

Integrative Structure Validation Report ?

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The following software was used in the production of this report:

Integrative Modeling Validation Version 2.0

Python-IHM Version 1.8

MolProbity Version 4.5.2

PDB ID	9A4V
PDB-Dev ID	PDBDEV_00000252
Structure Title	Integrative model of PT1-ODO2 by crosslinking MS and deep learning
Structure Authors	Kolja Stahl; Oliver Brock; Juri Rappsilber
Deposited on	2024-01-23

This is a PDB-IHM IM Structure Validation Report for a publicly released PDB-IHM entry.

We welcome your comments at helpdesk@pdb-ihm.org

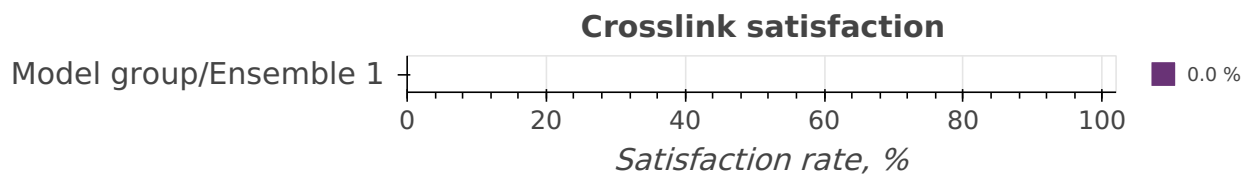
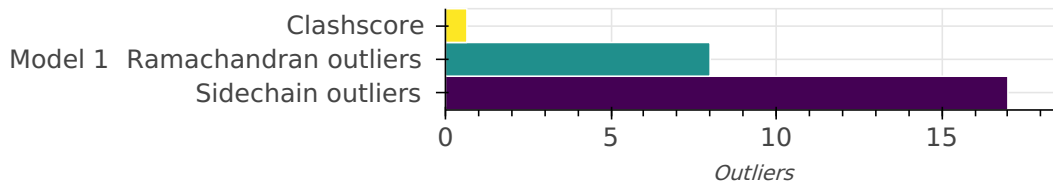
A user guide is available at https://pdb-ihm.org/validation_help.html with specific help available everywhere you see the ? symbol.

List of references used to build this report is available [here](#).

Overall quality ?

This validation report contains model quality assessments for all structures, data quality and fit to model assessments for SAS and crosslinking-MS datasets. Data quality and fit to model assessments for other datasets and model uncertainty are under development. Number of plots is limited to 256.

Model Quality: MolProbity Analysis



Ensemble information ?

This entry consists of 0 distinct ensemble(s).

Summary ?

This entry consists of 1 model(s). A total of 1 datasets were used to build this entry.

Representation ?

This entry has 1 representation(s).

ID	Model(s)	Entity ID	Molecule name	Chain(s) [auth]	Total residues	Rigid segments	Flexible segments	Model coverage/ Starting model coverage (%)	Scale
1	1	1	PT1_BACSU	A	570	-	1-570	100.00 / 0.00	Atomic
		2	ODO2_BACSU	B	417	-	1-417	100.00 / 0.00	Atomic

Datasets used for modeling ?

There is 1 unique dataset used to build the models in this entry.

ID	Dataset type	Database name	Data access code
1	Crosslinking-MS data	PRIDE	PXD035508

Methodology and software ?

This entry is a result of 1 distinct protocol(s).

Step number	Protocol ID	Method name	Method type	Method description	Number of computed models	Multi state modeling	Multi scale modeling
1	1	AlphaLink2	AlphaLink2	None	1	False	False

There is 1 software package reported in this entry.

ID	Software name	Software version	Software classification	Software location
1	AlphaLink2	1.00	model building	https://github.com/Rappsilber-Laboratory/AlphaLink2

Data quality ?

Crosslinking-MS

At the moment, data validation is only available for crosslinking-MS data deposited as a fully [compliant](#) dataset in the [PRIDE Crosslinking](#) database. Correspondence between crosslinking-MS and entry entities is established using [pyHMMER](#). Only residue pairs that passed the reported threshold are used for the analysis. The values in the report have to be interpreted in the context of the experiment (i.e. only a minor fraction of in-situ or in-vivo dataset can be used for modeling).

Crosslinking-MS dataset is not available in the [PRIDE Crosslinking](#) database.

Model quality ?

For models with atomic structures, MolProbity analysis is performed. For models with coarse-grained or multi-scale structures, excluded volume analysis is performed.

Standard geometry: bond outliers ?

There are no bond length outliers.

Standard geometry: angle outliers ?

There are 65 bond angle outliers in this entry (0.62% of 10477 assessed bonds). A summary is provided below.

Chain	Res	Type	Atoms	Z	Observed (Å)	Ideal (Å)	Model ID (Worst)	Models (Total)
B	101	LYS	C-N-CA	7.62	135.41	121.70	1	1
B	91	THR	C-N-CA	7.38	134.99	121.70	1	1
B	95	GLU	C-N-CA	6.58	133.54	121.70	1	1
B	98	LYS	C-N-CA	6.30	133.04	121.70	1	1
B	90	LYS	C-N-CA	6.27	132.98	121.70	1	1
B	112	GLU	C-N-CA	6.18	132.83	121.70	1	1
A	129	ASP	CA-CB-CG	5.82	118.42	112.60	1	1
A	523	PHE	CA-CB-CG	5.74	119.54	113.80	1	1
B	91	THR	O-C-N	5.60	114.04	123.00	1	1
B	91	THR	CA-C-N	5.56	127.32	116.20	1	1
B	412	GLN	OE1-CD-NE2	5.31	117.29	122.60	1	1
A	96	ASP	CA-CB-CG	5.24	117.84	112.60	1	1
B	392	ASP	CA-CB-CG	5.17	117.77	112.60	1	1
B	93	SER	C-N-CA	5.05	130.80	121.70	1	1
B	107	ALA	C-N-CA	5.03	130.75	121.70	1	1
B	268	GLN	OE1-CD-NE2	4.95	117.65	122.60	1	1
A	2	GLN	OE1-CD-NE2	4.85	117.75	122.60	1	1
B	102	GLN	OE1-CD-NE2	4.77	117.83	122.60	1	1
B	24	GLN	OE1-CD-NE2	4.77	117.83	122.60	1	1
B	297	ASP	CA-CB-CG	4.76	117.36	112.60	1	1
B	91	THR	CA-CB-OG1	4.72	116.69	109.60	1	1
B	197	GLN	OE1-CD-NE2	4.70	117.90	122.60	1	1
B	103	ALA	CA-C-N	4.69	125.58	116.20	1	1
B	56	GLN	OE1-CD-NE2	4.66	117.94	122.60	1	1
B	103	ALA	N-CA-CB	4.66	103.42	110.40	1	1
B	365	ARG	NE-CZ-NH1	4.61	126.11	121.50	1	1
B	103	ALA	O-C-N	4.60	115.64	123.00	1	1
B	353	GLN	OE1-CD-NE2	4.57	118.03	122.60	1	1
B	67	GLN	OE1-CD-NE2	4.56	118.04	122.60	1	1
A	249	GLN	OE1-CD-NE2	4.53	118.07	122.60	1	1
B	103	ALA	C-CA-CB	4.52	117.28	110.50	1	1
B	319	ASN	OD1-CG-ND2	4.52	118.08	122.60	1	1
B	93	SER	CA-C-N	4.51	125.23	116.20	1	1
B	238	GLN	OE1-CD-NE2	4.50	118.10	122.60	1	1

Chain	Res	Type	Atoms	Z	Observed (Å)	Ideal (Å)	Model ID (Worst)	Models (Total)
B	328	GLN	OE1-CD-NE2	4.50	118.10	122.60	1	1
B	207	GLN	OE1-CD-NE2	4.49	118.11	122.60	1	1
B	208	GLN	OE1-CD-NE2	4.45	118.15	122.60	1	1
B	95	GLU	CA-C-N	4.45	125.10	116.20	1	1
A	323	ARG	NH1-CZ-NH2	4.44	113.53	119.30	1	1
A	385	ASN	CA-CB-CG	4.44	117.04	112.60	1	1
A	346	GLN	OE1-CD-NE2	4.42	118.18	122.60	1	1
B	177	GLN	OE1-CD-NE2	4.41	118.19	122.60	1	1
B	103	ALA	C-N-CA	4.40	129.61	121.70	1	1
B	93	SER	O-C-N	4.37	116.02	123.00	1	1
A	543	GLN	OE1-CD-NE2	4.34	118.26	122.60	1	1
A	67	GLN	OE1-CD-NE2	4.34	118.26	122.60	1	1
B	270	ASP	CA-CB-CG	4.31	116.91	112.60	1	1
B	104	GLU	CA-C-N	4.30	123.36	116.90	1	1
A	60	HIS	CB-CG-CD2	4.28	125.64	131.20	1	1
A	455	ASP	CA-CB-CG	4.24	108.36	112.60	1	1
B	175	GLN	OE1-CD-NE2	4.23	118.37	122.60	1	1
B	170	GLN	OE1-CD-NE2	4.22	118.38	122.60	1	1
B	172	GLN	OE1-CD-NE2	4.21	118.39	122.60	1	1
B	181	GLN	OE1-CD-NE2	4.20	118.40	122.60	1	1
B	31	GLN	OE1-CD-NE2	4.20	118.40	122.60	1	1
A	455	ASP	C-CA-CB	4.19	118.06	110.10	1	1
A	140	HIS	CB-CG-CD2	4.16	125.79	131.20	1	1
B	141	GLN	OE1-CD-NE2	4.16	118.44	122.60	1	1
A	213	GLN	OE1-CD-NE2	4.14	118.46	122.60	1	1
B	173	GLN	OE1-CD-NE2	4.09	118.51	122.60	1	1
B	169	GLN	OE1-CD-NE2	4.09	118.51	122.60	1	1
B	112	GLU	O-C-N	4.08	116.48	123.00	1	1
A	447	ASP	CA-CB-CG	4.03	108.57	112.60	1	1
B	180	GLN	OE1-CD-NE2	4.01	118.59	122.60	1	1
A	267	HIS	CB-CG-CD2	4.00	126.00	131.20	1	1

Too-close contacts ?

The following all-atom clashscore is based on a MolProbity analysis. All-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The table below contains clashscores for all atomic models in this entry.

Model ID	Clash score	Number of clashes
1	0.65	10

There are 10 clashes. The table below contains the detailed list of all clashes based on a MolProbity analysis. Bad clashes are ≥ 0.4 Angstrom.

Atom 1	Atom 2	Clash(Å)	Model ID (Worst)	Models (Total)
B:107:ALA:HB1	B:108:GLN:HB2	0.61	1	1
A:253:TRP:CD1	A:535:THR:HG1	0.52	1	1
B:218:GLU:CD	B:365:ARG:HH22	0.49	1	1
A:318:LYS:HE2	A:383:TYR:CE1	0.48	1	1
B:365:ARG:HH21	B:379:MET:HG3	0.46	1	1
A:1:MET:HA	A:229:ILE:O	0.46	1	1
B:107:ALA:HB1	B:108:GLN:CB	0.43	1	1
A:5:LYS:HE2	A:226:ASP:OD1	0.42	1	1
A:30:LYS:HE3	A:98:VAL:HG11	0.41	1	1
A:400:LYS:HE2	A:445:GLU:OE1	0.41	1	1

Torsion angles: Protein backbone ?

In the following table, Ramachandran outliers are listed. The Analysed column shows the number of residues for which the backbone conformation was analysed.

Model ID	Analysed	Favored	Allowed	Outliers
1	983	926	49	8

There are 8 unique backbone outliers. Detailed list of outliers are tabulated below.

Chain	Res	Type	Models (Total)
A	206	LYS	1
B	99	GLU	1
B	104	GLU	1
B	105	PRO	1
B	164	SER	1
B	172	GLN	1
B	178	LYS	1
B	182	SER	1

Torsion angles : Protein sidechains ?

In the following table, sidechain rotameric outliers are listed. The Analysed column shows the number of residues for which the sidechain conformation was analysed.

Model ID	Analysed	Favored	Allowed	Outliers
1	834	799	18	17

There are 17 unique sidechain outliers. Detailed list of outliers are tabulated below.

Chain	Res	Type	Models (Total)
A	78	LEU	1
A	150	SER	1
A	190	SER	1
A	335	ASP	1
A	336	ILE	1
A	345	LEU	1
A	440	ASP	1
A	526	SER	1
A	556	SER	1
B	61	ASP	1
B	88	THR	1
B	91	THR	1
B	110	VAL	1
B	113	GLU	1
B	122	THR	1
B	155	ASP	1
B	272	LEU	1

Fit of model to data used for modeling ?

Fit of model(s) to crosslinking-MS data

Restraint types

Restraint types are summarized in the table below. Restraints assigned "*by-residue*" are interpreted as between CA atoms. Restraints between coarse-grained beads are indicated as "*coarse-grained*".

Restraint group represents a set of crosslinking restraints applied collectively in the modeling.

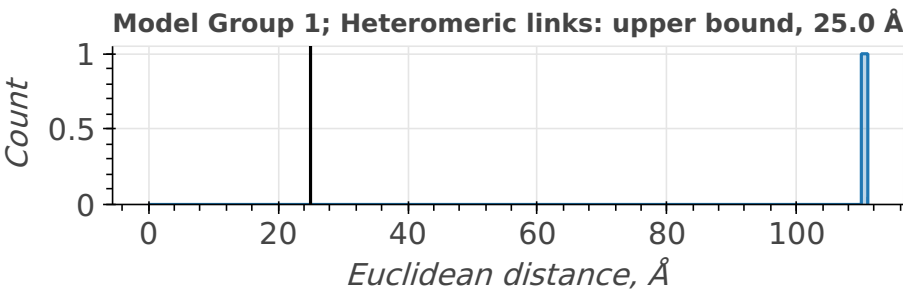
There are 1 crosslinking restraints combined in 1 restraint groups.

Linker	Residue 1	Atom 1	Residue 2	Atom 2	Restraint type	Distance, Å	Count
SDA	LYS	CA	LYS	CA	upper bound	25.0	1

Distograms of individual restraints

Restraints with identical thresholds are grouped into one plot. Only the best distance per restraint per

model group/ensemble is plotted. Inter- and intramolecular (including self-links) restraints are also grouped into one plot. Distance for a restraint between coarse-grained beads is calculated as a minimal distance between shells; if beads intersect, the distance will be reported as 0.0. A bead with the highest available resolution for a given residue is used for the assessment.



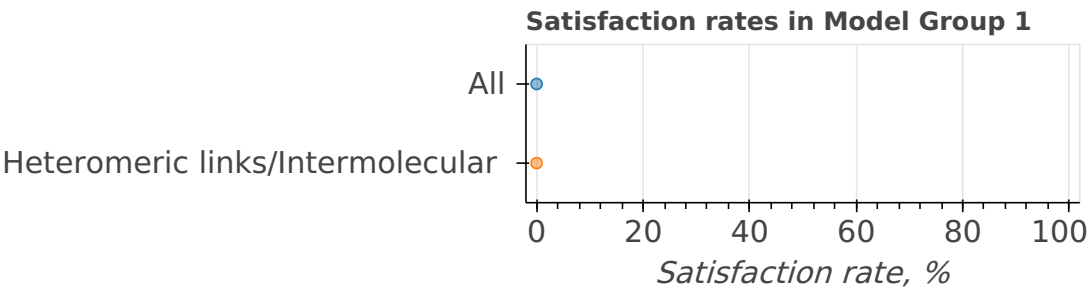
Satisfaction of restraints

Satisfaction of restraints is calculated on a [restraint group](#) (a set of crosslinking restraints applied collectively in the modeling) level. Satisfaction of a restraint group depends on satisfaction of individual restraints in the group and the conditionality (all/any). A restraint group is considered satisfied, if the condition was met in at least one model of the model group/ensemble. The number of measured restraints can be smaller than the total number of restraint groups if crosslinks involve non-modeled residues. Only deposited models are used for validation right now.

State group	State	Model group	# of Deposited models/Total	Restraint group type	Satisfied (%)	Violated (%)	Count (Total=1)
1	1	1	1/1	All	0.00	100.00	1
				Heteromeric links/ Intermolecular	0.00	100.00	1

Per-model satisfaction rates in ensembles

Every point represents one model in a model group/ensemble. Where possible, boxplots with quartile marks are also plotted.



Fit of model to data used for validation ?

Validation for this section is under development.

Acknowledgments

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