Integrative Structure Validation Report

September 10, 2024 - 10:14 PM PDT

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

Python-IHM Version 1.3
MolProbity Version 4.5.2
Integrative Modeling Validation Version 1.2

PDB ID	9A4L
PDB-Dev ID	PDBDEV_00000242
Structure Title	Integrative model of ALF-ACNA by crosslinking MS and deep learning
Structure Authors	Kolja Stahl; Oliver Brock; Juri Rappsilber

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

We welcome your comments at pdb-dev@mail.wwpdb.org

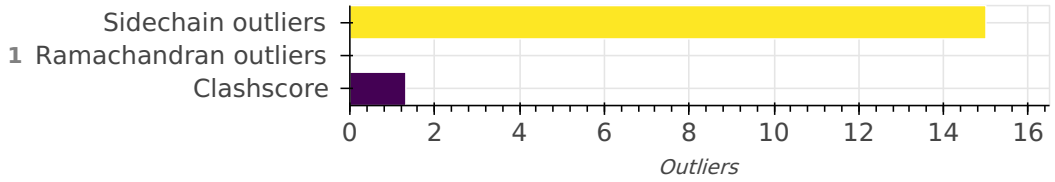
A user guide is available at https://pdb-dev.wwpdb.org/validation_help.html with specific help available everywhere you see the ? symbol.

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 assessment for SAS datasets and fit to model assessments for SAS 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 unique models, with 2 subunits in each model. A total of 1 datasets or restraints were used to build this entry. Each model is represented by 0 rigid bodies and 2 flexible or non-rigid units.

Entry composition

There is 1 unique type of models in this entry. This model is titled None/None.

Model ID	Subunit number	Subunit ID	Subunit name	Chain ID	Chain ID [auth]	Total residues
1	1	1	ALF_BACSU	A	A	285
1	2	2	ACNA_BACSU	B	B	909

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

Representation

This entry has only one representation and includes 0 rigid bodies and 2 flexible units.

Chain ID	Rigid bodies	Non-rigid segments

Chain ID	Rigid bodies	Non-rigid segments
A	-	1-285
B	-	1-909

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	1.0	model building	https://github.com/Rappsilber-Laboratory/AlphaLink2

Data quality

Crosslinking-MS

Validation for this section is under development.

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 9098 bond outliers in this entry. A summary is provided below, and a detailed list of outliers can be found here.

Bond type	Observed distance (Å)	Ideal distance (Å)	Number of outliers
CD2--HD22	1.09	0.97	97

Bond type	Observed distance (Å)	Ideal distance (Å)	Number of outliers
CE--HE2	1.09	0.97	108
CG1--HG11	1.09	0.97	111
CB--HB2	1.09	0.97	831
CD--HD2	1.09	0.97	180
CG--HG3	1.09	0.97	337
CB--HB3	1.09	0.97	831
CD1--HD13	1.09	0.97	161
CG--HG2	1.09	0.97	337
CG2--HG21	1.09	0.97	252
CE--HE1	1.09	0.97	27
CG2--HG22	1.09	0.97	252
CD2--HD23	1.09	0.97	97
CA--HA	1.09	0.97	1083
NZ--HZ3	1.01	0.89	81
CB--HB	1.09	0.97	252
CG1--HG13	1.09	0.97	175
CD1--HD11	1.09	0.97	161
NZ--HZ2	1.01	0.89	81
OG1--HG1	0.96	0.84	77
CG2--HG23	1.09	0.97	252
NZ--HZ1	1.01	0.89	81
CA--HA2	1.09	0.97	111
CB--HB1	1.09	0.97	99

Bond type	Observed distance (Å)	Ideal distance (Å)	Number of outliers
N--H3	1.01	0.89	2
CD2--HD21	1.09	0.97	97
CG1--HG12	1.09	0.97	175
OH--HH	0.96	0.84	38
CD--HD3	1.09	0.97	180
CD1--HD12	1.09	0.97	161
CG--HG	1.09	0.97	97
CA--HA3	1.09	0.97	111
CE--HE3	1.09	0.97	108
OG--HG	0.96	0.84	59
N--H2	1.01	0.89	2
N--H1	1.01	0.89	2
SG--HG	1.33	1.20	1
SG--HG	1.34	1.20	6
N--H	1.01	0.86	1131
NH2--HH21	1.01	0.86	38
CD1--HD1	1.08	0.93	91
ND2--HD21	1.01	0.86	57
CD2--HD2	1.08	0.93	104
NH1--HH12	1.01	0.86	38
NH1--HH11	1.01	0.86	38
CE2--HE2	1.08	0.93	83
NH2--HH22	1.01	0.86	38
NE2--HE21	1.01	0.86	31

Bond type	Observed distance (Å)	Ideal distance (Å)	Number of outliers
CE1--HE1	1.08	0.93	104
CE3--HE3	1.08	0.93	8
NE--HE	1.01	0.86	38
ND2--HD22	1.01	0.86	57
ND1--HD1	1.01	0.86	15
CZ3--HZ3	1.08	0.93	8
CZ--HZ	1.08	0.93	45
NE1--HE1	1.01	0.86	8
NE2--HE2	1.01	0.86	6
NE2--HE22	1.01	0.86	31
CZ2--HZ2	1.08	0.93	8
CH2--HH2	1.08	0.93	8

Standard geometry: angle outliers

There are 26 angle outliers in this entry. A summary is provided below, and a detailed list of outliers can be found here.

Angle type	Observed angle (°)	Ideal angle (°)	Number of outliers
OD1-CG-ND2	122.60	117.23	1
CA-CB-CG	112.60	117.94	1
CB-CG-CD2	131.20	124.29	1
OE1-CD-NE2	122.60	117.53	1
OE1-CD-NE2	122.60	117.58	1
OE1-CD-NE2	122.60	117.59	1
OE1-CD-NE2	122.60	117.80	1
CB-CG-CD2	131.20	125.23	1
OE1-CD-NE2	122.60	118.06	1

Angle type	Observed angle (°)	Ideal angle (°)	Number of outliers
OE1-CD-NE2	122.60	118.16	1
NE-CZ-NH2	119.20	123.16	1
OE1-CD-NE2	122.60	118.20	1
CB-CG-CD2	131.20	125.54	1
OE1-CD-NE2	122.60	118.29	1
CA-CB-CG	112.60	116.88	1
CB-CG-CD2	131.20	125.66	1
OE1-CD-NE2	122.60	118.34	1
CB-CG-CD2	131.20	125.67	1
CB-CG-CD2	131.20	125.78	1
OE1-CD-NE2	122.60	118.43	1
CA-CB-CG	112.60	108.48	1
CB-CG-CD2	131.20	125.87	1
C-CA-CB	110.10	117.89	1
CA-CB-CG	112.60	116.69	1
OE1-CD-NE2	122.60	118.55	1
OD1-CG-ND2	122.60	118.57	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 the models in this entry.

Model ID	Clash score	Number of clashes
1	1.32	24

All 24 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Model ID	Atom-1	Atom-2	Clash overlap (Å)

Model ID	Atom-1	Atom-2	Clash overlap (Å)
1	B:397:MET:HE1	B:563:TYR:CE1	0.673
1	B:304:LEU:HD22	B:333:TYR:CG	0.593
1	B:314:ALA:HA	B:317:TYR:CZ	0.586
1	B:491:VAL:HA	B:494:TYR:CE1	0.558
1	B:193:GLU:C	B:339:ARG:HH22	0.555
1	A:257:PRO:HA	A:260:TYR:CZ	0.531
1	B:679:ARG:HH21	B:758:TRP:CG	0.525
1	B:304:LEU:HD22	B:333:TYR:CD1	0.499
1	B:215:LEU:HD21	B:225:ILE:HG21	0.492
1	B:264:LYS:HZ3	B:301:GLU:CD	0.472
1	B:238:ILE:CD1	B:799:LYS:HE3	0.470
1	B:781:VAL:HG23	B:808:THR:HG23	0.469
1	A:257:PRO:HA	A:260:TYR:CE1	0.465
1	B:154:ASP:HA	B:635:ARG:HH12	0.447
1	B:117:VAL:HG21	B:906:MET:HE3	0.435
1	B:679:ARG:HH21	B:758:TRP:CD1	0.426
1	B:117:VAL:CG2	B:906:MET:HE3	0.418
1	B:104:VAL:HG12	B:166:PHE:CZ	0.408
1	B:314:ALA:HA	B:317:TYR:CE2	0.408
1	B:683:LYS:HE3	B:843:LEU:O	0.406
1	B:215:LEU:HD23	B:232:GLY:HA3	0.403
1	B:786:LYS:HE3	B:812:GLU:OE2	0.403
1	B:104:VAL:HG11	B:170:ALA:HB2	0.402

Model ID	Atom-1	Atom-2	Clash overlap (Å)
1	B:732:MET:HB3	B:796:TRP:CE3	0.400

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	Analyzed	Favored	Allowed	Outliers
1	1190	1157	33	0

Detailed list of outliers are tabulated below.

Torsion angles: Protein sidechains

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

Model ID	Analyzed	Favored	Allowed	Outliers
1	984	951	18	15

Detailed list of outliers are tabulated below.

Model ID	Chain	Residue ID	Residue type
1	A	1	MET
1	A	164	LEU
1	B	7	THR
1	B	52	LEU
1	B	202	ILE
1	B	299	ILE
1	B	304	LEU
1	B	371	VAL
1	B	405	LEU
1	B	430	LEU
1	B	435	THR
1	B	612	LEU

Model ID	Chain	Residue ID	Residue type
1	B	643	ASP
1	B	790	MET
1	B	890	ASP

Fit of model to data used for modeling

Crosslinking-MS

Validation for this section is under development.

Fit of model to data used for validation

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Acknowledgements

Development of integrative model validation metrics, implementation of a model validation pipeline, and creation of a validation report for integrative structures, are funded by NSF ABI awards (DBI-1756248, DBI-2112966, DBI-2112967, DBI-2112968, and DBI-1756250). The PDB-Dev team and members of Sali labcontributed model validation metrics and software packages.

Implementation of validation methods for SAS data and SAS-based models are funded by RCSB PDB (grant number DBI-1832184). Dr. Stephen Burley, Dr. John Westbrook, and Dr. Jasmine Young from RCSB PDB, Dr. Jill Trewhella, Dr. Dina Schneidman, and members of the SASBDB repository are acknowledged for their advice and support in implementing SAS validation methods.

Members of the wwPDB Integrative/Hybrid Methods Task Force provided recommendations and community support for the project.