

Forces between magnets and multipole arrays of magnets

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November 23, 2009

magnetforces

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1. About this file. This is a ‘literate programming’ approach to writing Matlab code using MATLABWEB¹. To be honest I don’t know if it’s any better than simply using the Matlab programming language directly. The big advantage for me is that you have access to the entire L^AT_EX document environment, which gives you access to vastly better tools for cross-referencing, maths typesetting, structured formatting, bibliography generation, and so on.

The downside is obviously that you miss out on Matlab’s IDE with its integrated M-Lint program, debugger, profiler, and so on. Depending on one’s work habits, this may be more or less of limiting factor to using literate programming in this way.

¹<http://tug.ctan.org/pkg/matlabweb>

2. This work consists of the source file `magnetforces.web` and its associated derived files. It is released under the Apache License v2.0.²

This means, in essence, that you may freely modify and distribute this code provided that you acknowledge your changes to the work and retain my copyright. See the License text for the specific language governing permissions and limitations under the License.

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3. Calculating forces between magnets. This is the source of some code to calculate the forces and/or stiffnesses between two cuboid-shaped magnets with arbitrary displacements and magnetisation direction. (A cuboid is like a three dimensional rectangle; its faces are all orthogonal but may have different side lengths.)

4. The main function is `magnetforces`, which takes three mandatory arguments: `magnet_fixed`, `magnet_float`, and `displ`. These will be described in more detail below.

Optional string arguments may be any combination of 'force', and/or 'stiffness' to indicate which calculations should be output. If no calculation is specified, 'force' is the default.

Inputs:	<code>magnet_fixed</code>	structure describing first magnet
	<code>magnet_float</code>	structure describing the second magnet
	<code>displ</code>	displacement between the magnets
	<code>[what to calculate]</code>	'force' and/or 'stiffness'
Outputs:	<code>forces</code>	forces on the second magnet
	<code>stiffnesses</code>	stiffnesses on the second magnet
Magnet properties:	<code>dim</code>	size of each magnet
	<code>magn</code>	magnetisation magnitude
	<code>magdir</code>	magnetisation direction

```

<magnetforces.m 4> ≡
function [varargout]=magnetforces(magnet_fixed, magnet_float, displ, varargin)
    < Matlab help text (forces) 35>
    < Parse calculation args 9>
    < Initialise main variables 6>
    < Precompute rotation matrices 31>
    < Decompose orthogonal superpositions 7>
    < Calculate everything 12>
    < Combine results and exit 10>
    < Functions for calculating forces and stiffnesses 19>
end

```

²<http://www.apache.org/licenses/LICENSE-2.0>

5. Variables and data structures.

6. First of all, address the data structures required for the input and output. Because displacement of a single magnet has three components, plus sizes of the faces another three, plus magnetisation strength and direction (two) makes nine in total, we use one of Matlab's structures to pass the information into the function. Otherwise we'd have an overwhelming number of input arguments.

The input variables *magnet.dim* should be the entire side lengths of the magnets; these dimensions are halved when performing all of the calculations. (Because that's just how the maths is.)

We use spherical coordinates to represent magnetisation angle, where *phi* is the angle from the horizontal plane ($-\pi/2 \leq \phi \leq \pi/2$) and θ is the angle around the horizontal plane ($0 \leq \theta \leq 2\pi$). This follows Matlab's definition; other conventions are commonly used as well. Remember:

$$\begin{aligned}(1, 0, 0)_{\text{cartesian}} &\equiv (0, 0, 1)_{\text{spherical}} \\ (0, 1, 0)_{\text{cartesian}} &\equiv (\pi/2, 0, 1)_{\text{spherical}} \\ (0, 0, 1)_{\text{cartesian}} &\equiv (0, \pi/2, 1)_{\text{spherical}}\end{aligned}$$

⟨ Initialise main variables 6 ⟩ \equiv

```
size1 = reshape(magnet_fixed.dim/2, [3 1]);
size2 = reshape(magnet_float.dim/2, [3 1]);
J1r = magnet_fixed.magn;
J2r = magnet_float.magn;
J1t = magnet_fixed.magdir(1);
J2t = magnet_float.magdir(1);
J1p = magnet_fixed.magdir(2);
J2p = magnet_float.magdir(2);
```

See also sections 18 and 28.

This code is used in section 4.

7. Superposition is used to turn an arbitrary magnetisation angle into a set of orthogonal magnetisations.

Each magnet can potentially have three components, which can result in up to nine force calculations for a single magnet.

We don't use Matlab's `sph2cart` here, because it doesn't calculate zero accurately (because it uses radians and $\cos(\pi/2)$ can only be evaluated to machine precision rather than symbolically).

```

⟨Decompose orthogonal superpositions 7⟩ ≡
    displ = reshape(displ, [3 1]);      % column vector
    J1 = [J1r*cosd(J1p)*cosd(J1t); ...
          J1r*cosd(J1p)*sind(J1t); ...
          J1r*sind(J1p)];
    J2 = [J2r*cosd(J2p)*cosd(J2t); ...
          J2r*cosd(J2p)*sind(J2t); ...
          J2r*sind(J2p)];

```

This code is used in section 4.

8. Wrangling user input and output.

9. We now have a choice of calculations to take based on the user input. Take the opportunity to bail out in case the user has requested more calculations than provided as outputs to the function.

This chunk and the next are used in both `magnetforces.m` and `multipoleforces.m`.

```
< Parse calculation args 9 > ≡
Nvarargin = length(varargin);
if ( Nvarargin ≠ 0 ∧ ∧ Nvarargin ≠ nargout )
    error('Must have as many outputs as calculations requested.')
end
calc_force_bool = false;
calc_stiffness_bool = false;
if Nvarargin ≡ 0
    calc_force_bool = true;
else
    for ii = 1 : Nvarargin
        switch varargin{ii}
            case 'force'
                calc_force_bool = true;
            case 'stiffness'
                calc_stiffness_bool = true;
            otherwise
                error(['Unknown calculation option ', varargin{ii}, ''])
            end
        end
    end
end
```

This code is used in sections 4 and 59.

10. After all of the calculations have occurred, they're placed back into `varargout`.

```
< Combine results and exit 10 > ≡
if Nvarargin ≡ 0
    varargout{1} = forces_out;
else
    for ii = 1 : Nvarargin
        switch varargin{ii}
            case 'force'
                varargout{ii} = forces_out;
            case 'stiffness'
                varargout{ii} = stiffnesses_out;
            end
        end
    end
end
```

This code is used in sections 4 and 59.

11. The actual mechanics.

12. The expressions we have to calculate the forces assume a fixed magnet with positive z magnetisation only. Secondly, magnetisation direction of the floating magnet may only be in the positive z - or y -directions.

The parallel forces are more easily visualised; if $J1z$ is negative, then transform the coordinate system so that up is down and down is up. Then proceed as usual and reverse the vertical forces in the last step.

The orthogonal forces require reflection and/or rotation to get the displacements in a form suitable for calculation.

Initialise a 9×3 array to store each force component in each direction, and then fill 'er up.

```
< Calculate everything 12 > ≡  
  < Print diagnostics 13 >  
  < Calculate  $x$  15 >  
  < Calculate  $y$  16 >  
  < Calculate  $z$  14 >  
  < Combine calculations 17 >
```

This code is used in section 4.

13. Let's print some information to the terminal to aid debugging. This is especially important (for me) when looking at the rotated coordinate systems.

```
< Print diagnostics 13 > ≡  
  debug_disp('␣␣')  
  debug_disp('CALCULATING␣THINGS')  
  debug_disp('=====')  
  debug_disp('Displacement:')  
  debug_disp(displ')  
  debug_disp('Magnetisations:')  
  debug_disp(J1')  
  debug_disp(J2')
```

This code is used in section 12.

14. The easy one first, where our magnetisation components align with the direction expected by the force functions.

⟨ Calculate z 14 ⟩ \equiv

```

if calc_force_bool
    debug_disp('z-z_force:')
    force_components(9, :) = forces_calc_z_z(size1, size2, displ, J1, J2);
    debug_disp('z-y_force:')
    force_components(8, :) = forces_calc_z_y(size1, size2, displ, J1, J2);
    debug_disp('z-x_force:')
    force_components(7, :) = forces_calc_z_x(size1, size2, displ, J1, J2);
end
if calc_stiffness_bool
    debug_disp('z-z_stiffness:')
    stiffness_components(9, :) = stiffnesses_calc_z_z(size1, size2, displ, J1,
        J2);
    debug_disp('z-y_stiffness:')
    stiffness_components(8, :) = stiffnesses_calc_z_y(size1, size2, displ, J1,
        J2);
    debug_disp('z-x_stiffness:')
    stiffness_components(7, :) = stiffnesses_calc_z_x(size1, size2, displ, J1,
        J2);
end

```

This code is used in section 12.

15. The other forces (i.e., x and y components) require a rotation to get the magnetisations correctly aligned. In the case of the magnet sizes, the lengths are just flipped rather than rotated (in rotation, sign is important). After the forces are calculated, rotate them back to the original coordinate system.

⟨ Calculate x 15 ⟩ \equiv

```

size1_rot = swap_x_z(size1);
size2_rot = swap_x_z(size2);
d_rot = rotate_x_to_z(displ);
J1_rot = rotate_x_to_z(J1);
J2_rot = rotate_x_to_z(J2);
if calc_force_bool
    debug_disp('Forces_x-x:')
    forces_x_x = forces_calc_z_z(size1_rot, size2_rot, d_rot, J1_rot, J2_rot);
    force_components(1, :) = rotate_z_to_x(forces_x_x);
    debug_disp('Forces_x-y:')
    forces_x_y = forces_calc_z_y(size1_rot, size2_rot, d_rot, J1_rot, J2_rot);
    force_components(2, :) = rotate_z_to_x(forces_x_y);
    debug_disp('Forces_x-z:')
    forces_x_z = forces_calc_z_y(size1_rot, size2_rot, d_rot, J1_rot, J2_rot);
    force_components(3, :) = rotate_z_to_x(forces_x_z);
end
if calc_stiffness_bool
    debug_disp('x-z_stiffness:')
    stiffness_components(3, :) = rotate_z_to_x(stiffnesses_calc_z_x(size1_rot,
        size2_rot, d_rot, J1_rot, J2_rot));
    debug_disp('x-y_stiffness:')
    stiffness_components(2, :) = rotate_z_to_x(stiffnesses_calc_z_y(size1_rot,
        size2_rot, d_rot, J1_rot, J2_rot));
    debug_disp('x-x_stiffness:')
    stiffness_components(1, :) = rotate_z_to_x(stiffnesses_calc_z_z(size1_rot,
        size2_rot, d_rot, J1_rot, J2_rot));
end

```

This code is used in section 12.

16. Same again, this time making y the ‘up’ direction.

⟨ Calculate y 16 ⟩ \equiv

```

size1_rot = swap_y_z(size1);
size2_rot = swap_y_z(size2);
d_rot = rotate_y_to_z(displ);
J1_rot = rotate_y_to_z(J1);
J2_rot = rotate_y_to_z(J2);
if calc_force_bool
    debug_disp('Forces_y-x:')
    forces_y_x = forces_calc_z_x(size1_rot, size2_rot, d_rot, J1_rot, J2_rot);
    force_components(4, :) = rotate_z_to_y(forces_y_x);
    debug_disp('Forces_y-y:')
    forces_y_y = forces_calc_z_z(size1_rot, size2_rot, d_rot, J1_rot, J2_rot);
    force_components(5, :) = rotate_z_to_y(forces_y_y);
    debug_disp('Forces_y-z:')
    forces_y_z = forces_calc_z_y(size1_rot, size2_rot, d_rot, J1_rot, J2_rot);
    force_components(6, :) = rotate_z_to_y(forces_y_z);
end
if calc_stiffness_bool
    debug_disp('y-z_stiffness:')
    stiffness_components(6, :) = rotate_z_to_y(stiffnesses_calc_z_y(size1_rot,
        size2_rot, d_rot, J1_rot, J2_rot));
    debug_disp('y-y_stiffness:')
    stiffness_components(5, :) = rotate_z_to_y(stiffnesses_calc_z_z(size1_rot,
        size2_rot, d_rot, J1_rot, J2_rot));
    debug_disp('y-x_stiffness:')
    stiffness_components(4, :) = rotate_z_to_y(stiffnesses_calc_z_x(size1_rot,
        size2_rot, d_rot, J1_rot, J2_rot));
end

```

This code is used in section 12.

17. Finally sum all the components in each direction to get the total forces.

⟨ Combine calculations 17 ⟩ \equiv

```

if calc_force_bool
    forces_out = sum(force_components);
end
if calc_stiffness_bool
    stiffnesses_out = sum(stiffness_components);
end

```

This code is used in section 12.

18. You might have noticed that the initialisation of the *force_components* (and other) variables has not yet been listed. That's because the code is boring.

⟨ Initialise main variables 6 ⟩ +≡

```
if calc_force_bool
    force_components = repmat(NaN, [9 3]);
end
if calc_stiffness_bool
    stiffness_components = repmat(NaN, [9 3]);
end
```

19. Functions for calculating forces and stiffnesses. The calculations for forces between differently-oriented cuboid magnets are all directly from the literature. The stiffnesses have been derived by differentiating the force expressions, but that's the easy part.

```

⟨ Functions for calculating forces and stiffnesses 19 ⟩ ≡
  ⟨ Parallel magnets force calculation 20 ⟩
  ⟨ Orthogonal magnets force calculation 21 ⟩
  ⟨ Parallel magnets stiffness calculation 24 ⟩
  ⟨ Orthogonal magnets stiffness calculation 25 ⟩
  ⟨ Helper functions 32 ⟩

```

This code is used in section 4.

20. The expressions here follow directly from Akoun and Yonnet [1].

Inputs:	<i>size1</i> =(<i>a</i> , <i>b</i> , <i>c</i>)	the half dimensions of the fixed magnet
	<i>size2</i> =(<i>A</i> , <i>B</i> , <i>C</i>)	the half dimensions of the floating magnet
	<i>displ</i> =(<i>dx</i> , <i>dy</i> , <i>dz</i>)	distance between magnet centres
	(<i>J</i> , <i>J2</i>)	magnetisations of the magnet in the z-direction
Outputs:	<i>forces_xyz</i> =(<i>Fx</i> , <i>Fy</i> , <i>Fz</i>)	Forces of the second magnet

```

⟨ Parallel magnets force calculation 20 ⟩ ≡
  function calc_out = forces_calc_z_z(size1, size2, offset, J1, J2)
    J1 = J1(3);
    J2 = J2(3);
    ⟨ Initialise subfunction variables 27 ⟩
    component_x = ...
    +0.5*(v.^2 - w.^2).*log(r - u)...
    +u.*v.*log(r - v)...
    +v.*w.*atan2(u.*v, r.*w)...
    +0.5*r.*u;
    component_y = ...
    +0.5*(u.^2 - w.^2).*log(r - v)...
    +u.*v.*log(r - u)...
    +u.*w.*atan2(u.*v, r.*w)...
    +0.5*r.*v;
    component_z = ...
    -u.*w.*log(r - u)...
    -v.*w.*log(r - v)...
    +u.*v.*atan2(u.*v, r.*w)...
    -r.*w;
    ⟨ Finish up 29 ⟩

```

This code is used in section 19.

21. Orthogonal magnets forces given by Yonnet and Allag [2].

⟨ Orthogonal magnets force calculation 21 ⟩ ≡

```

function calc_out = forces_calc_z_y(size1, size2, offset, J1, J2)
    J1 = J1(3);
    J2 = J2(2);
    ⟨ Initialise subfunction variables 27 ⟩
    component_x = ...
    -multiply_x_log_y(v.*w, r-u)...
    +multiply_x_log_y(v.*u, r+w)...
    +multiply_x_log_y(u.*w, r+v)...
    -0.5*u.^2.*atan1(v.*w, u.*r)...
    -0.5*v.^2.*atan1(u.*w, v.*r)...
    -0.5*w.^2.*atan1(u.*v, w.*r);
    component_y = ...
    0.5*multiply_x_log_y(u.^2-v.^2, r+w)...
    -multiply_x_log_y(u.*w, r-u)...
    -u.*v.*atan1(u.*w, v.*r)...
    -0.5*w.*r;
    component_z = ...
    0.5*multiply_x_log_y(u.^2-w.^2, r+v)...
    -multiply_x_log_y(u.*v, r-u)...
    -u.*w.*atan1(u.*v, w.*r)...
    -0.5*v.*r;
    allag_correction = -1;
    component_x = allag_correction*component_x;
    component_y = allag_correction*component_y;
    component_z = allag_correction*component_z;
    if 0
        ⟨ Test against Janssen results 22 ⟩
    end
    ⟨ Finish up 29 ⟩

```

See also section 23.

This code is used in section 19.

22. This is the same calculation with Janssen's equations instead. By default this code never runs, but if you like it can be enabled to prove that the equations are consistent.

```

⟨ Test against Janssen results 22 ⟩ ≡
    S = u;
    T = v;
    U = w;
    R = r;

    component_x_ii = ...
    (0.5*atan1(U, S) + 0.5*atan1(T.*U, S.*R)).*S.^2...
    +T.*S - 3/2*U.*S - multiply_x_log_y(S.*T, U + R) - T.^2.*atan1(S,
        T)...
    +U.*(U.*( ...
        0.5*atan1(S, U) + 0.5*atan1(S.*T, U.*R)...
        )...
        -multiply_x_log_y(T, S + R) + multiply_x_log_y(S, R - T)...
        )...
    +0.5*T.^2.*atan1(S.*U, T.*R)...
    ;

    component_y_ii = ...
    0.5*U.*(R - 2*S) + ...
    multiply_x_log_y(0.5*(T.^2 - S.^2), U + R) + ...
    S.*T.*(atan1(U, T) + atan1(S.*U, T.*R)) + ...
    multiply_x_log_y(S.*U, R - S)...
    ;

    component_z_ii = ...
    0.5*T.*(R - 2*S) + ...
    multiply_x_log_y(0.5*(U.^2 - S.^2), T + R) + ...
    S.*U.*(atan1(T, U) + atan1(S.*T, U.*R)) + ...
    multiply_x_log_y(S.*T, R - S)...
    ;

    if 1
        xx = index_sum .* component_x;
        xx_ii = index_sum .* component_x_ii;
        assert(abs(sum(xx(:)) - sum(xx_ii(:))) < 1 * 10-8)
    end

    if 1
        yy = index_sum .* component_y;
        yy_ii = index_sum .* component_y_ii;
        assert(abs(sum(yy(:)) - sum(yy_ii(:))) < 1 * 10-8)
    end

    if 1
        zz = index_sum .* component_z;
        zz_ii = index_sum .* component_z_ii;

```

```

    assert(abs(sum(zz(:)) - sum(zz_ii(:))) < 1 · 10-8)
end
if 1
    component_x = component_x_ii;
    component_y = component_y_ii;
    component_z = component_z_ii;
end

```

This code is used in section 21.

23. Don't need to swap $J1$ because it should only contain z components anyway. (This is assumption isn't tested because it it's wrong we're in more trouble anyway; this should all be taken care of earlier when the magnetisation components were separated out.)

⟨ Orthogonal magnets force calculation 21 ⟩ \equiv

```

function calc_out = forces_calc_z_x(size1, size2, offset, J1, J2)
    forces_xyz = forces_calc_z_y(...
        rotate_x_to_y(size1), rotate_x_to_y(size2), rotate_x_to_y(offset), ...
        J1, rotate_x_to_y(J2));
    calc_out = rotate_y_to_x(forces_xyz);
end

```

24. Stiffness calculations are derived³ from the forces.

⟨ Parallel magnets stiffness calculation 24 ⟩ \equiv

```

function calc_out = stiffnesses_calc_z_z(size1, size2, offset, J1, J2)
    J1 = J1(3);
    J2 = J2(3);
    ⟨ Initialise subfunction variables 27 ⟩
    component_x = ...
        -r...
        -(u.^2.*v)./(u.^2+w.^2)...
        -v.*log(r-v);
    component_y = ...
        -r...
        -(v.^2.*u)./(v.^2+w.^2)...
        -u.*log(r-u);
    component_z = -component_x - component_y;
    ⟨ Finish up 29 ⟩

```

This code is used in section 19.

³Literally.

25. Orthogonal magnets stiffnesses derived from Yonnet and Allag [2]. First the z - y magnetisation.

⟨ Orthogonal magnets stiffness calculation 25 ⟩ \equiv

```
function calc_out = stiffnesses_calc_z_y(size1, size2, offset, J1, J2)
    J1 = J1(3);
    J2 = J2(2);
    ⟨ Initialise subfunction variables 27 ⟩
    component_x = -((u.^2.*v)./(u.^2+v.^2))-(u.^2.*w)./(u.^2+w.^2)...
    +u.*atan1(v.*w,r.*u)-multiply_x_log_y(w,r+v)+...
    -multiply_x_log_y(v,r+w);
    component_y = v/2-(u.^2.*v)./(u.^2+v.^2)+(u.*v.*w)./(v.^2+w.^2)...
    +u.*atan1(u.*w,r.*v)+multiply_x_log_y(v,r+w);
    component_z = -component_x - component_y;
    allag_correction = -1;
    component_x = allag_correction*component_x;
    component_y = allag_correction*component_y;
    component_z = allag_correction*component_z;
    ⟨ Finish up 29 ⟩
```

See also section 26.

This code is used in section 19.

26. Now the z - x magnetisation, which is z - y rotated.

⟨ Orthogonal magnets stiffness calculation 25 ⟩ $+\equiv$

```
function calc_out = stiffnesses_calc_z_x(size1, size2, offset, J1, J2)
    stiffnesses_xyz = stiffnesses_calc_z_y(...
        rotate_x_to_y(size1), rotate_x_to_y(size2), rotate_x_to_y(offset), ...
        J1, rotate_x_to_y(J2));
    calc_out = rotate_y_to_x(stiffnesses_xyz);
end
```

27. Some shared setup code. First **return** early if either of the magnetisations are zero — that’s the trivial solution. Assume that the magnetisation has already been rounded down to zero if necessary; i.e., that we don’t need to check for $J1$ or $J2$ are less than $1 \cdot 10^{-12}$ or whatever.

```

< Initialise subfunction variables 27 > ≡
    if ( J1 ≡ 0 OR J2 ≡ 0 )
        debug_disp('Zero magnetisation.')
        calc_out = [0; 0; 0];
        return;
    end
    u = offset(1) + size2(1)*(-1).^index_j - size1(1)*(-1).^index_i;
    v = offset(2) + size2(2)*(-1).^index_l - size1(2)*(-1).^index_k;
    w = offset(3) + size2(3)*(-1).^index_q - size1(3)*(-1).^index_p;
    r = sqrt(u.^2 + v.^2 + w.^2);

```

This code is used in sections 20, 21, 24, and 25.

28. Here are some variables used above that only need to be computed once. The idea here is to vectorise instead of using **for** loops because it allows more convenient manipulation of the data later on.

```

< Initialise main variables 6 > +≡
    magconst = 1/(4*pi*(4*pi*1*10^-7));
    [index_i, index_j, index_k, index_l, index_p, index_q] = ndgrid([0 1]);
    [index2_j, index2_l, index2_q] = ndgrid([0 1]);
    index_sum = (-1).^(index_i+index_j+index_k+index_l+index_p+index_q);

```

29. And some shared finishing code.

```

< Finish up 29 > ≡
    component_x = index_sum .* component_x;
    component_y = index_sum .* component_y;
    component_z = index_sum .* component_z;
    calc_out = J1*J2*magconst.*...
    [sum(component_x(:));
     sum(component_y(:));
     sum(component_z(:))];
    debug_disp(calc_out')
end

```

This code is used in sections 20, 21, 24, and 25.

30. Setup code.

31. When the forces are rotated we use these rotation matrices to avoid having to think too hard. Use degrees in order to compute $\sin(\pi/2)$ exactly!

```
<Precompute rotation matrices 31> ≡  
    swap_x_y = @(vec) vec([2 1 3]);  
    swap_x_z = @(vec) vec([3 2 1]);  
    swap_y_z = @(vec) vec([1 3 2]);  
  
    Rx = @(θ) [1 0 0; 0 cosd(θ) - sind(θ); 0 sind(θ) cosd(θ)];  
    Ry = @(θ) [cosd(θ) 0 sind(θ); 0 1 0; -sind(θ) 0 cosd(θ)];  
    Rz = @(θ) [cosd(θ) - sind(θ) 0; sind(θ) cosd(θ) 0; 0 0 1];  
  
    Rx_180 = Rx(180);  
    Rx_090 = Rx(90);  
    Rx_270 = Rx(-90);  
    Ry_180 = Ry(180);  
    Ry_090 = Ry(90);  
    Ry_270 = Ry(-90);  
    Rz_180 = Rz(180);  
    Rz_090 = Rz(90);  
    Rz_270 = Rz(-90);  
  
    rotate_z_to_x = @(vec) Ry_090*vec;  
    rotate_x_to_z = @(vec) Ry_270*vec;  
  
    rotate_z_to_y = @(vec) Rx_090*vec;  
    rotate_y_to_z = @(vec) Rx_270*vec;  
  
    rotate_x_to_y = @(vec) Rz_090*vec;  
    rotate_y_to_x = @(vec) Rz_270*vec;
```

This code is used in section 4.

32. The equations contain two singularities. Specifically, the equations contain terms of the form $x \log(y)$, which becomes NaN when both x and y are zero since $\log(0)$ is negative infinity.

This function computes $x \log(y)$, special-casing the singularity to output zero, instead. (This is indeed the value of the limit.)

```
<Helper functions 32> ≡  
function out = multiply_x_log_y(x, y)  
    out = x .* log(y);  
    out(isnan(out)) = 0;  
end
```

See also sections 33 and 34.

This code is used in section 19.

33. Also, we're using `atan` instead of `atan2` (otherwise the wrong results are calculated — I guess I don't totally understand that), which becomes a problem when trying to compute `atan(0/0)` since `0/0` is `NaN`.

This function computes `atan` but takes two arguments.

⟨Helper functions 32⟩ +≡

```
function out = atan1(x, y)
    out = zeros(size(x));
    ind = x ≠ 0 ∧ y ≠ 0;
    out(ind) = atan(x(ind) ./ y(ind));
end
```

34. This function is for easy debugging; in normal use it gobbles its argument but will print diagnostics when required.

⟨Helper functions 32⟩ +≡

```
function debug_disp(str)
    %disp(str)
end
```

35. When users type `help magnetforces` this is what they see.

⟨Matlab help text (forces) 35⟩ ≡

```
%% MAGNETFORCES Calculate forces between two cuboid magnets
%
% Finish this off later.
%
```

This code is used in section 4.

36. Test files. The chunks that follow are designed to be saved into individual files and executed automatically to check for (a) correctness and (b) regression problems as the code evolves.

How do I know if the code produces the correct forces? Well, for many cases I can compare with published values in the literature. Beyond that, I'll be setting up some tests that I can logically infer should produce the same results (such as mirror-image displacements) and test that.

There are many Matlab unit test frameworks but I'll be using a fairly low-tech method. In time this test suite should be (somehow) useable for all implementations of `magnetocode`, not just Matlab. But I haven't thought about doing anything like that, yet.

37. Because I'm lazy, just run the tests manually for now. This script must be run twice if it updates itself.

```
<testall.m 37> ≡  
  clc;  
  unix('~/bin/mtangle_magnetforces');  
  magforce.test001a  
  magforce.test001b  
  magforce.test001c  
  magforce.test001d
```

38. Force testing.

39. This test checks that square magnets produce the same forces in the each direction when displaced in positive and negative x , y , and z directions, respectively. In other words, this tests the function *forces_calc_z_y* directly. Both positive and negative magnetisations are used.

```

<magforce_test001a.m 39> ≡
    disp('=====')
    fprintf('TEST_001a:␣')
    magnet_fixed.dim = [0.04 0.04 0.04];
    magnet_float.dim = magnet_fixed.dim;
    magnet_fixed.magn = 1.3;
    magnet_float.magn = 1.3;
    offset = 0.1;
    <Test z-z magnetisations 40>
    <Assert magnetisations tests 48>
    <Test x-x magnetisations 41>
    <Assert magnetisations tests 48>
    <Test y-y magnetisations 42>
    <Assert magnetisations tests 48>
    fprintf('passed\n')
    disp('=====')

```

40. Testing vertical forces.

```

<Test z-z magnetisations 40> ≡
    f = [];
    for ii = [1, -1]
        magnet_fixed.magdir = [0 ii*90];           % ±z
        for jj = [1, -1]
            magnet_float.magdir = [0 jj*90];
            for kk = [1, -1]
                displ = kk*[0 0 offset];
                f(:, end + 1) = magnetforces(magnet_fixed, magnet_float, displ);
            end
        end
    end
    dirforces = chop(f(3, :), 8);
    otherforces = f([1 2], :);

```

This code is used in section 39.

41. Testing horizontal x forces.

⟨ Test x - x magnetisations 41 ⟩ \equiv

```
f = [];
for ii = [1, -1]
    magnet_fixed.magdir = [90 + ii*90 0];      %  $\pm x$ 
    for jj = [1, -1]
        magnet_float.magdir = [90 + jj*90 0];
        for kk = [1, -1]
            displ = kk*[offset 0 0];
            f(:, end + 1) = magnetforces(magnet_fixed, magnet_float, displ);
        end
    end
end
dirforces = chop(f(1, :), 8);
otherforces = f([2 3], :);
```

This code is used in section 39.

42. Testing horizontal y forces.

⟨ Test y - y magnetisations 42 ⟩ \equiv

```
f = [];
for ii = [1, -1]
    magnet_fixed.magdir = [ii*90 0];          %  $\pm y$ 
    for jj = [1, -1]
        magnet_float.magdir = [jj*90 0];
        for kk = [1, -1]
            displ = kk*[0 offset 0];
            f(:, end + 1) = magnetforces(magnet_fixed, magnet_float, displ);
        end
    end
end
dirforces = chop(f(2, :), 8);
otherforces = f([1 3], :);
```

This code is used in section 39.

43. This test does the same thing but for orthogonally magnetised magnets.

```

<magforce_test001b.m 43> ≡
    disp('=====')
    fprintf('TEST_001b:␣')
    magnet_fixed.dim = [0.04 0.04 0.04];
    magnet_float.dim = magnet_fixed.dim;
    magnet_fixed.magn = 1.3;
    magnet_float.magn = 1.3;
    <Test ZYZ 44>
    <Assert magnetisations tests 48>
    <Test ZXZ 45>
    <Assert magnetisations tests 48>
    <Test ZXX 47>
    <Assert magnetisations tests 48>
    <Test ZYY 46>
    <Assert magnetisations tests 48>
    fprintf('passed\n')
    disp('=====')

```

44. z - y magnetisations, z displacement.

```

<Test ZYZ 44> ≡
    fzyz = [];
    for ii = [1, -1]
        for jj = [1, -1]
            for kk = [1, -1]
                magnet_fixed.magdir = ii*[0 90];           % ±z
                magnet_float.magdir = jj*[90 0];           % ±y
                displ = kk*[0 0 0.1];                       % ±z
                fzyz(:, end + 1) = magnetforces(magnet_fixed, magnet_float, displ);
            end
        end
    end
    dirforces = chop(fzyz(2, :), 8);
    otherforces = fzyz([1 3], :);

```

This code is used in section 43.

45. z - x magnetisations, z displacement.

⟨ Test ZXZ 45 ⟩ ≡

```
fzxx = [];
for ii = [1, -1]
    for jj = [1, -1]
        for kk = [1, -1]
            magnet_fixed.magdir = ii*[0 90];           % ±z
            magnet_float.magdir = [90 + jj*90 0];       % ±x
            displ = kk*[0.1 0 0];                       % ±x
            fzxx(:, end + 1) = magnetforces(magnet_fixed, magnet_float, displ);
        end
    end
end

dirforces = chop(fzxx(3, :), 8);
otherforces = fzxx([1 2], :);
```

This code is used in section 43.

46. z - y magnetisations, y displacement.

⟨ Test ZYY 46 ⟩ ≡

```
fzyy = [];
for ii = [1, -1]
    for jj = [1, -1]
        for kk = [1, -1]
            magnet_fixed.magdir = ii*[0 90];           % ±z
            magnet_float.magdir = jj*[90 0];           % ±y
            displ = kk*[0 0.1 0];                       % ±y
            fzyy(:, end + 1) = magnetforces(magnet_fixed, magnet_float,
                                              displ);
        end
    end
end

dirforces = chop(fzyy(3, :), 8);
otherforces = fzyy([1 2], :);
```

This code is used in section 43.

47. z - x magnetisations, x displacement.

⟨ Test ZXX 47 ⟩ \equiv

```
fzxx = [];
for ii = [1, -1]
    for jj = [1, -1]
        for kk = [1, -1]
            magnet_fixed.magdir = ii*[0 90];           %  $\pm z$ 
            magnet_float.magdir = [90 + jj*90 0];       %  $\pm x$ 
            displ = kk*[0 0 0.1];                       %  $\pm z$ 
            fzxx(:, end+1) = magnetforces(magnet_fixed, magnet_float, displ);
        end
    end
end
dirforces = chop(fzxx(1, :), 8);
otherforces = fzxx([2 3], :);
```

This code is used in section 43.

48. The assertions, common between directions.

⟨ Assert magnetisations tests 48 ⟩ \equiv

```
assert(...
    all(abs(otherforces(:)) < 1 · 10-11), ...
    'Orthogonal_forces_should_be_zero' ...
)
assert(...
    all(abs(dirforces) == abs(dirforces(1))), ...
    'Force_magnitudes_should_be_equal' ...
)
assert(...
    all(dirforces(1:4) == -dirforces(5:8)), ...
    'Forces_should_be_opposite_with_reversed_fixed_magnet_magnetisation' ...
)
assert(...
    all(dirforces([1 3 5 7]) == -dirforces([2 4 6 8])), ...
    'Forces_should_be_opposite_with_reversed_float_magnet_magnetisation' ...
)
```

This code is used in sections 39 and 43.

49. Now try combinations of displacements.

```

<magforce_test001c.m 49> ≡
    disp('=====')
    fprintf('TEST_001c:␣')
    magnet_fixed.dim = [0.04 0.04 0.04];
    magnet_float.dim = magnet_fixed.dim;
    magnet_fixed.magn = 1.3;
    magnet_float.magn = 1.3;
    <Test combinations ZZ 50>
    <Assert combinations tests 52>
    <Test combinations ZY 51>
    <Assert combinations tests 52>
    fprintf('passed\n')
    disp('=====')

```

50. Tests.

```

<Test combinations ZZ 50> ≡
    f = [];
    for ii = [-1 1]
        for jj = [-1 1]
            for xx = 0.12*[-1, 1]
                for yy = 0.12*[-1, 1]
                    for zz = 0.12*[-1, 1]
                        magnet_fixed.magdir = [0 ii*90];      % z
                        magnet_float.magdir = [0 jj*90];      % z
                        displ = [xx yy zz];
                        f(:, end + 1) = magnetforces(magnet_fixed, magnet_float,
                            displ);
                    end
                end
            end
        end
    end
    f = chop(f, 8);
    uniquedir = f(3, :);
    otherdir = f([1 2], :);

```

This code is used in section 49.

51. Tests.

⟨ Test combinations ZY 51 ⟩ ≡

```
f = [];
for ii = [-1 1]
    for jj = [-1 1]
        for xx = 0.12*[-1, 1]
            for yy = 0.12*[-1, 1]
                for zz = 0.12*[-1, 1]
                    magnet_fixed.magdir = [0 ii*90];           % ±z
                    magnet_float.magdir = [jj*90 0];           % ±y
                    displ = [xx yy zz];
                    f(:, end + 1) = magnetforces(magnet_fixed, magnet_float,
                                                    displ);
                end
            end
        end
    end
end
f = chop(f, 8);
uniquedir = f(1, :);
otherdir = f([2 3], :);
```

This code is used in section 49.

52. Shared tests, again.

⟨ Assert combinations tests 52 ⟩ ≡

```
test1 = abs(diff(abs(f(1, :)))) < 1 · 10-10;
test2 = abs(diff(abs(f(2, :)))) < 1 · 10-10;
test3 = abs(diff(abs(f(3, :)))) < 1 · 10-10;
assert( all(test1) ∧ ∧ all(test2) ∧ ∧ all(test3), ...
        'All forces in a single direction should be equal' )
test = abs(diff(abs(otherdir))) < 1 · 10-11;
assert(all(test), 'Orthogonal forces should be equal')
test1 = f(:, 1:8) ≡ f(:, 25:32);
test2 = f(:, 9:16) ≡ f(:, 17:24);
assert( all(test1(:)) ∧ ∧ all(test2(:)), ...
        'Reverse magnetisation shouldn't make a difference' )
```

This code is used in section 49.

53. Now we want to try non-orthogonal magnetisation.

```

<magforce_test001d.m 53> ≡
    disp('=====')
    fprintf('TEST_001d: ')
    magnet_fixed.dim = [0.04 0.04 0.04];
    magnet_float.dim = magnet_fixed.dim;

    % Fixed parameters:
    magnet_fixed.magn = 1.3;
    magnet_float.magn = 1.3;
    magnet_fixed.magdir = [0 90];      % z
    displ = 0.12*[1 1 1];
    <Test XY superposition 54>
    <Assert superposition 57>
    <Test XZ superposition 55>
    <Assert superposition 57>
    <Test planar superposition 56>
    <Assert superposition 57>
    fprintf('passed\n')
    disp('=====')

```

54. Test with a magnetisation unit vector of $(1, 1, 0)$.

```

<Test XY superposition 54> ≡
    magnet_float.magdir = [45 0];      %  $\vec{e}_x + \vec{e}_y$ 
    f1 = magnetforces(magnet_fixed, magnet_float, displ);

    % Components:
    magnet_float.magdir = [0 0];      %  $\vec{e}_x$ 
    fc1 = magnetforces(magnet_fixed, magnet_float, displ);
    magnet_float.magdir = [90 0];     %  $\vec{e}_y$ 
    fc2 = magnetforces(magnet_fixed, magnet_float, displ);
    f2 = (fc1 + fc2)/sqrt(2);

```

This code is used in section 53.

55. Test with a magnetisation unit vector of $(1, 0, 1)$.

```

< Test XZ superposition 55 > ≡
    magnet_float.magdir = [0 45];           %  $\vec{e}_y + \vec{e}_z$ 
    f1 = magnetforces(magnet_fixed, magnet_float, displ);

    % Components:
    magnet_float.magdir = [0 0];           %  $\vec{e}_x$ 
    fc1 = magnetforces(magnet_fixed, magnet_float, displ);
    magnet_float.magdir = [0 90];          %  $\vec{e}_z$ 
    fc2 = magnetforces(magnet_fixed, magnet_float, displ);
    f2 = (fc1 + fc2)/sqrt(2);

```

This code is used in section 53.

56. Test with a magnetisation unit vector of $(1, 1, 1)$. This is about as much as I can be bothered testing for now. Things seem to be working.

```

< Test planar superposition 56 > ≡
    [t p r] = cart2sph(1/sqrt(3), 1/sqrt(3), 1/sqrt(3));
    magnet_float.magdir = [t p]*180/π;      %  $\vec{e}_y + \vec{e}_z + \vec{e}_z$ 
    f1 = magnetforces(magnet_fixed, magnet_float, displ);

    % Components:
    magnet_float.magdir = [0 0];           %  $\vec{e}_x$ 
    fc1 = magnetforces(magnet_fixed, magnet_float, displ);
    magnet_float.magdir = [90 0];          %  $\vec{e}_y$ 
    fc2 = magnetforces(magnet_fixed, magnet_float, displ);
    magnet_float.magdir = [0 90];          %  $\vec{e}_z$ 
    fc3 = magnetforces(magnet_fixed, magnet_float, displ);
    f2 = (fc1 + fc2 + fc3)/sqrt(3);

```

This code is used in section 53.

57. The assertion is the same each time.

```

< Assert superposition 57 > ≡
    assert(...
        isequal(chop(f1, 4), chop(f2, 4)), ...
        'Components should sum due to superposition' ...
    )

```

This code is used in section 53.

58. Forces between (multipole) magnet arrays.

59. This function uses `magnetforces.m` to compute the forces between two multipole magnet arrays. As before, we can calculate either force and/or stiffness in all three directions.

Inputs:	<i>fixed_array</i>	structure describing first magnet array
	<i>float_array</i>	structure describing the second magnet array
	<i>displ</i>	displacement between first magnet of each array
	[<i>what to calculate</i>]	'force' and/or 'stiffness'
Outputs:	<i>forces</i>	forces on the second array
	<i>stiffnesses</i>	stiffnesses on the second array
Array properties:	<i>mcount</i>	[<i>i j k</i>] magnets in each direction
	<i>msize</i>	size of each magnet
	<i>mgap</i>	gap between successive magnets
	<i>magn</i>	magnetisation magnitude
	<i>magdir_fn</i>	function to calculate the magnetisation direction

⟨multipoleforces.m 59⟩ ≡

function [varargout] = multipoleforces(*fixed_array*, *float_array*, *displ*, varargin)

⟨ Matlab help text (multipole) 67 ⟩

⟨ Parse calculation args 9 ⟩

⟨ Create arrays of magnets 61 ⟩

⟨ Calculate array forces 60 ⟩

⟨ Combine results and exit 10 ⟩

60. To calculate the forces between the magnet arrays, let's assume that we have two large arrays enumerating the positions and magnetisations of each individual magnet in each magnet array.

```

⟨ Calculate array forces 60 ⟩ ≡
    if calc_force_bool
        array_forces = repmat(NaN, [M N 3]);
    end
    if calc_stiffness_bool
        array_stiffnesses = repmat(NaN, [M N 3]);
    end
    for mm = 1 : M
        fixed_magnet = struct(...
            'dim', fixed_magnet_dim(mm, :), ...
            'magn', fixed_magnet_magn(mm), ...
            'magdir', fixed_magnet_magdir(mm, :) ...
        );
        for nn = 1 : N
            float_magnet = struct(...
                'dim', float_magnet_dim(nn, :), ...
                'magn', float_magnet_magn(nn), ...
                'magdir', float_magnet_magdir(nn, :) ...
            );
            displ = fixed_magnet_loc(mm, :) + float_magnet_loc(nn, :);
            if calc_force_bool
                array_forces(mm, nn, :) = ...
                    magnetforces(fixed_magnet, float_magnet, displ, 'force');
            end
            if calc_stiffness_bool
                array_stiffnesses(mm, nn, :) = ...
                    magnetforces(fixed_magnet, float_magnet, displ, 'stiffness');
            end
        end
    end
    if calc_force_bool
        forces_out = squeeze(sum(sum(array_forces, 1), 2));
    end
    if calc_stiffness_bool
        stiffnesses_out = squeeze(sum(sum(array_stiffnesses, 1), 2));
    end
end

```

This code is used in section 59.

61. We separate the force calculation from transforming the inputs into an intermediate form used for that purpose. This will hopefully allow us a little more flexibility later on.

```

⟨ Create arrays of magnets 61 ⟩ ≡
    M = prod(fixed_array.mcount);
    N = prod(float_array.mcount);
    fixed_magnet_loc = repmat(NaN, [M 3]);
    float_magnet_loc = repmat(NaN, [N 3]);
    fixed_magnet_magdir = repmat(NaN, [M 2]);
    float_magnet_magdir = repmat(NaN, [N 2]);

```

See also sections 62, 63, 64, 65, and 66.

This code is used in section 59.

62. Size of each magnet.

```

⟨ Create arrays of magnets 61 ⟩ +≡
    if length(fixed_array.msize) == 3
        fixed_magnet_dim_array = ...
            repmat(reshape(fixed_array.msize, [1 1 1 3]), fixed_array.mcount);
        fixed_magnet_dim = reshape(fixed_magnet_dim_array, [M 3]);
    else
        error('Not yet implemented.')
    end
    if length(float_array.msize) == 3
        float_magnet_dim_array = ...
            repmat(reshape(float_array.msize, [1 1 1 3]), float_array.mcount);
        float_magnet_dim = reshape(float_magnet_dim_array, [N 3]);
    else
        error('Not yet implemented.')
    end
end

```

63. Magnetisation strength of each magnet.

```

⟨ Create arrays of magnets 61 ⟩ +≡
    if length(fixed_array.magn) == 1
        fixed_magnet_magn = repmat(fixed_array.magn, [M 1]);
    else
        error('Not yet implemented.')
    end
    if length(float_array.magn) == 1
        float_magnet_magn = repmat(float_array.magn, [N 1]);
    else
        error('Not yet implemented.')
    end
end

```

64. Gaps, if any, between each magnet.

```

⟨ Create arrays of magnets 61 ⟩ +=
    if length(fixed_array.mgap) == 3
        fixed_gaps = fixed_array.mgap;
    elseif length(fixed_array.mgap) == 1
        fixed_gaps = repmat(fixed_array.mgap, [3 1]);
    else
        error('Not yet implemented.')
    end

    if length(float_array.mgap) == 3
        float_gaps = float_array.mgap;
    elseif length(float_array.mgap) == 1
        float_gaps = repmat(float_array.mgap, [3 1]);
    else
        error('Not yet implemented.')
    end
end

```

65. Location of each magnet.

```

⟨ Create arrays of magnets 61 ⟩ +=
    ii = 0;
    for xx = 1 : fixed_array.mcount (1)
        for yy = 1 : fixed_array.mcount (2)
            for zz = 1 : fixed_array.mcount (3)
                ii = ii + 1;
                fixed_magnet_loc(ii, :) = ...
                    [xx - 1; yy - 1; zz - 1] .* (squeeze(fixed_magnet_dim_array(xx,
                        yy, zz, :)) + fixed_gaps);
            end
        end
    end

    ii = 0;
    for xx = 1 : float_array.mcount (1)
        for yy = 1 : float_array.mcount (2)
            for zz = 1 : float_array.mcount (3)
                ii = ii + 1;
                float_magnet_loc(ii, :) = ...
                    [xx - 1; yy - 1; zz - 1] .* (squeeze(float_magnet_dim_array(xx,
                        yy, zz, :)) + float_gaps);
            end
        end
    end
end

```


66. Magnetisation direction of each magnet.

⟨ Create arrays of magnets **61** ⟩ \equiv

```

ii = 0;
for xx = 1 : fixed_array.mcount (1)
    for yy = 1 : fixed_array.mcount (2)
        for zz = 1 : fixed_array.mcount (3)
            ii = ii + 1;
            fixed_magnet_magdir(ii, :) = fixed_array.magdir_fn(xx, yy, zz);
        end
    end
end

ii = 0;
for xx = 1 : float_array.mcount (1)
    for yy = 1 : float_array.mcount (2)
        for zz = 1 : float_array.mcount (3)
            ii = ii + 1;
            float_magnet_magdir(ii, :) = float_array.magdir_fn(xx, yy, zz);
        end
    end
end
end

```

67. When users type `help multipoleforces` this is what they see.

⟨ Matlab help text (multipole) **67** ⟩ \equiv

```

%% MULTIPOLEFORCES Calculate forces between two multipole arrays of magnets
%
% Finish this off later.
%

```

This code is used in section **59**.

68. Test files for multipole arrays.

```

<multiforce_test001a.m 68> ≡
  W = 4;
  fixed_array = ...
  struct(...
    'mcount', [5 1 1], ...
    'msize', [0.01 0.01 0.01], ...
    'mgap', 0, ...
    'magn', 1, ...
    'magdir_fn', @(ii, jj, kk) [0 + 90 + (ii - 1)/W*360] ...
  );
  float_array = ...
  struct(...
    'mcount', [5 1 1], ...
    'msize', [0.01 0.01 0.01], ...
    'mgap', 0, ...
    'magn', 1, ...
    'magdir_fn', @(ii, jj, kk) [0 - 90 - (ii - 1)/W*360] ...
  );
  displ = [0 0 0.04];
  multipoleforces(fixed_array, float_array, displ)

```

69. These are MATLABWEB declarations to improve the formatting of this document. Ignore unless you're editing `magnetforces.web`.

```

define end ≡ end
format END TeX

```

Index of magnetforces

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