

Overview: Landmark

Landmarking Program to Create/Edit Landmark Points Between Samples

<http://open2dprot.sourceforge.net/Landmark>

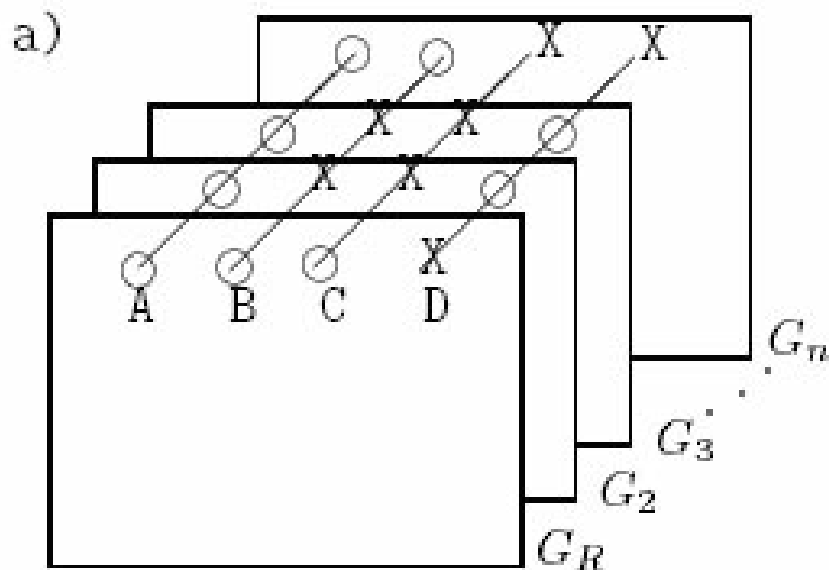
Introduction

- Spot pairing between spots from two samples may require a small set of corresponding spots known to be the same in the two samples.
- Some pairing programs require landmark data.
- Set of landmark points must be created prior to spot pairing and are stored in a landmark database.

Introduction (cont.)

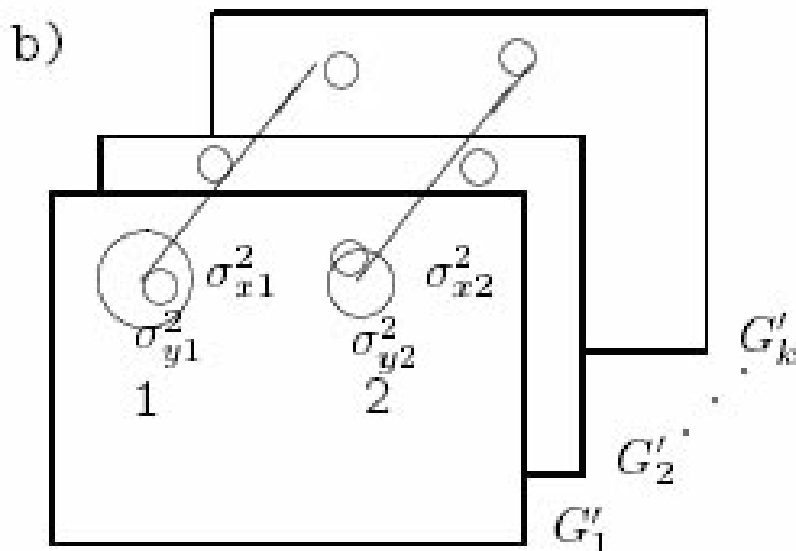
- Landmark program is a step [3] pipeline module used for creating and editing a set of landmarks between a reference sample and sample.
- It uses a graphical user interface to let the user interactively assign corresponding spots.
- It saves the landmark set data for a pair of samples in an XML landmark database used by the other Open2Dprot pipeline modules.

Composite Samples Database (CSD) Paradigm



Proteomic composite samples database (CSD) consisting of a set of n samples G_1, G_2, \dots, G_n with representative sample $G_r = G_1$

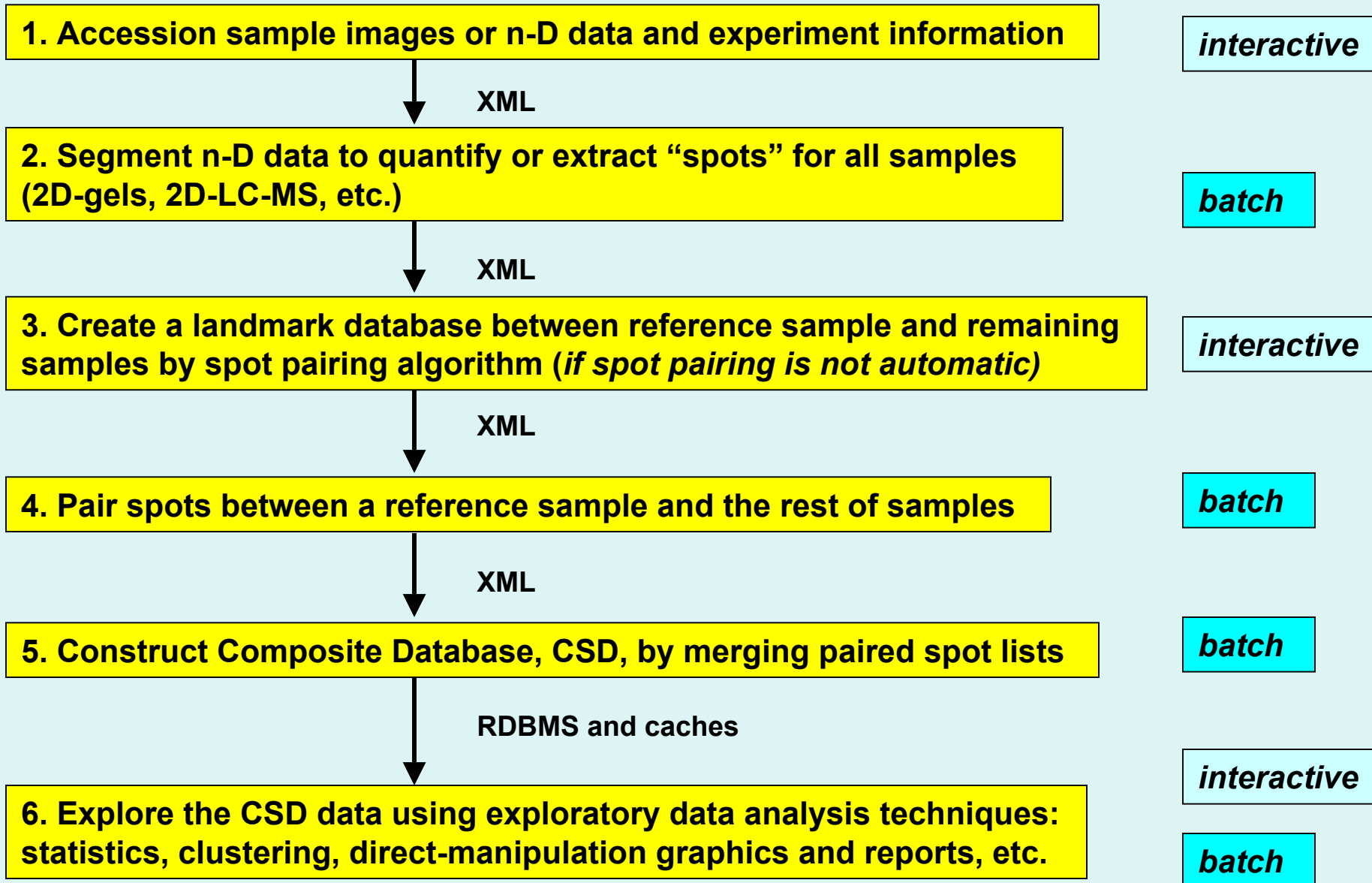
Expression profiles A, B, C, ...



A canonical sample database is a statistical representation of the CSD spot geometry and quantification that could be used for data mining

in Lemkin *et al.*,
*Computers Biomedical
Research*, 1981

Basic Open n-D Analysis Pipeline



Initial Open n-D Data-Mining Tools

- **Accession n-D sample images or n-D data and experiment data**
- **Quantify 'spots' from sample images or peptide clusters**
- **Pair spots between samples and a reference sample**
- **Construct composite sample database for exploratory data analysis**
- **Manage subsets of proteins in the database**
- **Manage replicate samples and condition sets of samples**
- **Analyze expression profiles for multiple conditions**
- **Data-filter protein sets by statistics, clustering, set membership**
- **Direct-manipulation of data in graphics, spreadsheets**
- **Integrate R language statistical, clustering, classifiers, class prediction, and other methods**
- **Integrate access to Internet proteomic/genomic/function data servers for user-specified protein sets**

Open2Dprot Pipeline Subprojects



Open2Dprot - Subprojects - Netscape

File Edit View Go Bookmarks Tools Window Help

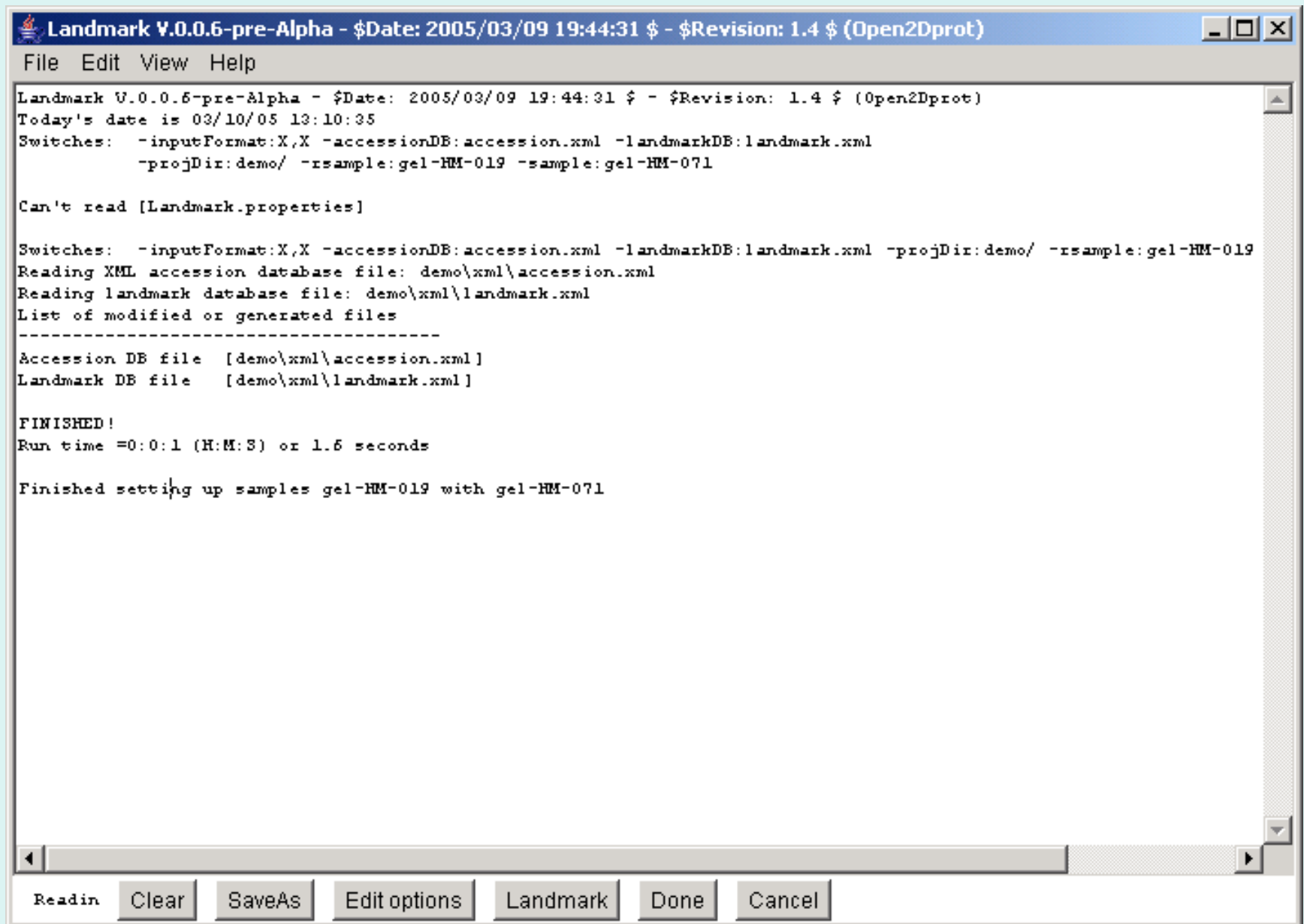
Open2Dprot pipeline subprojects

[Open2Dprot](#) consists of a series of coordinated [Open2Dprot pipeline processing modules](#). The scheduler program, also called Open2Dprot, will schedule and run the modules in the pipeline after doing a data-dependency analysis. By using XML as the "glue" between modules, it is possible to substitute alternate modules at the various pipeline steps. As pipeline modules and alternate modules become available, they will be added to this table. *We encourage the donation of alternate pipeline processing modules which will be added to this table.*

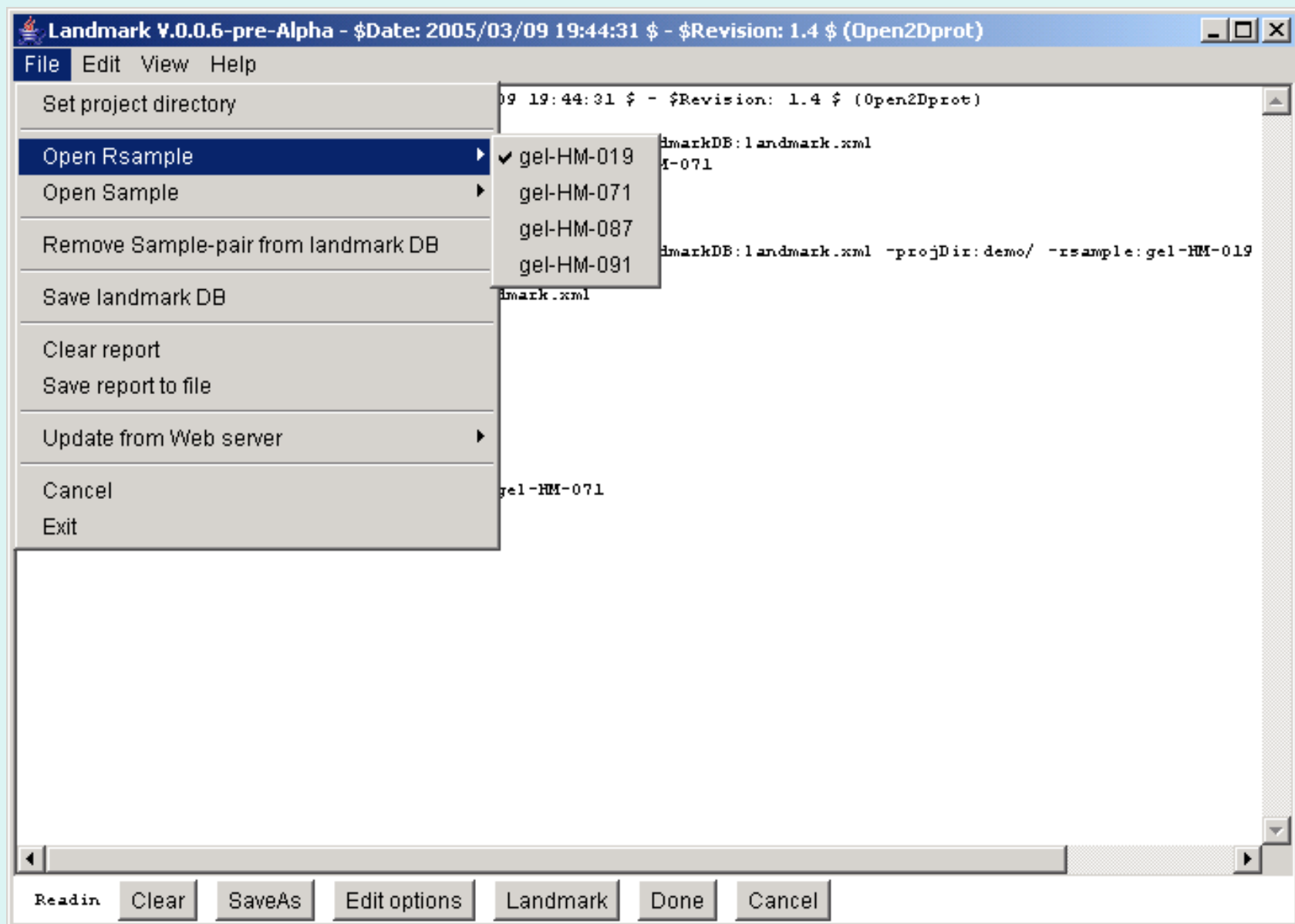
We will be using a common [O2Plib library](#) in the Open2Dprot pipeline modules. This will help ensure that they use the same conventions, data structures and XML data interchange formats.

Subproject Home	Download	Documentation	Overview (PDF)	PDF documents	Version	Revision history	Status	Pipeline step
Open2Dprot project	(see below)	Open2Dprot	Open2Dprot project	Open2Dprot project	Open2Dprot project	Open2Dprot project	<i>Open2Dprot design prototype</i>	
Open2Dprot program	Open2Dprot program	Open2Dprot program	Open2Dprot program	Open2Dprot program	Open2Dprot program	Open2Dprot program	<i>Open2Dprot pre-alpha program</i>	[scheduler]
 Ace	Ace	Ace	Ace	Ace	Ace	Ace	<i>Acession pre-alpha</i>	[1]
 SG2	SG2	SG2	SG2	SG2	SG2	SG2	<i>Seg2Dgel pre-alpha</i>	[2]
Landmark	Landmark	Landmark	Landmark	Landmark	Landmark	Landmark	<i>Landmark pre-alpha</i>	[3]
 CMP	CMP	CMP	CMP	CMP	CMP	CMP	<i>CmpSpots pre-alpha</i>	[4]
BuildCSD	BuildCSD	BuildCSD	BuildCSD	BuildCSD	BuildCSD	BuildCSD	<i>BuildCSD design prototype</i>	[5]
CSDminer	CSDminer	CSDminer	CSDminer	CSDminer	CSDminer	CSDminer	<i>CSDminer design prototype</i>	[6]
O2Plib	O2Plib.jar	O2Plib	O2Plib	O2Plib	O2Plib	O2Plib	<i>O2Plib pre-alpha</i>	--common--

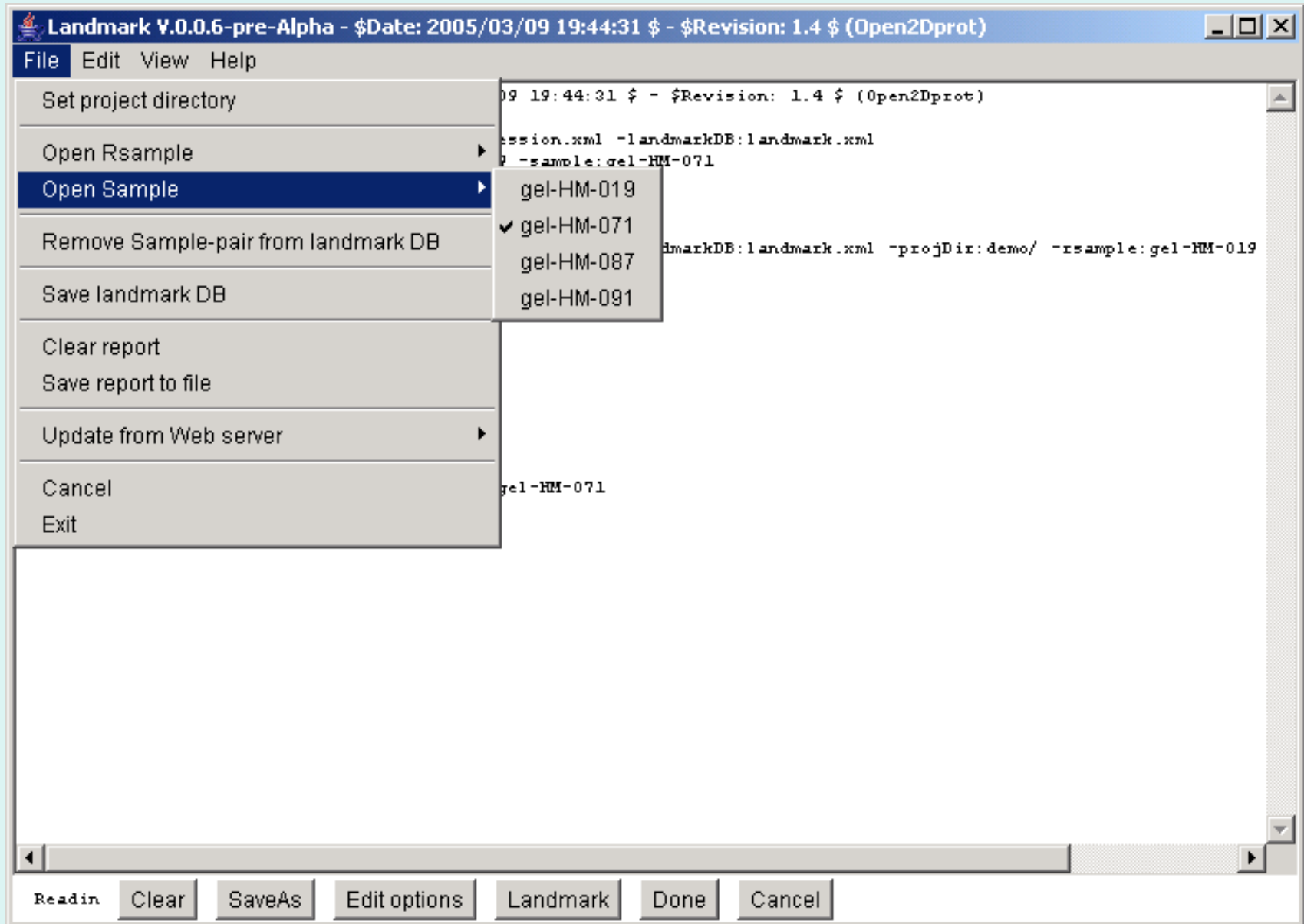
Main Window



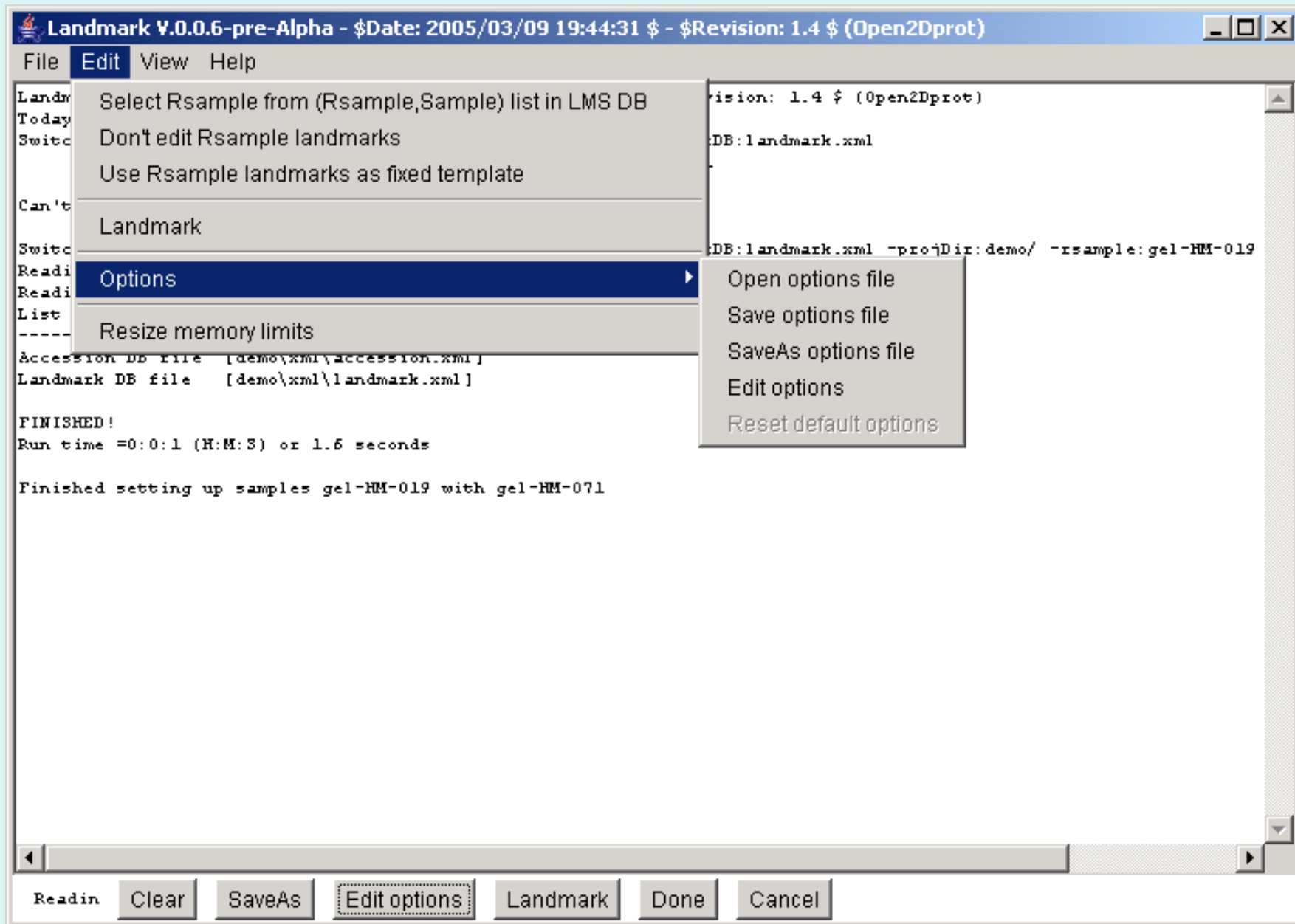
File Menu - Select Reference Sample



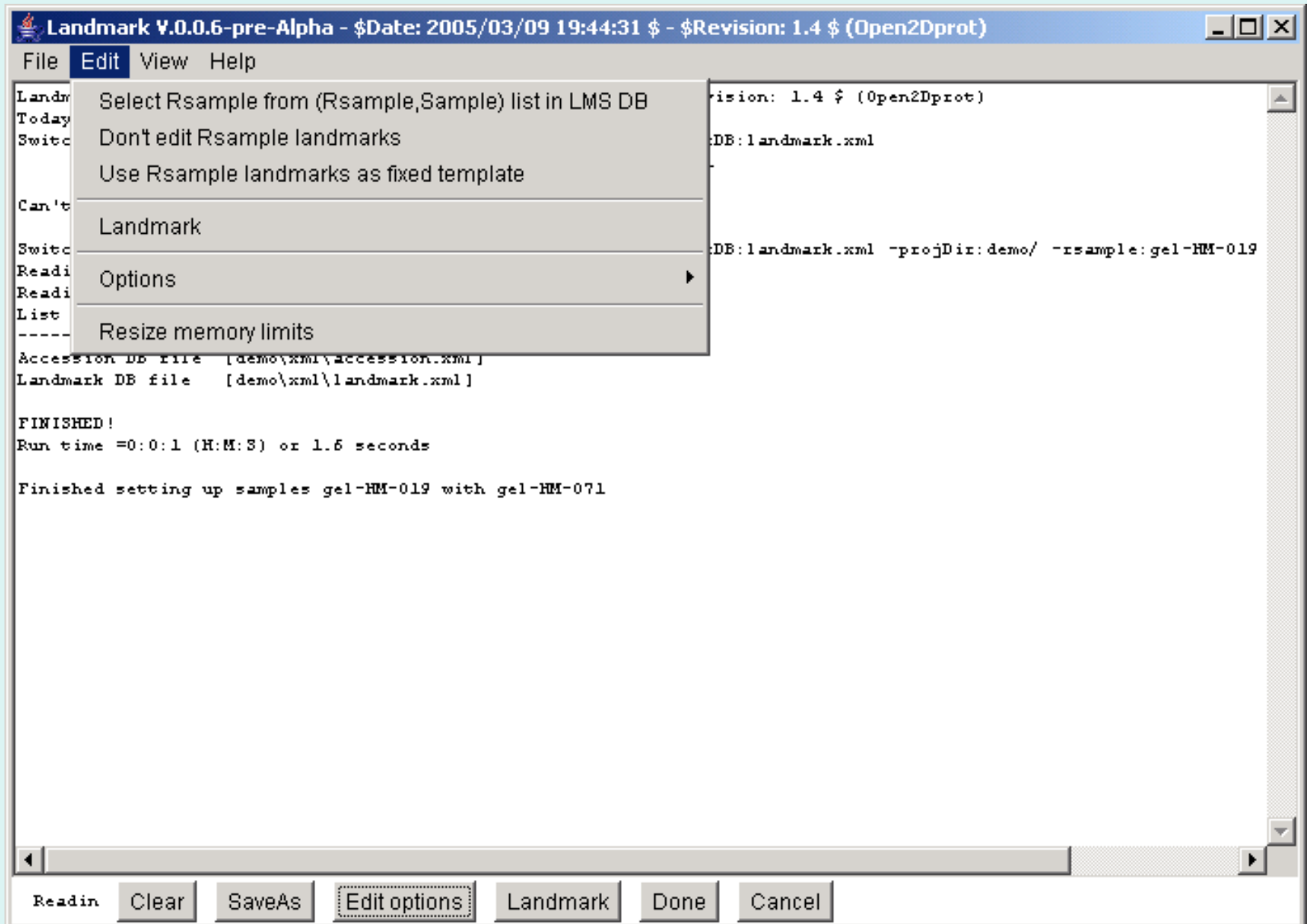
File Menu - Select Current Sample



File Menu - Update From Web Server



Edit Menu



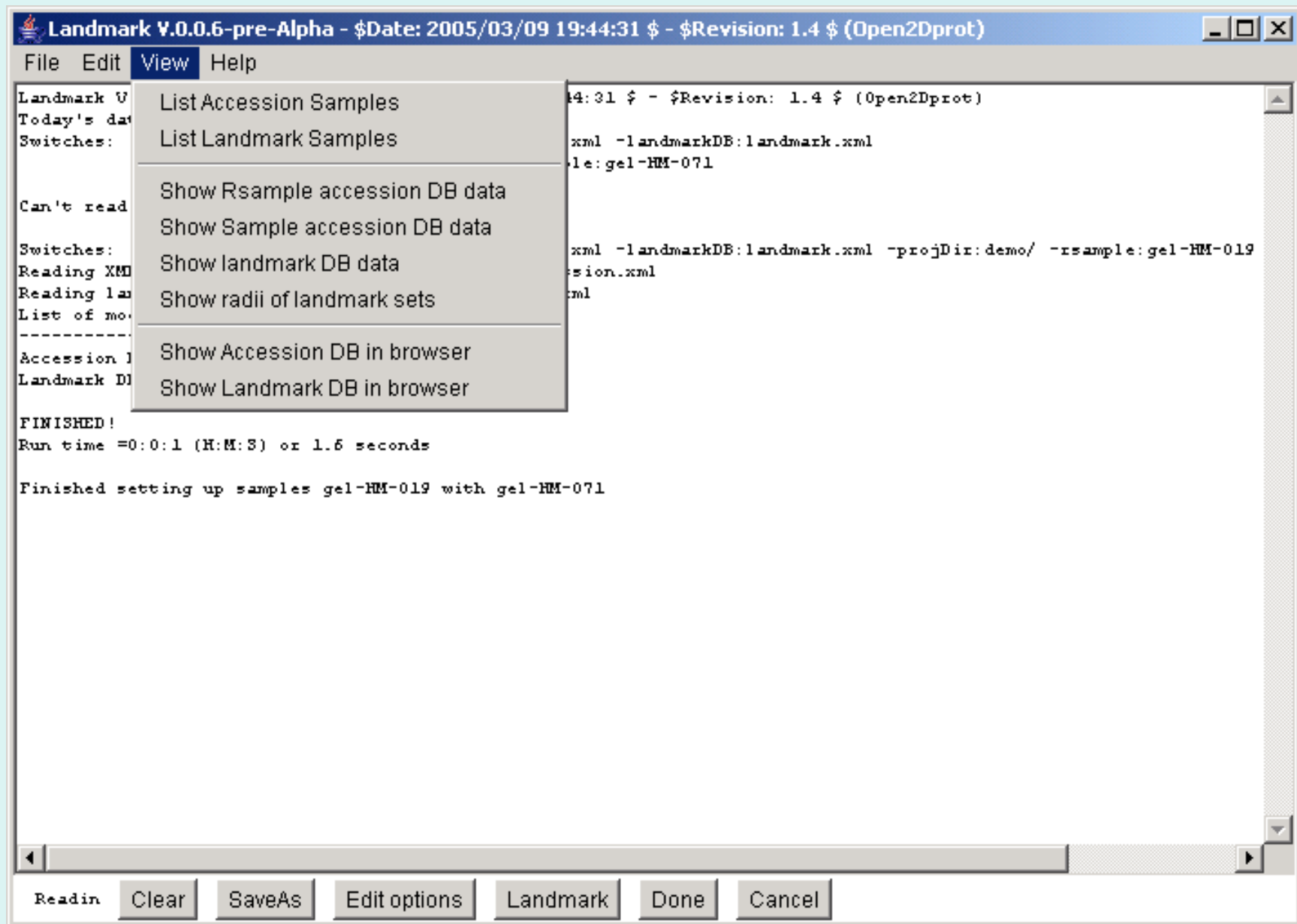
Edit Menu - command line options

Enter sample image and select switch options

Change the startup parameters, then press 'Set new options' button to save them.
At that point, you can press 'Pair spots' to pair the Rsample and Sample with the new parameters. You may also enter new Rsample and Sample data file names and/or edit switch options and threshold sliders.

<input checked="" type="checkbox"/> -accessionFile	accession.xml
<input checked="" type="checkbox"/> -backupDatabases	---
<input type="checkbox"/> -commutativeLMS	---
<input type="checkbox"/> -debugBits	0,0
<input type="checkbox"/> -default	---
<input checked="" type="checkbox"/> -demo	---
<input type="checkbox"/> -dtd	---
<input checked="" type="checkbox"/> -inputFormat	X,X
<input checked="" type="checkbox"/> -landmarkFile	landmark.xml
<input checked="" type="checkbox"/> -projDir Project directory <input type="button" value="Browse dir"/>	demo\
<input type="checkbox"/> -propertiesFile Properties file <input type="button" value="Browse file"/>	Landmark.properties
<input checked="" type="checkbox"/> -rsample Rsample file <input type="button" value="Browse file"/>	gel-HM-019
<input checked="" type="checkbox"/> -sample Sample file <input type="button" value="Browse file"/>	gel-HM-071

View Menu



Landmark popup window

LMS edit Rsample: gel-HM-019, Sample: gel-HM-071 - 03/10/05 13:18:27

File View Edit Help

```
[1] x=187, y=226 gray=175  
Min R[E]=11 distNLM[22,22] to NLMs[D,D], distNNLM[41,38] to NNLMs[A,A]
```

Rsample: gel-HM-019

Sample: gel-HM-071

Flicker Sample: gel-HM-071

Rsample Delay:0.3Sec Sample Delay:0.3Sec > Sample:gel-HM-071

1X Add LM Delete LM Select LM Change LM Save image Done Cancel

Landmark XML DB in browser

```
<?xml version="1.0" ?>
- <LandmarkDatabase>
  <DatabaseName>HM</DatabaseName>
  <Date>08/10/04 19:36:18</Date>
  - <LandmarkSet>
    <Rsample>gel-HM-019</Rsample>
    <Sample>gel-HM-071</Sample>
    <lmNbr>1</lmNbr>
    <xRsample>207</xRsample>
    <yRsample>190</yRsample>
    <xSample>227</xSample>
    <ySample>176</ySample>
  </LandmarkSet>
  - <LandmarkSet>
    <Rsample>gel-HM-019</Rsample>
    <Sample>gel-HM-071</Sample>
    <lmNbr>2</lmNbr>
    <xRsample>176</xRsample>
    <yRsample>151</yRsample>
    <xSample>196</xSample>
    <ySample>140</ySample>
  </LandmarkSet>
  - <LandmarkSet>
    <Rsample>gel-HM-019</Rsample>
    <Sample>gel-HM-071</Sample>
    <lmNbr>3</lmNbr>
    <xRsample>158</xRsample>
    <yRsample>190</yRsample>
    <xSample>178</xSample>
    <ySample>173</ySample>
  </LandmarkSet>
</LandmarkDatabase>
```


Summary

- Landmark is a fully open-source landmark spot set editing program and is being developed at <http://open2dprot.sourceforge.net/>
- It uses a graphical user interface to let the user interactively assign corresponding spots.
- It saves the landmark set data for a pair of samples in an XML landmark database used by the other Open2Dprot pipeline modules.