

SLAP 2

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What happened since Victoria

- A draft of the document sent to DAL group chairs
- A list of questions and corrections was submitted
- Document was updated
- Next slides will introduce the modifications introduced in the standard

SLAP 2 interface (reminder)

| Resource type | Resource name | Required |
|-------------------|---------------|----------|
| DALI-sync | {lines} | yes |
| DALI-sync | {species} | no |
| DALI-examples | /examples | no |
| VOSI-availability | /availability | yes |
| VOSI-capabilities | /capabilities | yes |

- Returns a VOTable containing a list of lines
- Wavelength range is a mandatory query parameter for all implementations
- New parameters introduced :
 - MAXREC : maximum number of lines returned
 - RESPONSEFORMAT : default is text/xml, other possible value is application/x-votable+xml
 - LEVEL_ENERGY : range of energy for the levels of a line (upper or lower), in addition to already existing UPPER_LEVEL_ENERGY and LOWER_LEVEL_ENERGY

- Update of CHEMICAL_ELEMENT parameter definition in document :

"Atom can be specified exactly by symbol. Molecules can be specified by conventional molecular name (CO₂, CH₄ ...) which might no be unique."

- Recommended search pattern when looking for data related to a given species :

- query species resource to get list of species (with names)
- use found name to query the line resource

- Question : should we introduce a INCHIKEY parameter (or accept INCHIKEY value in CHEMICAL_ELEMENT) ?

- INCHIKEY is a hash of an inchi description of a Species

Example for Ozone molecule :

– Inchi : 1S/O3/c1-3-2

– Inchikey : CBENFWSGALASAD-UHFFFAOYSA-N

- advantage : more precise than molecular name to do a query

- but : data producer should provide inchikey for their species (not always trivial)
and probably update their Dbs (already done for VAMDC services)

Next step

- Presentation of Simple Spectral Lines Data Model in DM session
- Submission of document to WG