Nipype: A Flexible, Lightweight and Extensible Neuroimaging Data Processing Framework in Python

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<tr>
<td>Publisher</td>
<td>Frontiers Research Foundation</td>
</tr>
<tr>
<td>Version</td>
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</tr>
<tr>
<td>Accessed</td>
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Nipype: a flexible, lightweight and extensible neuroimaging data processing framework in Python

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INTRODUCTION
Over the past 20 years, advances in non-invasive in vivo neuroimaging have resulted in an explosion of studies investigating human cognition in health and disease. Current imaging studies acquire multi-modal image data (e.g., structural, diffusion, functional) and combine these with non-imaging behavioral data, patient and/or treatment history, and demographic and genetic information. Several sophisticated software packages (e.g., AFNI, BrainVoyager, FSL, FreeSurfer, Nipy, R, SPM) are used to process and analyze such extensive data. In a typical analysis, algorithms from these packages, each with its own set of parameters, process the raw data. However, data collected for a single study can be diverse (highly multi-dimensional) and large, and algorithms suited for one dataset may not be optimal for another. This complicates analysis methods and makes data exploration and inference challenging, and comparative analysis of new algorithms difficult.

CURRENT PROBLEMS
Here we outline issues that hinder replicable, efficient, and optimal use of neuroimaging analysis approaches.

No uniform access to neuroimaging analysis software and usage information
For current multi-modal datasets, researchers typically resort to using different software packages for different components of the analysis. However, these different software packages are accessed, and interfaced with, in different ways, such as: shell scripting (FSL, AFNI, Camino), MATLAB (SPM), and Python (Nipy). This has resulted in a heterogeneous set of software with no uniform way to use these tools or execute them. With the primary focus on algorithmic improvement, academic software development often lacks a rigorous software engineering framework that involves extensive testing and documentation and ensures compatibility with other
tools. This often necessitates extensive interactions with the authors of the software to understand their parameters, their quirks, and their usage.

No framework for comparative algorithm development and dissemination

Except for some large software development efforts (e.g., SPM, FSL, FreeSurfer), most algorithm development happens in-house and stays within the walls of a lab, without extensive exposure or testing. Furthermore, testing comparative efficacy of algorithms often requires significant effort (Klein et al., 2010). In general, developers create software for a single package (e.g., VBM8 for SPM), create a standalone cross-platform tool (e.g., Mricron), or simply do not distribute the software or code (e.g., normalization software used for registering architectonic atlases to MNI single subject template – Hömke, 2006).

Personnel turnover in laboratories often limits methodological continuity and training new personnel takes time

Most cognitive neuroscience laboratories aim to understand some aspect of cognition. Although a majority of such laboratories gather and analyze neuroimaging data, very few of them have the personnel with the technical expertise to understand methodological development and modify laboratory procedures to adopt new tools. Lab personnel with no prior imaging experience often learn by following online tutorials, taking organized courses or, as is most often the case, by learning from existing members of the lab. While this provides some amount of continuity, understanding different aspects of neuroimaging has a steep learning curve, and steeper when one takes into account the time and resources needed to learn the different package interfaces and algorithms.

Neuroimaging software packages do not address computational efficiency

The primary focus of neuroimaging analysis algorithms is to solve problems (e.g., registration, statistical estimation, tractography). While some developers focus on algorithmic or numerical efficiency, most developers do not focus on efficiency in the context of running multiple algorithms on multiple subjects, a common scenario in neuroimaging analysis. Creating an analysis workflow for a particular study is an iterative process dependent on the quality of the data and participant population (e.g., neurotypical, presurgical, etc.). Researchers usually experiment with different methods and their parameters to create a workflow suitable for their application, but no suitable framework currently exists to make this process efficient. Furthermore, very few of the available neuroimaging tools take advantage of the growing number of parallel hardware configurations (multi-core, clusters, clouds, and supercomputers).

Method sections of journal articles are often inadequate for reproducing results

Several journals (e.g., PNAS, Science, PLoS) require mandatory submission of data and scripts necessary to reproduce results of a study. However, most current method sections do not have sufficient details to enable a researcher knowledgeable in the domain to reproduce the analysis process. Furthermore, as discussed above, typical neuroimaging analyses integrate several tools and current analysis software do not make it easy to reproduce all the analysis steps in the proper order. This leaves a significant burden on the user to satisfy these journal requirements as well as ensure that analysis details are preserved with the intent to reproduce.

CURRENT SOLUTIONS

There have been several attempts to address these issues by creating pipeline systems (for comparison see Table 1). Taverna (Oinn et al., 2006), VisTrails (Callahan et al., 2006) are general pipelining systems with excellent support for web-services, but they do not address problems specific to neuroimaging. BrainVisa (Cointepas et al., 2001), MIPAV (McAuliffe et al., 2001), SPM include their own batch processing tools, but do not allow mixing components from other packages. Fiswidents (Fissell et al., 2003), a promising initial approach, appears to have not been developed and does not support state of the art methods. A much more extensive and feature rich solution is the LONI Pipeline (Rex et al., 2003; Dinov et al., 2009, 2010). It provides an easy to use graphical interface for choosing processing steps or nodes from a predefined library and defining their dependencies and parameters. Thanks to an advanced client–server architecture, it also has extensive support for parallel execution on an appropriately configured cluster (including data transfer, pausing execution, and combining local and remote software). Additionally, the LONI Pipeline saves information about executed steps (such as software origin, version, and architecture) providing provenance information (Mackenzie-Graham et al., 2008).

However, the LONI Pipeline does not come without limitations. Processing nodes are defined using eXtensible Markup Language (XML). This “one size fits all” method makes it easy to add new nodes as long as they are well-behaved command lines. However, many software packages do not meet this criterion. For example, SPM, written in MATLAB, does not provide a command line interface. Furthermore, for several command line programs, arguments are not easy to describe in the LONI XML schema (e.g., ANTS – Avants and Gee, 2004). Although it provides a helpful graphical interface, the LONI Pipeline environment does not provide an easy option to script a workflow or for rapidly exploring parametric variations within a workflow (e.g., VisTrails). Finally, due to restrictive licensing, it is not straightforward to modify and redistribute the modifications.

To address issues with existing workflow systems and the ones described earlier, we present Nipype (Neuroimaging in Python: Pipelines and Interfaces), an open-source, community-developed, Python-based software package that easily interfaces with existing software for efficient analysis of neuroimaging data and rapid comparative development of algorithms. Nipype uses a flexible, efficient, and general purpose programming language – Python – as its foundation. Processing modules and their inputs and outputs are described in an object-oriented manner providing the flexibility to interface with any type of software (not just well-behaved command lines). The workflow execution engine has a plug-in architecture and supports both local execution on multi-core machines and remote execution on clusters. Nipype is distributed with a Berkeley Software Distribution (BSD) license allowing anyone to make changes and redistribute it. Development is done openly with collaborators from many different labs, allowing adaptation to the varied needs of the neuroimaging community.
Table 1 | Feature comparison of selected pipeline frameworks.

<table>
<thead>
<tr>
<th></th>
<th>Local multi-processing</th>
<th>Grid engine</th>
<th>Scripting support</th>
<th>XNAT</th>
<th>Web-services</th>
<th>Platforms</th>
<th>Graphical user interface</th>
<th>Designed for neuroimaging</th>
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<td>PBS</td>
<td>Java, R</td>
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<td>DRMAA</td>
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<td>Mac, Unix, Windows</td>
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<td></td>
</tr>
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BrainVisa, MIPAV, and SPM were not included due to their inability to combine software from different packages.

1Without additional dependencies.

2Support for executing processing steps defined as web-services.

IMPLEMENTATION DETAILS

Nipype consists of three components (see Figure 1): (1) interfaces to external tools that provide a unified way for setting inputs, executing, and retrieving outputs; (2) a workflow engine that allows creating analysis pipelines by connecting inputs and outputs of interfaces as a directed acyclic graph (DAG); and (3) plug-ins that execute workflows either locally or in a distributed processing environment (e.g., Torque, SGE/OGE). In the following sections, we describe key architectural components and features of this software.

INTERFACES

Interfaces form the core of Nipype. The goal of Interfaces is to provide a uniform mechanism for accessing analysis tools from neuroimaging software packages (e.g., FreeSurfer, FSL, SPM). Interfaces can be used directly as a Python object, incorporated into custom Python scripts or used interactively in a Python console. For example, there is a Realign Interface that exposes the SPM re-alignment routine, while the MCFLIRT Interface exposes the FSL realignment routine. In addition, one can also implement an algorithm in Python within Nipype and expose it as an Interface. Interfaces are flexible and can accommodate the heterogeneous software that needs to be supported, while providing unified and uniform access to these tools for the user. Since, there is no need for the underlying software to be changed (re-compiled or adjusted to conform to a certain standard), developers can continue to create software using the computer language of their choice.

An Interface definition consists of: (a) input parameters, their types (e.g., file, floating point value, list of integers, etc.) and dependencies (e.g., does input “a” require input “b”); (b) outputs and their types, (c) how to execute the underlying software (e.g., run a MATLAB script, or call a command line program); and (d) a mapping which defines the outputs that are produced given a particular set of inputs. Using an object-oriented approach, we minimize redundancy in interface definition by creating a hierarchy of base Interface classes (see Figure 2) to encapsulate common functionality (e.g., Interfaces that call command line programs are derived from the CommandLine class, which provides methods to translate Interface inputs into command line parameters and for calling the command). Source code of an example Interface is shown in Listing 1.

We use Enthought Traits to create a formal definition for Interface inputs and outputs, to define input constraints (e.g., type, dependency, whether mandatory) and to provide validation (e.g., file existence). This allows malformed or underspecified inputs to be detected prior to executing the underlying program. The input definition also allows specifying relations between inputs. Often, some input options should not be set together (mutual exclusion) while other inputs need to be set as a group (mutual inclusion). Part of the input specification for the “bet” (Brain Extraction Tool) program from FSL is shown in Listing 2.

Currently, Nipype (version 0.4) is distributed with a wide range of interfaces (see Table 2). Adding new Interfaces is simply a matter of writing a Python class definition as was shown in Listing 1. When a formal specification of inputs and outputs are provided by the underlying software, Nipype can support these programs automatically. For example, the Slicer command line execution modules come with an XML specification that allows Nipype to wrap them without creating individual interfaces.

NODES, MAPNODES, AND WORKFLOWS

Nipype provides a framework for connecting Interfaces to create a data analysis Workflow. In order for Interfaces to be used in a Workflow they need to be encapsulated in either Node or MapNode objects. Node and MapNode objects provide additional functionality to Interfaces. For example, creating a hash of the input state, caching of results, and the ability to iterate over inputs. Additionally, they execute the underlying interfaces in their own uniquely named directories (almost like a sandbox), thus providing a mechanism to isolate and track the outputs resulting from execution of the Interfaces. These mechanisms allow not only for provenance tracking, but aid in efficient pipeline execution.

The MapNode class is a sub-class of Node that implements a MapReduce-like architecture (Dean and Ghemawat, 2008). Encapsulating an Interface within a MapNode allows Interfaces that normally operate on a single input to execute the Interface on multiple inputs. When a MapNode executes, it creates a separate instance of the underlying Interface for every value of an input list and executes these instances independently. When all instances finish running, their results are collected into a list and exposed through the MapNode’s outputs (see Figure 4D). This approach improves...
together within a Workflow. By connecting outputs of some Nodes to inputs of others, the user implicitly specifies dependencies. These are represented internally as a DAG. The current semantics of Workflow do not allow conditionals and hence the graph needs to be acyclic. Workflows themselves can be a node of the Workflow graph (see Figure 1). This enables a hierarchical architecture and encourages Workflow reuse. The Workflow engine validates that all nodes have unique names, ensures that there are no cycles, and prevents connecting multiple outputs to a given input. For example in an fMRI processing Workflow, preprocessing, model fitting, and visualization of results can be implemented as individual Workflows connected together in a main Workflow. This not only improves clarity of designed Workflows but also enables easy exchange of whole subsets. Common Workflows can be shared across different studies within and across laboratories thus reducing redundancy and increasing consistency.

While a neuroimaging processing pipeline could be implemented as a Bash, MATLAB, or a Python script, Nipype explicitly implements a pipeline as a graph. This makes it easy to follow what steps are being executed and in what order. It also makes it easier to go back and change things by simply reconnecting different outputs and inputs or by inserting new Nodes/MapNodes. This alleviates the tedious component of scripting where one has to manually ensure that the inputs and outputs of different processing calls match and that operations do not overwrite each other’s outputs.

A Workflow provides a detailed description of the processing steps and how data flows between Interfaces. Thus it is also a source of provenance information. We encourage users to provide Workflow definitions (as scripts or graphs) as supplementary material when submitting articles. This ensures that at least the data processing part of the published experiment is fully reproducible. Additionally, exchange of Workflows between researchers stimulates efficient use of methods and experimentation.
from nipype.interfaces.base import (TraitedSpec, CommandLineInputSpec, CommandLine, File)

import os

class GZipInputSpec(CommandLineInputSpec):
    input_file = File(desc = "File", exists = True, mandatory = True, argstr = "%s")

class GZipOutputSpec(TraitedSpec):
    output_file = File(desc = "Zip file", exists = True)

class GZipTask(CommandLine):
    input_spec = GZipInputSpec
    output_spec = GZipOutputSpec
    cmd = 'gzip'

    def _list_outputs(self):
        outputs = self.output_spec().get()
        outputs['output_file'] = os.path.abspath(self.inputs.input_file + "\.gz")
        return outputs

if __name__ == '__main__':
    zipper = GZipTask(input_file='an_existing_file')
    print zipper.cmdline
    zipper.run()

**LISTING 1** | An example interface wrapping the *gzip* command line tool and a usage example. This Interface takes a file name as an input, calls gzip to compress it and returns a name of the compressed output file.

```python
class BETInputSpec(FSLCommandInputSpec):
    in_file = File(exists=True,
                   desc = 'input file to skull strip',
                   argstr='\%s', position=0, mandatory=True)

    out_file = File(desc = 'name of output skull stripped image',
                    argstr='\%s', position=1, genfile=True)

    mask = traits.Bool(desc = 'create binary mask image', argstr='\-m')

    functional = traits.Bool(argstr='\-F', xor=('functional', 'reduce_bias'),
                              desc="apply to 4D fMRI data")
```

**LISTING 2** | Part of the input specification for the Brain Extraction Tool (BET) Interface. Full specification covers 18 different arguments. Each attribute of this class is a Traits object which defines an input and its data type (i.e., list of integers), constraints (i.e., length of the list), dependencies (when for example setting one option is mutually exclusive with another, see the xor parameter), and additional parameters (such as argstr and position which describe how to convert an input into a command line argument).

**EXAMPLE – BUILDING A WORKFLOW FROM SCRATCH**

In this section, we describe how to create and extend a typical fMRI processing Workflow. A typical fMRI Workflow can be divided into two sections: (1) preprocessing and (2) modeling. The first one deals with cleaning data from confounds and noise and the second one fits a model to the cleaned data based on the experimental design. The preprocessing stage in this Workflow will consist of only two steps: (1) motion correction (aligns all volumes in a functional run to the mean realigned volume) and (2) smoothing (convolution with a 3D Gaussian kernel). We use SPM Interfaces to define the processing Nodes.

```python
from nipype.pipeline.engine import Node, Workflow
realign = Node(interface=spm.Realign(), name="realign")
realign.inputs.register_to_mean = True
smooth = Node(interface=spm.Smooth(), name="smooth")
smooth.inputs.fwhm = 4
```
DataSink on the other side provides means for storing selected results in a specified location. It supports automatic creation of folders, simple substitutions, and regular expressions to alter target filenames. In this example we store the statistical (T maps) resulting from contrast estimation.

A Workflow defined this way (see Figure 3, for full code see Supplementary Material) is ready to run. This can be done by calling `run()` method of the master Workflow.

If the `run()` method is called twice, the Workflow input hashing mechanism ensures that none of the Nodes are executed during the second run if the inputs remain the same. If, however, a highpass filter parameter of `specify_model` is changed, some of the Nodes (but not all) would have to rerun. Nipype automatically determines which Nodes require rerunning.

**Iterables – parameter space exploration**

Nipype provides a flexible approach to prototype and experiment with different processing strategies, through the unified and uniform access to a variety of software packages (Interfaces) and creating data flows (Workflows). However, for various neuroimaging

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List of software packages fully or partially supported by Nipype. For more details see http://nipy.org/nipype/interfaces.
tasks, there is often a need to explore the impact of variations in parameter settings (e.g., how do different amounts of smoothing affect group statistics, what is the impact of spline interpolation over trilinear interpolation). To enable such parametric exploration, Nodes have an attribute called *iterables*.

When an *iterable* is set on a Node input, the Node, and its subgraph are executed for each value of the iterable input (see Figure 4 *iterables_vs_mapnode*). Iterables can also be set on multiple inputs of a Node (e.g., `somenode.iterables = [("input1","1,2,3"), ("input2","a","b")])`. In such cases, every combination of those values is used as a parameter set (the prior example would result in the following parameter sets: (1, "a"), (1, "b"), (2, "a"), etc...). This feature is especially useful to investigate interactions between parameters of intermediate stages with respect to the final results of a workflow. A common use-case of *iterables* is to execute the same Workflow for many subjects in an fMRI experiment and to simultaneously look at the impact of parameter variations on the results of the Workflow.

It is important to note that unlike MapNode, which creates copies of the underlying interface for every element of an input of type list, *iterables* operate on the subgraph of a node and create copies not only of the node but also of all the nodes dependent on it (see Figure 4).

**PARALLEL DISTRIBUTION AND EXECUTION PLUG-INS**

Nipype supports executing Workflows locally (in series or parallel) or on load-balanced grid-computing clusters (e.g., SGE, Torque, or even via SSH) through an extensible plug-in interface. No change is needed to the Workflow to switch between these execution modes. One simply calls the Workflow’s run function with a different plug-in and its arguments. Very often different components of a Workflow can be executed in parallel and even more so when the same Workflow is being repeated on multiple parameters (e.g., subjects). Adding support for additional cluster management systems does not require changes in Nipype, but simply writing a plug-in extension conforming to the plug-in API.

The Workflow engine sends an execution graph to the plug-in. Executing the Workflow in series is then simply a matter of performing a topological sort on the graph and running each node in the sorted order. However, Nipype also provides additional plug-ins that use Python’s multi-processing module, use IPython (includes SSH-based, SGE, LSF, PBS, among others) and provide native interfaces to SGE or PBS/Torque clusters. For all of these, the graph structure defines the dependencies as well as which nodes can be executed in parallel at any given stage of execution.

One of the biggest advantages of Nipype’s execution system is that parallel execution using local multi-processing plug-in does not require any additional software (such as cluster managers like SGE) and therefore makes prototyping on a local multi-core workstations easy. However for bigger studies and complex Workflows, a high-performance computing cluster can provide substantial improvements in execution time. Since there is a clear separation between definition of the Workflow and its execution, Workflows can be executed in parallel (locally or on a cluster) without any modification. Transitioning from developing a processing pipeline on a single subject on a local workstation to executing it on a bigger cohort on a cluster is therefore seamless.

Rerunning workflows has also been optimized. When a Node or MapNode is run, the framework will actually execute the underlying interface only if inputs have changed relative to prior execution. If not, it will simply return cached results.

**THE FUNCTION INTERFACE**

One of the Interfaces implemented in Nipype requires special attention: The Function Interface. Its constructor takes as arguments Python function pointer or code, list of inputs, and list of outputs. This allows running any Python code as part of a Workflow. When combined with libraries such as Nibabel (neuroimaging data input and output), Numpy/Scipy (array representation and processing) and scikit-learn or PyMVPA (machine learning and data mining) the Function Interface provides means for rapid prototyping of complex data processing methods. In addition, by using the

![Figure 4](https://example.com/figure4.png)

**FIGURE 4 | Workflow modification using *iterables* and MapNodes.** If we take the processing pipeline (A) and set *iterables* parameter of DataGrabber to a list of two subjects, Nipype will effectively execute graph (B). Identical processing will be applied to every subject from the list. Iterables can be used in a graph on many levels. For example, setting *iterables* on Smooth FWHM to a list of 4 and 8 mm will result in graph (C). In contrast to *iterables*, MapNode branches within a node of the graph and also merges the results of the branches, effectively performing a MapReduce operation (D).
Function Interface users can avoid writing their own Interfaces which is especially useful for ad hoc solutions (e.g., calling an external program that has not yet been wrapped as an Interface).

**WORKFLOW VISUALIZATION**

To be able to efficiently manage and debug Workflows, one has to have access to a graphical representation. Using graphviz (Ellson et al., 2002), Nipype generates static graphs representing Nodes and connections between them. In the current version four types of graphs are supported: *orig* – does not expand inner Workflows, *flat* – expands inner workflows, *exec* – expands workflows and iterables, and *hierarchical* – expands workflows but maintains their hierarchy.

Graphs can be saved in a variety of file formats including Scalable Vector Graphics (SVG) and Portable Network Graphics (PNG; see Figures 3 and 6).

**CONFIGURATION OPTIONS**

Certain options concerning verbosity of output and execution efficiency can be controlled through configuration files or variables. These include, among others, `hash_method` and `remove_unnecessary_outputs`. As explained before, rerunning a Workflow only recomputes those Nodes whose inputs have changed since the last run. This is achieved by recording a hash of the inputs. For files there are two ways of calculating the hash (controlled by the `hash_method` configuration option): *timestamp* – based only on the size and modification time and *content* – based on the content of the file. The first one is faster, but does not deal with the situation when an identical copy overwrites the file. The second one can be slower especially for big files, but can tell that two files are identical even if they have different modification times. To allow efficient recomputation Nipype has to store outputs of all Nodes. This can generate a significant amount of data for typical neuroimaging studies. However, not all outputs of every Node are used as inputs to other Nodes or relevant to the final results. Users can decide to remove those outputs (and save some disk space) by setting the `remove_unnecessary_outputs` to `True`. These and other configuration options provide a mechanism to streamline the use of Nipype for different applications.

**DEPLOYMENT**

Nipype supports GNU/Linux and Mac OS X operating systems (OS). A recent Internet survey based study (Hanke and Halchenko, 2011) showed that GNU/Linux is the most popular platform in the neuroimaging community and together with Mac OS X is used by over 70% of neuroimagers. There are not theoretical reasons why Nipype should not work on Windows (Python is a cross-platform language), but since most of the supported software (for example FSL) requires a Unix based OS, Nipype has not been tested on this platform.

We currently provide three ways of deploying Nipype on a new machine: manual installation from sources, PyPI repository, and from package repositories on Debian-based systems. Manual installation involves downloading a source code archive and running a standard Python installation script (distutils). This way the user has to take care of installing all of the dependencies. Installing from PyPI repository lifts this constraint by providing dependency information and automatically installing required packages. Nipype is available from standard repositories on recent Debian and Ubuntu releases. Moreover, NeuroDebian (Hanke et al., 2010) repository provides the most recent releases of Nipype for Debian-based systems and a NeuroDebian Virtual Appliance making it easy to deploy Nipype and other imaging tools in a virtual environment on other OS, e.g., Windows. In addition to providing all core dependencies and automatic updates NeuroDebian also provides many of the software packages supported by Nipype (AFNI, FSL, Micron, etc.), making deployment of heterogeneous Nipype pipelines more straightforward.

**DEVELOPMENT**

Nipype is trying to address the problem of interacting with the ever changing universe of neuroimaging software in a sustainable manner. Therefore the way its development is managed is a part of the solution. Nipype is distributed under BSD license which allows free copying, modification, and distribution and additionally meets all the requirements of open-source definition (see Open-Source Initiative) and Debian Free Software Guidelines. Development is carried out openly through distributed version control system (git via GitHub) in an online community. The current version of the source code together with complete history is accessible to everyone. Discussions between developers and design decisions are done on an open access mailing list. This setup encourages a broader community of developers to join the project and allows sharing of the development resources (effort, money, information, and time).

In these previous paragraphs, we presented key features of Nipype that facilitate rapid development and deployment of analysis procedures in laboratories, and address all of the issues described earlier. In particular, Nipype provides: (1) uniform access to neuroimaging analysis software and usage information; (2) a framework for comparative algorithm development and dissemination; (3) an environment for methodological continuity and paced training of new personnel in laboratories; (4) computationally efficient execution of neuroimaging analysis; and (5) a mechanism to capture the data processing details in compact scripts and graphs. In the following section, we provide examples to demonstrate these solutions.

**USAGE EXAMPLES**

**UNIFORM ACCESS TO TOOLS, THEIR USAGE, AND EXECUTION**

Users access Interfaces by importing them from Nipype modules. Each neuroimaging software distribution such as FSL, SPM, Camino, etc., has a corresponding module in the `nipype.interfaces` namespace.

```python
from nipype.interfaces.camino import DTIFit
```

The `help()` function for each interface prints the inputs and the outputs associated with the interface.

---

1[http://nipy.org/nipype/]
2[http://pypi.python.org/pypi/nipype/]
3[http://neuro.debian.net]
4[http://www.opensource.org/docs/osd]
5[http://www.debian.org/social_contract#guidelines]
6[http://github.com/nipy/nipype]
The output of the `help()` function is standardized across all Interfaces. It is automatically generated based on the traited input and output definitions and includes information about required inputs, types, and default value. Alternatively, extended information is available in the form of auto-generated HTML documentation on the Nipype website (see Figure 5). This extended information includes examples that demonstrate how the interface can be used.

For every Interface, input values are set through the `inputs` field:

```python
fit.inputs.scheme_file = 'A.scheme'
fit.inputs.in_file = 'tensor_fitted_data.Bfloat'
```

When trying to set an invalid input type (for example a non-existing input file, or a number instead of a string) the Nipype framework will display an error message. Input validity checking before actual Workflow execution saves time. To run an Interface user needs to call `run()` method:

```python
fit.run()
```

At this stage the framework checks if all mandatory inputs are set and all input dependencies are satisfied, generating an error if either of these conditions are not met.

---

**DTIFit**

Wraps command `dtfit`

Reads diffusion MRI data, acquired using the acquisition scheme detailed in the scheme file, from the data file. Use non-linear fitting instead of the default linear regression to the log measurements. The data file stores the diffusion MRI data in voxel order with the measurements stored in big-endian format and ordered as in the scheme file. The default input data type is four-byte float. The default output data type is eight-byte double. See `modelfit` and `camino` for the format of the data file and scheme file. The program fits the diffusion tensor to each voxel and outputs the results, in voxel order and as big-endian eight-byte doubles, to the standard output. The program outputs eight values in each voxel: [exit code, In(S(0)), Dxx, Dxy, Dxz, Dyw, Dwz, Dzz]. An exit code of zero indicates no problems. For a list of other exit codes, see `modelfit()`. The entry S(0) is an estimate of the signal at q=0.

**Example**

```python
>>> import nipype.interfaces.camino as camino
>>> fit = camino.DTIFit()
>>> fit.inputs.scheme_file = 'A.scheme'
>>> fit.inputs.in_file = 'tensor_fitted_data.Bfloat'
>>> fit.run()
```

**Inputs**

<table>
<thead>
<tr>
<th>Mandatory</th>
<th>(an existing file name)</th>
</tr>
</thead>
<tbody>
<tr>
<td><code>in_file</code> : voxel-order data filename</td>
<td></td>
</tr>
<tr>
<td><code>scheme_file</code> : Camino scheme file (b values / vectors, see camino.fsl2scheme)</td>
<td></td>
</tr>
</tbody>
</table>

<table>
<thead>
<tr>
<th>Optional</th>
<th>(a string)</th>
</tr>
</thead>
<tbody>
<tr>
<td><code>args</code> : Additional parameters to the command</td>
<td></td>
</tr>
<tr>
<td><code>environ</code> : (a dictionary with keys which are a value of type 'str' and with values which are a value of type 'str') Environment variables</td>
<td></td>
</tr>
<tr>
<td><code>ignore_exception</code> : (a boolean)</td>
<td></td>
</tr>
<tr>
<td><code>non_linear</code> : Use non-linear fitting instead of the default linear regression to the log measurements.</td>
<td></td>
</tr>
<tr>
<td><code>out_file</code> : (a file name)</td>
<td></td>
</tr>
</tbody>
</table>

**Outputs**

<table>
<thead>
<tr>
<th>(an existing file name)</th>
</tr>
</thead>
<tbody>
<tr>
<td><code>tensor_fitted</code> : path/name of 4D volume in voxel order</td>
</tr>
</tbody>
</table>

**Figure 5** HTML help page for dtfit command from Camino. This was generated based on the Interface code: description and example was taken from the class docstring and inputs/outputs were list was created using traited input/output specification.
Nipype standardizes running and accessing help information irrespective of whether the underlying software is a MATLAB program, a command line tool, or Python module. The framework deals with translating inputs into appropriate form (e.g., command line arguments or MATLAB scripts) for executing the underlying tools in the right way, while presenting the user with a uniform interface.

**A FRAMEWORK FOR COMPARATIVE ALGORITHM DEVELOPMENT AND DISSEMINATION**

Uniform semantics for interfacing with a wide range of processing methods not only opens the possibility for richer Workflows, but also allows comparing algorithms that are designed to solve the same problem across and within such diverse Workflows. Typically, such an exhaustive comparison can be time-consuming, because of the need to deal with interfacing different software packages. Nipype simplifies this process by standardizing the access to the software. Additionally, the iterables mechanism allows users to easily extend such comparisons by providing a simple mechanism to test different parameter sets.

Accuracy or efficiency of algorithms can be determined in an isolated manner by comparing their outputs or execution time or memory consumption on a given set of data. However, researchers typically want to know how different algorithms used at earlier stages of processing might influence the final output or statistics they are interested in. As an example of such use, we have compared voxelwise isotropic, voxelwise anisotropic, and surface based smoothing all for two levels of FWHM – 4 and 8 mm. First one is the standard convolution with Gaussian kernel as implemented in SPM. Second one involves smoothing only voxels of similar intensity in attempt to retain structure. This was implemented in SUSAN from FSL (Smith, 1992). Third method involves reconstructing surface of the cortex and smoothing along it (Hagler et al., 2006). This avoids bleeding of signal over sulci.

Establishing parameters from data and smoothing using SUSAN is already built into Nipype as a Workflow. It can be created using create_susan_smooth() function. It has similar inputs and outputs as SPM Smooth Interface. Smoothing on a surface involves doing a full cortical reconstruction from T1 volume using FreeSurfer (Fischl et al., 1999) followed by coregistering functional images to the reconstructed surface using BBRegister (Greve and Fischl, 2009). Finally, the surface smoothing algorithm from FreeSurfer is called.

Smoothed EPI volumes (direct/local influence) and statistical maps (indirect/global influence), along with the pipeline used to generate them can be found in Figures 6 and 7. Full code used to generate this data can be found in the Supplementary Material. This comparison serves only to demonstrate Nipype capabilities; a comparison between smoothing methods is outside of the scope of this paper.

**FIGURE 6** | Graph showing the workflow used for the smoothing methods and parameters comparison. The gray shaded nodes have iterables parameter set. This allows easy iteration over all combinations of FWHM and smoothing algorithms used in the comparison.
AN ENVIRONMENT FOR METHODOLOGICAL CONTINUITY AND PACED TRAINING OF NEW PERSONNEL IN LABORATORIES

Neuroimaging studies in any laboratory typically use similar data processing methods with possibly different parameters. Nipype Workflows can be very useful in dividing the data processing into reusable building blocks. This not only improves the speed of building new Workflows but also reduces the number of potential errors, because a well tested piece of code is being reused (instead of being reimplemented every time). Since a Workflow definition is an abstract and simplified representation of the data processing stream, it is much easier to describe and hand over to new project personnel. Furthermore, a data independent Workflow definition (see Figure 8) enables sharing Workflows within and across research laboratories. Nipype provides a high-level abstraction mechanism for exchanging knowledge and expertise between researchers focused on methods in neuroimaging and those interested in applications.

The uniform access to Interfaces and the ease of use of Workflows in Nipype helps with training new staff. Composition provided by Workflows allows users to gradually increase the level of details when learning how to perform neuroimaging analysis. For example user can start with a “black box” Workflow that does analysis from A to Z, and gradually learn what the sub-components (and their sub-components) do. Playing with Interfaces in an interactive console is also a great way to learn how different algorithms work with different parameters without having to understand how to set them up and execute them.

COMPUTATIONALLY EFFICIENT EXECUTION OF NEUROIMAGING ANALYSIS

A computationally efficient execution allows for multiple rapid iterations to optimize a Workflow for a given application. Support for optimized local execution (running independent processes in parallel, rerunning only those steps that have been influenced by the changes in parameters or dependencies since the last run) and exploration of parameter space eases Workflow development. The Nipype package provides a seamless and flexible environment for executing Workflows in parallel on a variety of environments from local multi-core workstations to high-performance clusters. In the SPM workflow for single subject functional data analysis (see Figure 9), only a few components can be parallelized. However, running this Workflow across several subjects provides room for embarrassingly parallel execution. Running this Workflow in distributed mode for 69 subjects on a compute cluster (40 cores distributed across 6 machines) took 1 h and 40 min relative to the 32-min required to execute the analysis steps in series for a single subject on the same cluster. The difference from the expected runtime of 64 min (32 min for the first 40 subjects and another 32 min for the remaining 29 subjects) stems from disk I/O and other network and processing resource bottlenecks.

CAPTURES DETAILS OF ANALYSIS REQUIRED TO REPRODUCE RESULTS

The graphs and code presented in the examples above capture all the necessary details to rerun the analysis. Any user, who has the same versions of the tools installed on their machine and access to the data and scripts, will be able to reproduce the results of the study. For example, running Nipype within the NeuroDebian framework can...
provide access to specific versions of the underlying tools. This provides an easy mechanism to be compliant with the submitting data and scripts/code mandates of journals such as PNAS and Science.

**DISCUSSION**

Current neuroimaging software offer users an incredible opportunity to analyze their data in different ways, with different underlying assumptions. However, this heterogeneous collection of specialized applications creates several problems: (1) No uniform access to neuroimaging analysis software and usage information; (2) No framework for comparative algorithm development and dissemination; (3) Personnel turnover in laboratories often limit methodological continuity and training new personnel takes time; (4) Neuroimaging software packages do not address computational efficiency; and (5) Method sections of journal articles are often inadequate for reproducing results.

We addressed these issues by creating Nipype, an open-source, community-developed initiative under the umbrella of Nipy. Nipype solves these issues by providing uniform Interfaces to existing neuroimaging software and by facilitating interaction between these packages within Workflows. Nipype provides an environment that encourages interactive exploration of algorithms from different packages (e.g., SPM, FSL), eases the design of Workflows within and between packages, and reduces the learning curve necessary to use different packages. Nipype is addressing limitations of existing pipeline systems and creating a collaborative platform for neuroimaging software development in Python, a high-level scientific computing language.

We use Python for several reasons. It has extensive scientific computing and visualization support through packages such as SciPy, NumPy, Matplotlib, and Mayavi (Pérez et al., 2010; Millman and Aivazis, 2011). The Nibabel package provides support for reading and writing common neuroimaging file formats (e.g., NIFTI, ANALYZE, and DICOM). Being a high-level language, Python supports rapid prototyping, is easy to learn and adopt and is available across all major OS. At the same time Python allows to seamlessly bind with C code (using Weave package) for improved efficiency of critical subroutines.

Python is also known to be a good choice for the first programming language to learn (Zelle, 1999) and is chosen as the language for introductory programming at many schools and universities.\(^1\)

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\(^1\)http://wiki.python.org/moin/SchoolsUsingPython
FIGURE 9 | Single subject fMRI Workflow used for benchmarking parallel execution.
Being a generic and free language, with various extensions available “out of the box,” it has allowed many researchers to start implementing and sharing their ideas with minimal knowledge of Python, while learning more of the language and programming principles along the way. Many such endeavors later on became popular community-driven FOSS projects, attracting users and contributors, and even outlasting the involvement of the original authors. Python has already been embraced by the neuroscientific community and is rapidly gaining popularity (Bednar, 2009; Goodman and Brette, 2009). The Connectome Viewer Toolkit (Gerhard et al., 2011), Dipy (Garyfallidis et al., 2011), NiBabel13, Nipy15, NiTime (Rokem et al., 2009), PyMVPA (Hanke et al., 2009), PyXNAT (Schwartz et al., 2011), and Scikits-Learn14 are just a few examples of neuroimaging related software written in Python. Nipype, based on Python, thus has immediate access to this extensive community and its software, technological resources and support structure.

Nipype provides a formal and flexible framework to accommodate the diversity of imaging software. Within the neuroimaging community, not all software is limited to well-behaved command line tools. Furthermore, a number of these tools do not have well defined inputs, outputs, or usage help. Although, currently we use Enthought Traits to define inputs and outputs of interfaces, such definitions could be easily translated into instances of XML schemas compatible with other pipeline frameworks. On the other hand, when a tool provides a formal XML description of their inputs and outputs (e.g., Slicer 3D, BRAINS), it is possible to take these definitions and automatically generate Nipype wrappers for those classes.

Nipype development welcomes input and contributions from the community. The source code is freely distributed under a BSD license allowing anyone any use of the software and Nipype conforms to the Open Software Definition of the Open-Source Initiative. Development process is fully transparent and encourages contributions from users from all around the world. The diverse and geographically distributed user and developer base makes Nipype a flexible project that takes into account needs of many scientists.

Improving openness, transparency, and reproducibility of research has been a goal of Nipype since its inception. A Workflow definition is, in principle, sufficient to reproduce the analysis. Since it was used to analyze the data, it is more detailed and accurate than a typical methods description in a paper, but also has the advantage of being reused and shared within and across laboratories. Accompanying a publication with a formal definition of the processing pipeline (such as a Nipype script) increases reproducibility and transparency of research. The Interfaces and Workflows of Nipype capture neuroimaging analysis knowledge and the evolution of methods. Although, at the execution level, Nipype already captures a variety of provenance information, this aspect can be improved by generating provenance reports defined by a standardized XML schema (Mackenzie-Graham et al., 2008).

Increased diversity of neuroimaging data processing software has made systematic comparison of performance and accuracy of underlying algorithms essential (for examples, see Klein et al., 2009, 2010). However, a platform for comparing algorithms, either by themselves or in the context of an analysis workflow, or determining optimal workflows in a given application context (e.g., Churchill et al., 2011), does not exist. Furthermore, in this context of changing hardware and software, traditional analysis approaches may not be suitable in all contexts (e.g., data from 32-channel coils which show a very different sensitivity profile, or data from children). Nipype can make such evaluations, design of optimal workflows, and investigations easier (as demonstrated via the smoothing example above), resulting in more efficient data analysis for the community.

**SUMMARY**

We presented Nipype, an extensible Python library and framework that provides interactive manipulation of neuroimaging data through uniform Interfaces and enables reproducible, distributed analysis using the Workflow system. Nipype has encouraged the scientific exploration of different algorithms and associated parameters, eased the development of Workflows within and between packages and reduced the learning curve associated with understanding the algorithms, APIs, and user interfaces of disparate packages. An open, community-driven development philosophy provides the flexibility required to address the diverse needs in neuroimaging analysis. Overall, Nipype represents an effort toward collaborative, open-source, reproducible, and efficient neuroimaging software development and analysis.

**ACKNOWLEDGMENTS**

A list of people who have contributed code to the project is available at http://github.com/nipy/nipype/contributors. We thank Fernando Perez, Matthew Brett, Gael Varoquaux, Jean-Baptiste Poline, Bertrand Thirion, Alexis Roche, and Jarrod Millman for technical and social support and for design discussions. We would like to thank Prof. John Gabrieli’s laboratory at MIT for testing Nipype through its evolutionary stages, in particular, Tyler Perrachione and Gretchen Reynolds. We would also like to thank the developers of FreeSurfer, FSL, and SPM for being supportive of the project and providing valuable feedback on technical issues. We would like to thank James Bednar, Stephan Gerhard, and Helen Ramsden for providing feedback during the preparation of the manuscript. Satrajit Ghosh would like to acknowledge support from NIBIB R03 EB008673 (PI: Ghosh and Whitfield-Gabrieli), the Ellison Medical Foundation, Katrien Vander Straeten, and Amie Ghosh. Krzysztof Gorgolewski would like to thank his supervisors Mark Bastin, Cyril Pernet, and Amos Storkey for their support during this project.

**SUPPLEMENTARY MATERIAL**

The Supplementary Material for this article can be found online at http://www.frontiersin.org/neuroinformatics/10.3389/fninf.2011.00013/abstract
REFERENCES


Conflict of Interest Statement: The authors declare that the research was conducted in the absence of any commercial or financial relationships that could be construed as a potential conflict of interest.

Received: 23 June 2011; accepted: 23 July 2011; published online: 22 August 2011.


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