

Integrative Structure Validation Report ?

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

Python-IHM Version 1.3

MolProbit Version 4.5.2

Integrative Modeling Validation Version 1.2

PDB ID	9A4J
PDB-Dev ID	PDBDEV_00000240
Structure Title	Integrative model of ACNA-IDH by crosslinking MS and deep learning
Structure Authors	Kolja Stahl; Oliver Brock; Juri Rappaport

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

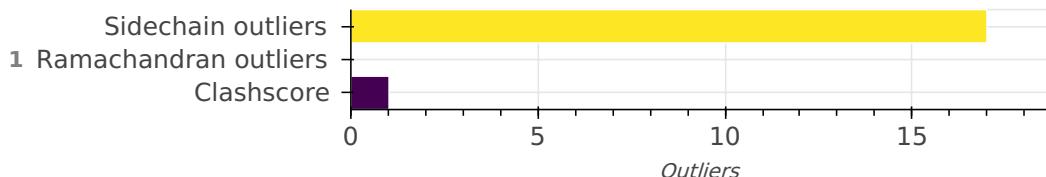
A user guide is available at https://pdb-dev.wwpdb.org/validation_help.html with specific help available everywhere you see the ? symbol.

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: MolProbit 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	ACNA_BACSU	A	A	909
1	2	2	IDH_BACSU	B	B	423

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	-	1-909
B	-	1-423

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 10237 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
CB--HB2	1.09	0.97	928

Bond type	Observed distance (Å)	Ideal distance (Å)	Number of outliers
CA--HA	1.09	0.97	1212
CG2--HG22	1.09	0.97	284
CB--HB3	1.09	0.97	928
CG--HG3	1.09	0.97	372
CD1--HD12	1.09	0.97	190
CG2--HG21	1.09	0.97	284
CD1--HD13	1.09	0.97	190
CG1--HG11	1.09	0.97	117
CA--HA2	1.09	0.97	120
CG--HG2	1.09	0.97	372
OG--HG	0.96	0.84	61
CG1--HG12	1.09	0.97	201
CD--HD3	1.09	0.97	204
CD--HD2	1.09	0.97	204
CG2--HG23	1.09	0.97	284
NZ--HZ2	1.01	0.89	93
CE--HE2	1.09	0.97	118
CA--HA3	1.09	0.97	120
CD2--HD22	1.09	0.97	106
CG1--HG13	1.09	0.97	201
CD2--HD23	1.09	0.97	106
CD2--HD21	1.09	0.97	106
CD1--HD11	1.09	0.97	190

Bond type	Observed distance (Å)	Ideal distance (Å)	Number of outliers
NZ--HZ1	1.01	0.89	93
CB--HB	1.09	0.97	284
CB--HB1	1.09	0.97	111
CG--HG	1.09	0.97	106
CE--HE3	1.09	0.97	118
OG1--HG1	0.96	0.84	83
NZ--HZ3	1.01	0.89	93
N--H2	1.01	0.89	2
OH--HH	0.96	0.84	46
CE--HE1	1.09	0.97	25
N--H3	1.01	0.89	2
N--H1	1.01	0.89	2
SG--HG	1.34	1.20	6
CE2--HE2	1.08	0.93	97
NE--HE	1.01	0.86	46
ND2--HD21	1.01	0.86	69
NE2--HE21	1.01	0.86	34
N--H	1.01	0.86	1265
CE3--HE3	1.08	0.93	14
NE2--HE22	1.01	0.86	34
ND2--HD22	1.01	0.86	69
CD2--HD2	1.08	0.93	114
NH1--HH11	1.01	0.86	46
CE1--HE1	1.08	0.93	114

Bond type	Observed distance (Å)	Ideal distance (Å)	Number of outliers
NE1--HE1	1.01	0.86	14
NH1--HH12	1.01	0.86	46
CZ--HZ	1.08	0.93	51
CD1--HD1	1.08	0.93	111
NH2--HH22	1.01	0.86	46
CZ3--HZ3	1.08	0.93	14
NH2--HH21	1.01	0.86	46
CZ2--HZ2	1.08	0.93	14
ND1--HD1	1.01	0.86	13
CH2--HH2	1.08	0.93	14
NE2--HE2	1.01	0.86	4

Standard geometry: angle outliers?

There are 35 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	117.04	1
CA-CB-CG	112.60	118.11	1
CB-CG-CD2	131.20	124.16	1
CA-CB-CG	112.60	117.91	1
OE1-CD-NE2	122.60	117.45	1
OE1-CD-NE2	122.60	117.71	2
OD1-CG-ND2	122.60	117.76	1
OD1-CG-ND2	122.60	117.84	1
NE-CZ-NH2	119.20	123.44	1
OE1-CD-NE2	122.60	117.93	1

Angle type	Observed angle (°)	Ideal angle (°)	Number of outliers
OE1-CD-NE2	122.60	118.04	1
OE1-CD-NE2	122.60	118.05	1
OE1-CD-NE2	122.60	118.14	1
OE1-CD-NE2	122.60	118.19	1
C-CA-CB	110.10	118.36	1
CD-NE-CZ	124.40	130.42	1
NE-CZ-NH2	119.20	123.04	1
CB-CG-CD2	131.20	125.65	1
CB-CG-CD2	131.20	125.66	1
CA-CB-CG	112.60	116.84	1
NH1-CZ-NH2	119.30	113.78	1
OE1-CD-NE2	122.60	118.37	1
CA-CB-CG	113.80	117.93	1
CB-CG-CD2	131.20	125.83	1
OE1-CD-NE2	122.60	118.48	2
CA-CB-CG	112.60	108.48	1
OD1-CG-ND2	122.60	118.50	1
OE1-CD-NE2	122.60	118.50	1
OE1-CD-NE2	122.60	118.51	1
OE1-CD-NE2	122.60	118.56	1
OE1-CD-NE2	122.60	118.57	1
CB-CG-CD2	131.20	125.98	1
OE1-CD-NE2	122.60	118.60	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	1.02	21

All 21 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Model ID	Atom-1	Atom-2	Clash overlap (Å)
1	A:117:VAL:HG21	A:906:MET:HE2	0.768
1	A:215:LEU:HD21	A:225:ILE:HG21	0.721
1	A:397:MET:HE1	A:563:TYR:CE1	0.640
1	A:117:VAL:HG21	A:906:MET:CE	0.626
1	A:304:LEU:HD22	A:333:TYR:CG	0.598
1	A:314:ALA:HA	A:317:TYR:CZ	0.595
1	B:350:LYS:HE3	B:351:TYR:CE2	0.588
1	A:491:VAL:HA	A:494:TYR:CE1	0.567
1	A:192:LEU:HD12	A:225:ILE:HD12	0.541
1	A:679:ARG:HH21	A:758:TRP:CG	0.538
1	A:193:GLU:C	A:339:ARG:HH22	0.523
1	A:304:LEU:HD22	A:333:TYR:CD1	0.508
1	B:190:LYS:HE3	B:306:MET:HE1	0.497
1	A:781:VAL:HG23	A:808:THR:HG23	0.492
1	B:326:ILE:HG22	B:360:SER:HB2	0.480
1	B:116:PHE:CE2	B:117:VAL:HG23	0.478
1	A:245:LEU:HD13	A:901:VAL:HG11	0.441
1	A:104:VAL:HG12	A:166:PHE:CZ	0.425
1	A:732:MET:HB3	A:796:TRP:CE3	0.409

Model ID	Atom-1	Atom-2	Clash overlap (Å)
1	A:314:ALA:HA	A:317:TYR:CE2	0.407
1	A:307:ARG:HH12	A:327:ASP:CG	0.401

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	1328	1294	34	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	1101	1057	27	17

Detailed list of outliers are tabulated below.

Model ID	Chain	Residue ID	Residue type
1	A	7	THR
1	A	52	LEU
1	A	73	LEU
1	A	192	LEU
1	A	202	ILE
1	A	304	LEU
1	A	371	VAL
1	A	405	LEU
1	A	435	THR
1	A	540	LEU
1	A	612	LEU

Model ID	Chain	Residue ID	Residue type
1	A	643	ASP
1	A	691	ASP
1	A	890	ASP
1	B	10	SER
1	B	362	VAL
1	B	397	TYR

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 ?

Validation for this section is under development.

Acknowledgements

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