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The screenshot displays the Spark software interface. At the top, the menu bar includes File, Edit, Project, View, Display, Run, Database, Window, and Help. Below the menu bar is a toolbar with icons for New, Open, Save, Add Prot, Constraints, Process, and Reference And Results. The main window is divided into several sections:

- Results Table:** A table with columns for Rank, Fav, Structure, BIF%, Score, and other metrics. The first four rows are highlighted in green. A right-click context menu is open over the table, showing options like Copy, Delete Selected Results, Export Selected Results, Mark as Favorite, Tag Results, View Log for Selected Results, and View Parent Structure for Selected Result.
- 3D Molecular Model:** A central 3D model of a molecule with various atoms and fields. A context menu is open over the model, showing options like Show Fields as Spheres, Tetrahedra, Cubes, Octahedra, Dodecahedra, and Icosahedra. Another context menu is open over an atom, showing options like Full screen (F11), Save Image As..., and Capture Scene (F4).
- Reference And Results Panel:** A panel on the left showing a list of references and results. A context menu is open over this panel, showing options like All References, Protein, All Results, and Selected Results.
- Toolbars and Menus:** Various toolbars and menus are visible, including the Reference And Results dropdown, the Fields menu, the Display menu, and the Atom menu.

At the bottom of the interface, there is a status bar with the text: "Left mouse rotate, right mouse translate, middle mouse/wheel/Alt-left scales. Shift-mouse z-clip."

| Shortcut | File operation |
|--------------|---------------------------|
| F1 | Show manual |
| Ctrl+Shift+S | Save project as |
| Ctrl+T | Open project notes editor |
| Ctrl+P | Print all results |
| Ctrl+Shift+P | Print selected results |
| Alt+F4 | Exit |

| Shortcut | Molecule operation |
|--------------|---------------------------------|
| Ctrl+F | Copy Favorites 3D |
| Ctrl+Alt+F | Copy Favorites 2D |
| Ctrl+V | Paste Reference from Clipboard |
| Ctrl+Space | Clear Favorites |
| Ctrl+Shift+V | Paste Excluded Volume Molecule |
| Ctrl+L | View log for selected molecules |

| Shortcut | View/ Display operation |
|--------------|-------------------------|
| Ctrl+Shift+R | Reset Display |
| Ctrl+Alt+R | Rock Display |
| Ctrl+Alt+S | Spin Display |
| F11 | Full Screen |
| F10 | Stereo |
| F4 | Capture scene |

| Right click to access menu |
|-----------------------------|
| Atom |
| Molecule in table |
| Toolbars and window headers |
| Column title |