -2.3171e-01 - 6.3413e-01i
   2.5977e-01 - 3.7616e-02i  -3.8020e-02 + 1.6325e+00i
   2.5977e-01 + 3.7616e-02i  -3.8020e-02 - 1.6325e+00i
   4.9256e-01 - 4.4575e-01i   7.0879e-01 + 4.5038e+00i
   4.9256e-01 + 4.4575e-01i   7.0879e-01 - 4.5038e+00i
  -2.0006e-01 - 1.1931e+00i   4.2572e+00 + 1.1415e+01i
  -2.0006e-01 + 1.1931e+00i   4.2572e+00 - 1.1415e+01i
  -7.5452e-01 + 2.8741e-01i   1.8285e+01 + 2.5425e+01i
  -7.5452e-01 - 2.8741e-01i   1.8285e+01 - 2.5425e+01i
```

6 Comparison with contour-based approach

We now compare RKFIT with the accuracy of the contour-based rational approximants derived in [5]. As discussed there, this approach leads to approximants which are very accurate near $t \approx 1$, but their accuracy degrades rapidly as one moves away from this parameter.

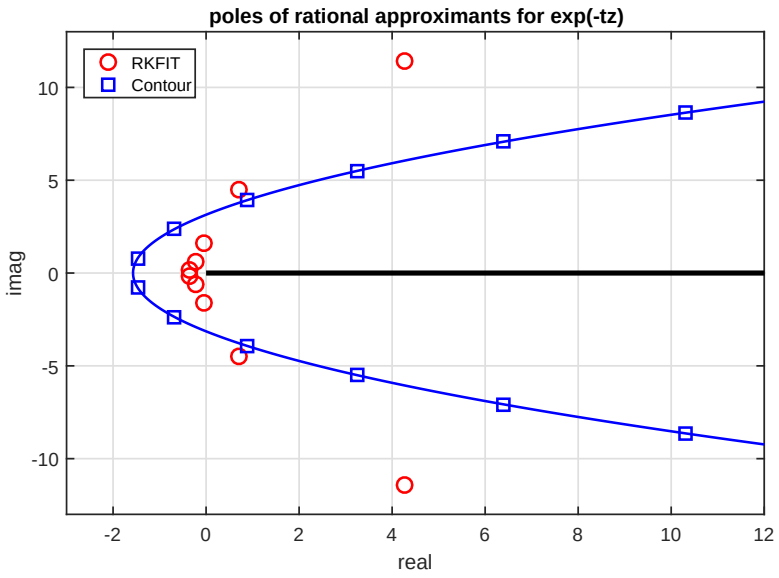
```
% Contour integral code from [5].
NN = 12; theta = pi*(1:2:NN-1)/NN;    % quad pts in (0, pi)
z  = NN*(.1309-.1194*theta.^2);       % quad pts on contour
z  = z + NN*.2500i*theta;
w  = NN*(-.1194*2*theta+.2500i);      % derivatives
for j = 1:length(t)
  c = (1i/NN)*exp(t(j)*z).*w;         % quadrature weights
  appr = zeros(size(v));
  for k = 1:NN/2,                     % sparse linear solves
    appr = appr - c(k)*((z(k)*speye(size(L))+L)\v);
  end
  appr = 2*real(appr);                % exploit symmetry
  err_cont(j) = norm(appr-exac(:,j));
end
loglog(t,err_cont,'b--s')
legend('RKFIT Bound', 'RKFIT PFE', 'Contour PFE', ...
       'Location', 'NorthWest')
```



7 Plot of the poles

Finally, the $m = 12$ poles of the rational functions $r^{[j]}$ are shown in the following plot. We can see that the "optimal" RKFIT poles do not seem to lie on a parabolic contour.

```
figure
hh1 = plot(xi, 'ro');
axis([-3, 12, -13, 13])
hold on

% Also plot the contour.
theta = linspace(0, 2*pi, 300);
zz = -NN*(.1309-.1194*theta.^2+.2500i*theta);
plot(zz, 'b-')
plot(conj(zz), 'b-')
hh2 = plot(-[z, conj(z)], 'bs');
plot([0, 1e3], [0, 0], 'k-', 'LineWidth', 3)
xlabel('real'), ylabel('imag')
title('poles of rational approximants for exp(-tz)')
grid on
legend([hh1, hh2], 'RKFIT', 'Contour', 'Location', 'NorthWest')
```



8 References

[1] Advanpix LLC., *Multiprecision Computing Toolbox for MATLAB,* ver 3.8.3.8882, Tokyo, Japan, 2015. http://www.advanpix.com/.

[2] M. Berljafa and S. Güttel. *A Rational Krylov Toolbox for MATLAB,* MIMS EPrint 2014.56 (http://eprints.ma.man.ac.uk/2390/), Manchester Institute for Mathematical Sciences, The University of Manchester, UK, 2014.

[3] M. Berljafa and S. Güttel. *The RKFIT algorithm for nonlinear rational approximation,* SIAM J. Sci. Comput., 39(5):A2049–A2071, 2017.

[4] R.-U. Börner, O. G. Ernst, and S. Güttel. *Three-dimensional transient electromagnetic modeling using rational Krylov methods,* Geophys. J. Int., 202(3):2025–2043, 2015. Available also as MIMS EPrint 2014.36 (http://eprints.ma.man.ac.uk/2219/).

[5] L. N. Trefethen, J. A. C. Weideman, and T. Schmelzer. *Talbot quadratures and rational approximations,* BIT, 46(3):653–670, 2006.