



The LXCat Project

Wouter Graef on behalf of the LXCat team

LT1.00008

What Is LXCat

- An open-access, web-based platform for storing, exchanging and manipulating data for modeling the electron and ion components of low-temperature, non-equilibrium plasmas.
- We support electron and ion scattering cross sections and rate coefficients, electron and ion swarm/transport parameters, ion-neutral interaction potentials, and optical oscillator strengths.
- Online tools allow for searching, graphical display, and downloading of data, and an online Boltzmann solver allows users to calculate electron transport and rate coefficients in arbitrary gas mixtures if "complete" sets of cross sections for the individual components are available in the databases.
- Visit us at www.lxcat.net

LXCat History

- 2009, Toulouse: Researchers at LAPLACE initiated the LXCat project
- 2010 GEC: Involvement by the community
- 2013, Eindhoven: Mirror server
- 2015: Founding of Association: Data for modeling plasmas
- 2017: Formation of technical support team
- 2019, Drake: Mirror server

LXCat Contributors

- Database contributors retain ownership and are responsible for the contents and maintenance of the individual databases
- New contributors are welcome and can request an account and receive instructions for setting up a password-protected database
- LXCat does not recommend data
- Contributors:

Leanne C. Pitchford, Luís L. Alves, Klaus Bartschat, Stephen F. Biagi, Marie-Clause Bordage, Igor Bray, Chris E. Brion, Michael J. Brunger, Stephen Buckman, Laurence Campbell, Alise Chachereau, Bhaskar Chaudhury, Zhi-Wen Cheng, Sanchita Chowdhury, Loucas G. Christophorou, Emile Carbone, Uwe Czarnetzki, Jan van Dijk, J. Dutton, Nikolay A. Dyatko, Christian M. Franck, Dmitry V. Fursa, Reetesh K. Gangwar, Vasco Guerra, Pascal Haefliger, Gerjan J.M. Hagelaar, Malte Hildebrandt, Andreas Hoesl, Felix Jacob, Yukikazu Itikawa, Indrek Jogi, Igor V. Kochetov, Vincenzo Laporta, Robert P. McEachran, W. Lowell Morgan, Anatoly P. Napartovich, Sergey Pancheshnyi, Yi-Kang Pu, Vincent Puech, Mohamed Rabie, Lalita Sharma, Rajesh Srivastava, Allan D. Stauffer, Jacob Stephens, Jonathan Tennyson, Tsanko Vaskov Tsankov, Jaime de Urquijo, Larry A. Viehland, Mark C. Zammit, Oleg Zatsarinny, Yan-Fei Wang, Yang Wang, Da-Ren Yu, Xi-Ming Zhu

LXCat Team

Outreach Team

Emile Carbone (Germany), Leanne Pitchford (France),
Klaus Bartschat (USA), Yi-Kang Pu (China)

Tech Team

Sergey Pancheshnyi (France, 2009-2017)
Wouter Graef, Diana Mihailova, Jan van Dijk (Netherlands),
Matthew Hopkins, Benjamin Yee, Jacob Stephens (USA)

LXCat Statistics

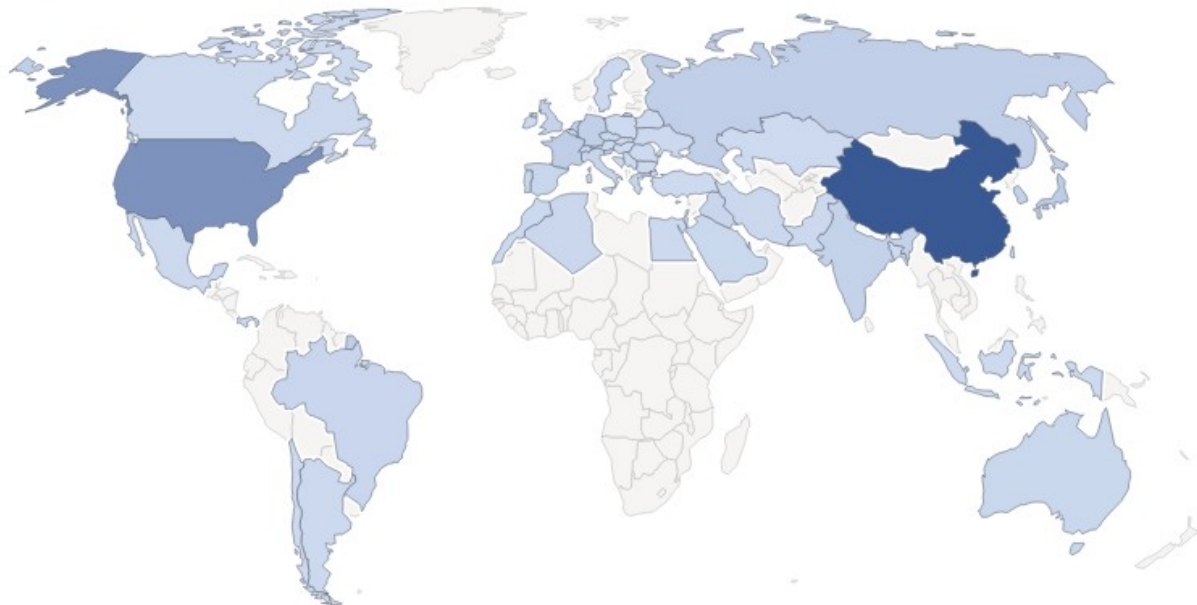
- 30 contributor databases:
 - Scattering cross sections: 26 databases | 103 x 517 species | 24.2k records
 - Differential scattering cross sections: 4 databases | 29 species | 517 records
 - Interaction potentials: 1 database | 91 x 7 species | 660 records
 - Oscillator strengths: 1 database | 65 species | 150 records
 - Swarm / transport data: 16 databases | 430 x 114 species | 193k records
 - Publications, notes and reports: 5 databases | 35 records
- 1280 citations

LXCat Statistics

- 3 servers: Eindhoven University of Technology & Drake University
- 150 daily visits



4,265 visits



How It Works

Data Retrieval

Select data type and database

STEP 1: SELECT DATA TYPE & CLICK NEXT

« PREV NEXT »

SCATTERING CROSS SECTIONS

DIFFERENTIAL SCATTERING CROSS SECTIONS

INTERACTION POTENTIALS

OSCILLATOR STRENGTHS

SWARM / TRANSPORT DATA

GLOBAL SPECIES FILTERING

ELECTRONS

IONS

STEP 2: SELECT DATABASES & CLICK NEXT

« PREV SORT BY NEXT »

SELECT ALL & FILTER TOOL no filter - +

Biagi (transcription of data from SF Biagi's Fortran code, Magboltz.)

Biagi-v7.1 (Magboltz version 7.1)

Bordage database

BSR (Quantum-mechanical calculations by D. Zatsarinny and K. Bartschat)

CCC database

Christophorou database

COP (Complex Optical Potential)

How It Works

Data Retrieval

Select species

Ground states

| | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|-----|-------|----------|------|------|-----|----|----|------|------|------|-------|------|------|------|------|-----|--------|------|----|-----|-----|-----|----|-----|-----|-----|------|-----|----|-----|-------|-----|----|----------------|-----|----|-----|-----|-----|-----|-----|
| Ar | BCl | BCl2 | BCl3 | BF3 | Be | C | C2 | C2H2 | C2H4 | C2H6 | C2OH6 | C3 | C3H4 | C3H6 | C3H8 | C3N | CCl2F2 | CCl4 | CF | CF2 | CF3 | CF4 | CH | CH2 | CH3 | CH4 | CHF3 | CNH | CO | CO2 | CONH3 | CO2 | CS | CaF | Cl2 | | | | | | |
| Cu | D2 | DT | F | F2 | F2O | H | H2 | H2O | H2S | H4C | HBr | HCHO | HCN | HCP | HCl | HD | HT | He | Hg | Kr | Mg | N | N2 | N2O | NF3 | NH3 | NO | NO2 | Na | Ne | O | O2 | O3 | O ⁻ | PH3 | SF | SF2 | SF3 | SF4 | SF5 | SF6 |
| SO2 | Si2H6 | Si(CH3)4 | SiF2 | SiH4 | SiO | T2 | Xe | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |

State-specific and gas mixtures

| | | | | | | | | | | | | |
|-----------------------|----------------------------|----------------------------|----------------------------|----------------------------|----------------------------|----------------------------|----------------------------|----------------------------|----------------------------|----------------------------|----------------------------|----------------------|
| Ar* | Ar(3d ³ [3/2]2) | Ar(3d ⁵ [5/2]2) | Ar(3d ⁵ [5/2]3) | Ar(3d ¹ [1/2]0) | Ar(3d ¹ [1/2]1) | Ar(3d ³ [3/2]1) | Ar(3d ³ [3/2]2) | Ar(3d ⁵ [5/2]2) | Ar(3d ⁵ [5/2]3) | Ar(3d ⁷ [7/2]3) | Ar(3d ⁷ [7/2]4) | Ar(3p5 4p J = 0 2p5) |
| Ar(3p5 4p J = 1 2p10) | Ar(3p5 4p J = 1 2p2) | Ar(3p5 4p J = 1 2p4) | Ar(3p5 4p J = 1 2p7) | Ar(3p5 4p J = 2 2p3) | Ar(3p5 4p J = 2 2p6) | Ar(3p5 4p J = 2 2p8) | Ar(3p5 4p J = 3 2p9) | Ar(3p5 4s J = 0 1s3) | | | | |

Select processes

Biagi (transcription of data from SF Biagi's Fortran code, Magboltz.)

Data Group [Ar]: Transcribed from S.F. Biagi's Fortran Magboltz version 8.97 (Sept 2011).
Data are based in part on the calculations of Zatsariny and Bartschat. See BSR database on this site.


e / Ar

- Elastic E + Ar → E + Ar (m/M = 0.0000136, complete set) | elastic momentum transfer from Magboltz 8.97 Sept 2011. Note that the energy data. Updated: 3 October 2014.
- Excitation E + Ar → E + Ar(1S5) (E = 11.548 eV, complete set) | from Magboltz 8.97 Sept 2011. Updated: 20 October 2011.
- Excitation E + Ar → E + Ar(1S4) (E = 11.624 eV, complete set) | from Magboltz 8.97 Sept 2011. Updated: 20 October 2011.
- Excitation E + Ar → E + Ar(1S3) (E = 11.723 eV, complete set) | from Magboltz 8.97 Sept 2011. Updated: 20 October 2011.
- Excitation E + Ar → E + Ar(1S2) (E = 11.828 eV, complete set) | from Magboltz 8.97 Sept 2011. Updated: 20 October 2011.
- Excitation E + Ar → E + Ar(2P10) (E = 12.907 eV, complete set) | from Magboltz 8.97 Sept 2011. Updated: 20 October 2011.
- Excitation E + Ar → E + Ar(2P9) (E = 13.076 eV, complete set) | from Magboltz 8.97 Sept 2011. Updated: 20 October 2011.
- Excitation E + Ar → E + Ar(2P8) (E = 13.095 eV, complete set) | from Magboltz 8.97 Sept 2011. Updated: 20 October 2011.
- Excitation E + Ar → E + Ar(2P7) (E = 13.153 eV, complete set) | from Magboltz 8.97 Sept 2011. Updated: 20 October 2011.
- Excitation E + Ar → E + Ar(2P6) (E = 13.172 eV, complete set) | from Magboltz 8.97 Sept 2011. Updated: 20 October 2011.
- Excitation E + Ar → E + Ar(2P5) (E = 13.272 eV, complete set) | from Magboltz 8.97 Sept 2011. Updated: 20 October 2011.

How It Works

Online BOLSIG+ Calculation

Select from available complete sets

| | CH4 | N2 |
|--------------|---|---|
| Hayashi | <input type="checkbox"/> | |
| IST-Lisbon | <input type="checkbox"/> | <input type="checkbox"/> |
| Morgan | <input checked="" type="checkbox"/> | <input checked="" type="checkbox"/> |
| TRINITY | <input type="checkbox"/> | <input type="checkbox"/> |
| Biagi | | <input type="checkbox"/> |
| Phelps | | <input type="checkbox"/> |
| SIGLO | | <input type="checkbox"/> |
| Publications |  |  |

Select settings

BASIC PARAMETERS

non-Maxwellian | Maxwellian EEDF

E/N = - Td

of points =

▾


T_{gas} = K


Include super-elastic collisions


SPECIES MOLE FRACTIONS

CH4 =


N2 =

AC FIELD 

ADVANCED OPTIONS 

BROWSE RESULTS 

Electron cross sections (new window)

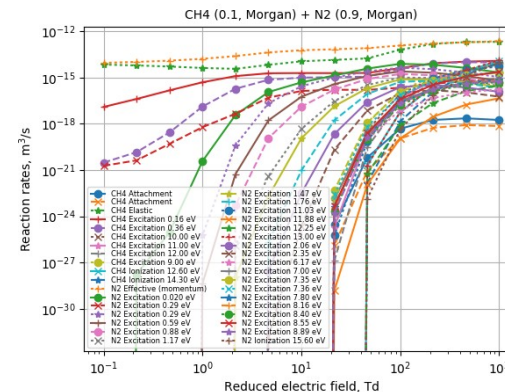
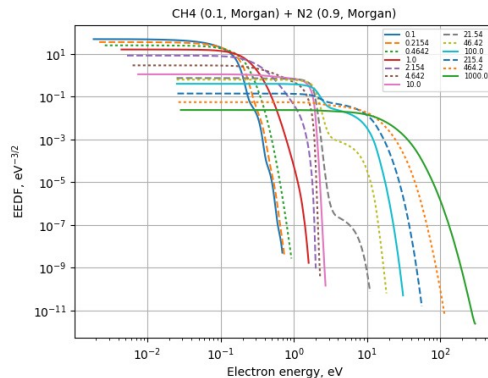
DOWNLOAD 

Cross section input file

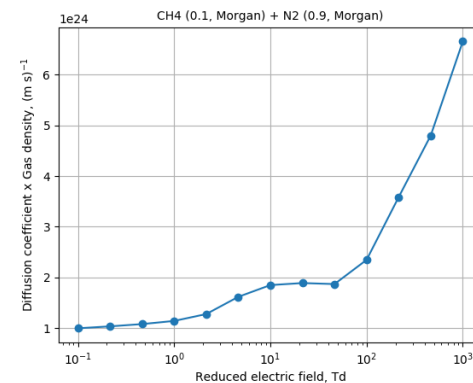
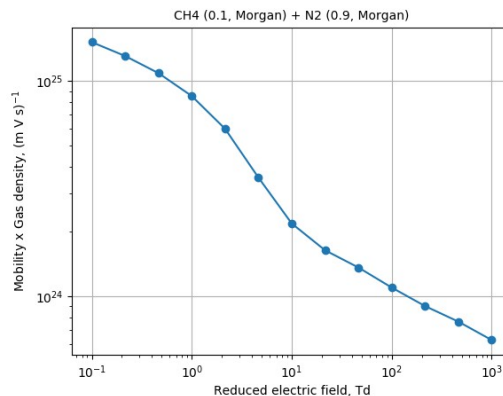
How It Works

Online BOLSIG+ Calculation

EEDF's &
Rate coefficients



Mobility &
Diffusion coefficient



Results can be compared to database

Referencing

Users of LXCat **MUST** properly reference used material

Website:

Terms of use

Users acknowledge understanding that LXCat is a community-based project with open-access databases being freely provided by individual contributors.

Proper referencing of material retrieved from this site is essential for the survival of the project.

Users further accept that the databases on this site remain property of their respective contributors and are not to be distributed by third parties or used for commercial purposes. All questions regarding copyright should be addressed to the [LXCat team](#).

Output files:

HOW TO REFERENCE: Fortran program, MAGBOLTZ, S.F. Biagi. The version number of MAGBOLTZ from which the cross sections were transcribed is provided in the comment for each species. This version number must be included in the reference.

Resources

- Online quick start guides
- Discussion board hosted on Google groups
- Links to various software related to LXCat and software by and for the LTP community:

BOLSIG+, ZDPlasKin, EEDF, BOLOS, PumpKin, METHES, MultiBolt, LoKI-B, THERMCAT, ELEM, MOBION

LXCat In The Future

- Species naming scheme, allowing for proper categorization of data
- In addition to current output format, output in XML/JSON with schemas
- API for direct data access
- Additional online services, such as cross section calculator for interaction potentials

Get Involved

- Become a contributor
- Share your ideas and wishes
- Help future development



<https://www.lxcat.net>

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