

The following is the code for evaluating a definite integral of a given function by a Numerical Method called Simpson's 1/3rd Rule.

DOWNLOAD:[simpson](#)

```
funcprot(0);
function ans=simpson(a,b,n,g)
    h=(b-a)/n;
    sum=0;
    for i=1:n-1
        x=a+i*h;
        if modulo(i,2)==0
            sum=sum+2*g(x);
        else
            sum=sum+4*g(x);
        end
    end
    ans=(h/3)*(g(a)+g(b)+sum);
endfunction
```

You can either copy the code above and save it as a .sci file or download the file [simpson](#) . Once you run the code, the function '*simpson(a,b,n,f)*' can be called by other programs or even in the console.

Function syntax:

simpson(a,b,n,f)

where,

a=initial limit(real no.)

b=final limit(real no.)

n=no. of sub-intervals(the higher the value of 'n' the better is the result.

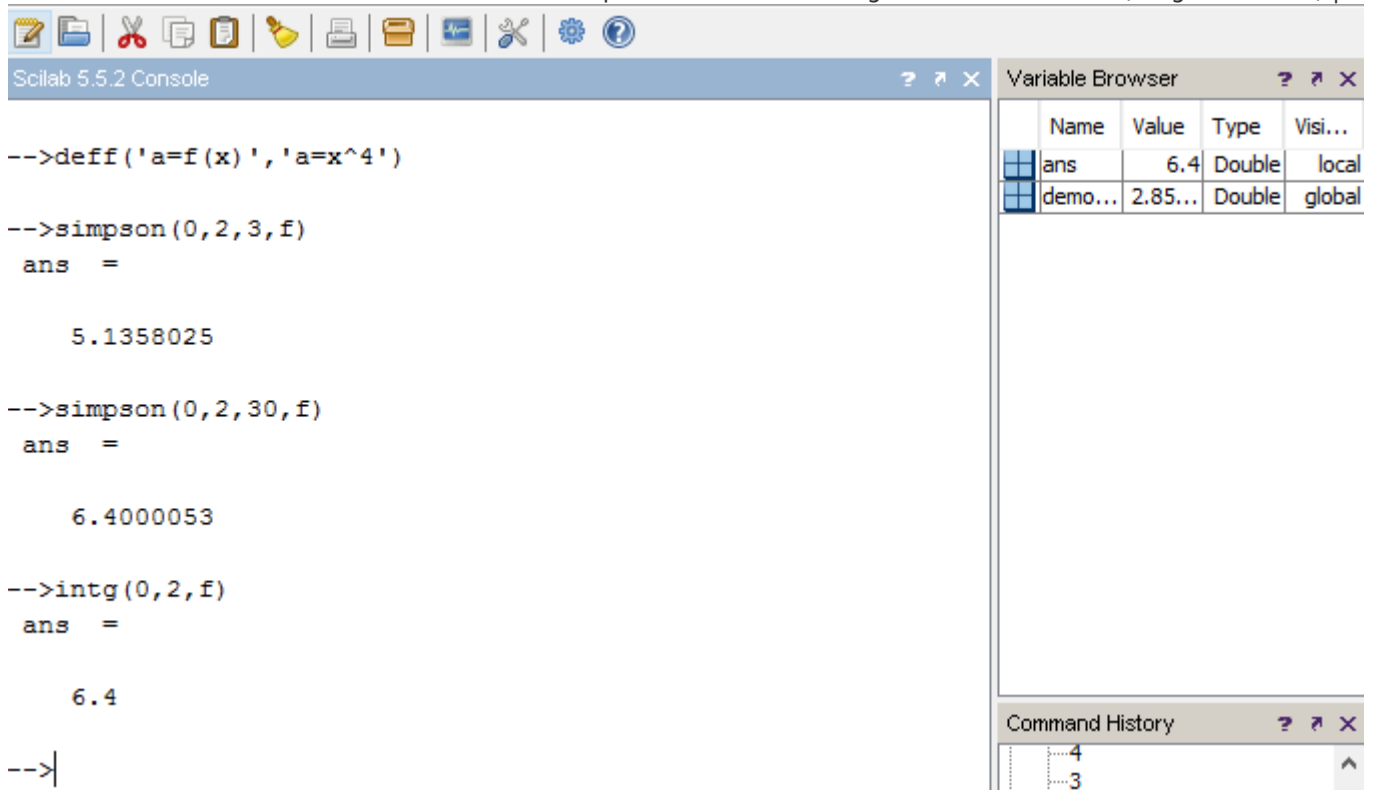
NOTE: n should be an even no.

Example:

The following code snippet evaluates the integral of x^4 from 0 to 2.

```
deff('a=f(x)', 'a=x^4');
integral=simpson(0,2,30,f);
```

Here is a comparison of the result with the inbuilt function 'intg'.



```

-->deff('a=f(x)', 'a=x^4')

-->simpson(0,2,3,f)
ans =

    5.1358025

-->simpson(0,2,30,f)
ans =

    6.4000053

-->intg(0,2,f)
ans =

    6.4

-->|

```

Name	Value	Type	Visi...
ans	6.4	Double	local
demo...	2.85...	Double	global

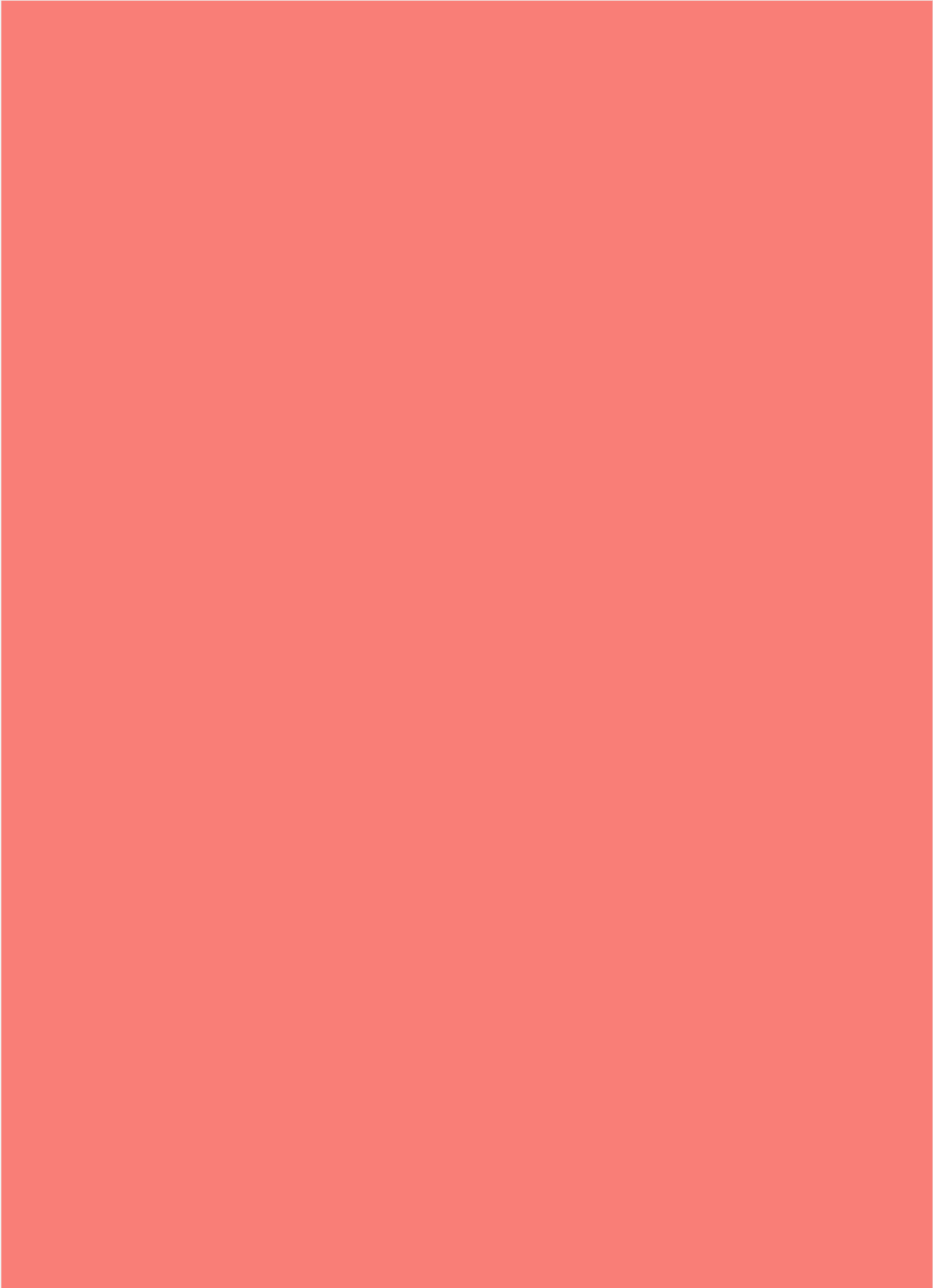
Command History	
4	↑
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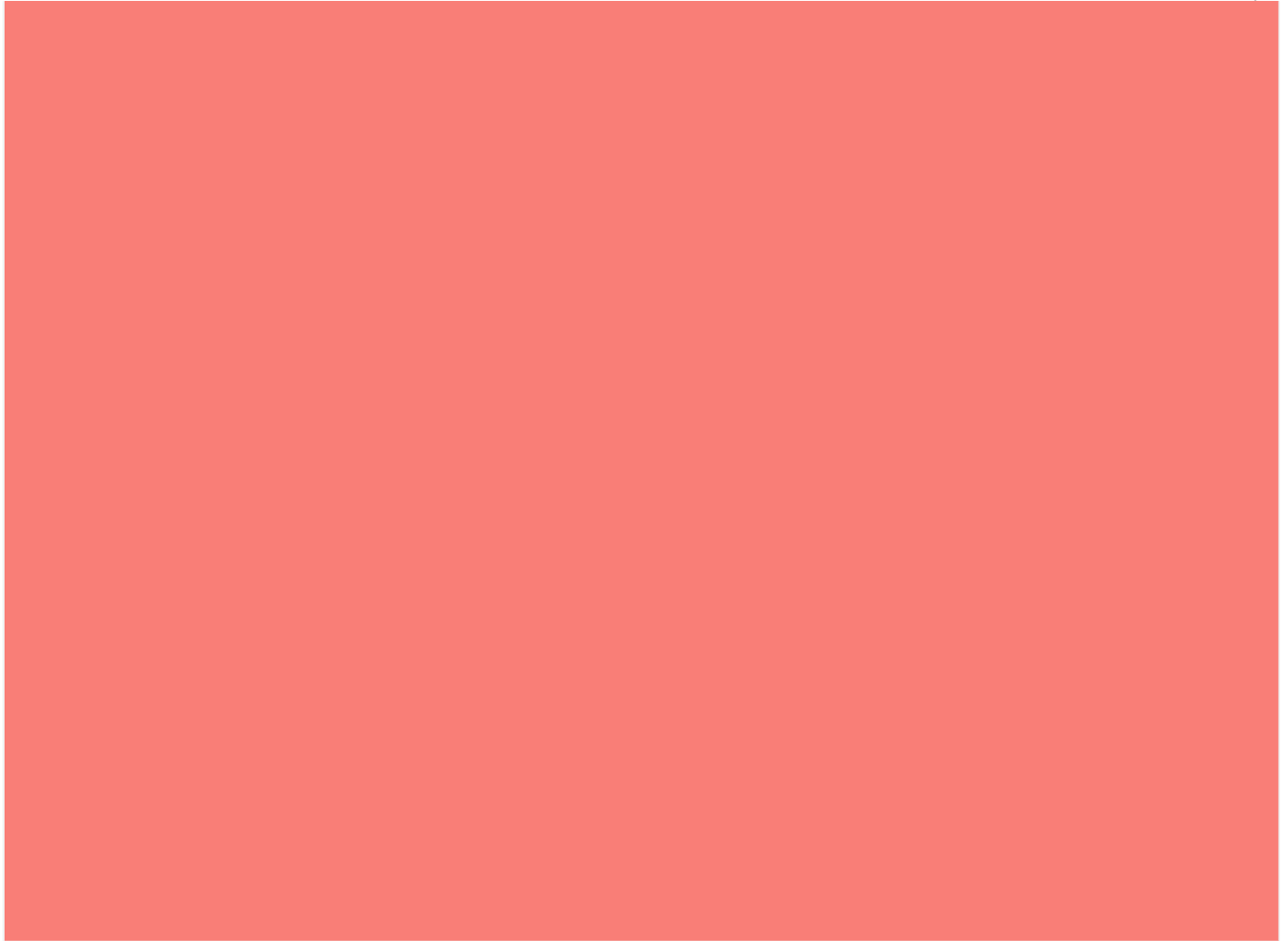
I'm a physicist specializing in computational material science with a PhD in Physics from Friedrich-Schiller University Jena, Germany. I write efficient codes for simulating light-matter interactions at atomic scales. I like to develop Physics, DFT, and Machine Learning related apps and software from time to time. Can code in most of the popular languages. I like to share my knowledge in Physics and applications using this Blog and a YouTube channel.

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