Integrative Structure Validation Report

July 22, 2024 - 05:22 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	9A3M
PDB-Dev ID	PDBDEV_00000207
Structure Title	Parathyroid hormone receptor type 1 in complex with a long-acting parathyroid hormone analog and arrestin 2 (6pwc-based template)
Structure Authors	Aydin, Y.; Bottke, T.; Lam, J.H.; Ernicke, S.; Fortmann, A.; Tretbar, M.; Zarzycka, B.; Gurevich, V.V.; Katritch, V.; Coin, I.

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

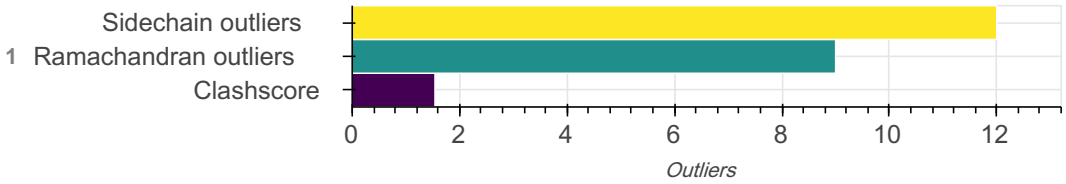
A user guide is available at https://pdb-dev.wwpdb.org/validation_help.html with specific help available everywhere you see the ? symbol.

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 3 subunits in each model. A total of 6 datasets or restraints were used to build this entry. Each model is represented by 0 rigid bodies and 3 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	Arrestin2	A	A	357
1	2	2	Long-acting parathyroid hormone analog	B	B	32
1	3	3	PTH1R	C	P	504

Datasets used for modeling

There are 6 unique datasets used to build the models in this entry.

ID	Dataset type	Database name	Data access code
1	Crosslinking-MS data	Not available	Not available
2	Experimental model	PDB	6NBF
3	Experimental model	PDB	6PWC

ID	Dataset type	Database name	Data access code
4	Comparative model	Not available	Not available
5	Comparative model	Not available	Not available
6	De Novo model	Not available	Not available

Representation

This entry has only one representation and includes 0 rigid bodies and 3 flexible units

Chain ID	Rigid bodies	Non-rigid segments
C	-	1-504
A	-	1-357
B	-	1-32

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	None	None	None	None	False	False

There is 1 software package reported in this entry.

ID	Software name	Software version	Software classification	Software location
1	ICM-Pro	v.3.9.2c	Model building	https://www.molsoft.com/icm_pro.html

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 7097 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
CA--HA	1.09	0.98	4
CA--HA2	1.09	0.97	51
CD--HD3	1.09	0.97	162
CB--HB2	1.09	0.97	663
CG--HG2	1.09	0.97	255
CG2--HG23	1.09	0.97	166
CG--HG3	1.09	0.97	255
CD1--HD11	1.09	0.97	132
CG2--HG21	1.09	0.97	166
CB--HB1	1.09	0.97	71
CB--HB3	1.09	0.97	663
CA--HA	1.09	0.97	830
CD2--HD22	1.09	0.97	93
CG--HG	1.09	0.97	93
CD--HD2	1.09	0.97	162
CB--HB	1.09	0.97	167
CD1--HD12	1.09	0.97	132
CG1--HG11	1.09	0.97	74
CA--HA3	1.09	0.97	51

Bond type	Observed distance (Å)	Ideal distance (Å)	Number of outliers
CE--HE3	1.09	0.97	73
CG1--HG12	1.09	0.97	113
CE--HE1	1.09	0.97	16
CE--HE2	1.09	0.97	73
CD1--HD13	1.09	0.97	132
CG2--3HG2	1.09	0.97	1
CG2--HG22	1.09	0.97	166
CD2--HD23	1.09	0.97	93
CD2--HD21	1.09	0.97	93
CG1--HG13	1.09	0.97	113
CB--2HB	1.09	0.97	4
CG2--1HG2	1.09	0.97	1
CG2--2HG2	1.09	0.97	1
NZ--HZ3	1.01	0.89	57
NZ--HZ1	1.01	0.89	57
NZ--HZ2	1.01	0.89	57
SG--HG	1.33	1.20	13
N--H	1.00	0.86	828
NH1--HH12	1.00	0.86	56
NH2--HH21	1.00	0.86	56
ND2--HD21	1.00	0.86	30
NH1--HH11	1.00	0.86	56
NE--HE	1.00	0.86	56

Bond type	Observed distance (Å)	Ideal distance (Å)	Number of outliers
ND1--HD1	1.00	0.86	15
ND2--HD22	1.00	0.86	30
NH2--HH22	1.00	0.86	56
NE2--HE21	1.00	0.86	23
NE2--HE22	1.00	0.86	23
NE1--HE1	1.00	0.86	15
CH2--HH2	1.09	0.93	15
OG--HG	1.00	0.84	47
CE3--HE3	1.09	0.93	15
CD2--HD2	1.09	0.93	94
CD1--HD1	1.09	0.93	94
CE2--HE2	1.09	0.93	79
CZ--HZ	1.09	0.93	42
OG1--HG1	1.00	0.84	53
CE1--HE1	1.09	0.93	94
OH--HH	1.00	0.84	37
CZ2--HZ2	1.09	0.93	15
CZ3--HZ3	1.09	0.93	15

Standard geometry: angle outliers

Bond angle outliers do not exist or can not be evaluated for this model

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

Model ID	Clash score	Number of clashes
1	1.54	22

All 22 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	C:165:TYR:C	C:165:TYR:CD2	0.669
1	C:165:TYR:CD2	C:166:THR:N	0.596
1	C:71:ALA:N	C:72:PRO:CD	0.577
1	C:62:TYR:N	C:63:PRO:HD2	0.571
1	A:224:THR:OG1	A:225:ASN:N	0.536
1	C:339:VAL:N	C:340:PRO:CD	0.531
1	C:305:LEU:N	C:306:PRO:CD	0.524
1	A:243:LEU:C	A:243:LEU:HD23	0.509
1	A:87:PHE:HA	A:88:PRO:C	0.493
1	C:62:TYR:HB3	C:63:PRO:HD3	0.487
1	A:203:LEU:HD22	A:349:PHE:CE2	0.481
1	C:321:ALA:O	C:322:ASN:C	0.459
1	C:388:MET:HB2	C:389:PRO:HD3	0.456
1	C:62:TYR:N	C:63:PRO:CD	0.453
1	C:123:ASP:C	C:123:ASP:OD1	0.447
1	C:1:ASP:C	C:1:ASP:OD1	0.437
1	C:490:LEU:HA	C:491:PRO:C	0.436
1	A:95:LYS:N	A:96:PRO:HD2	0.425
1	C:47:SER:O	C:48:THR:C	0.421
1	C:388:MET:N	C:389:PRO:CD	0.417

Model ID	Atom-1	Atom-2	Clash overlap (Å)
1	C:70:GLU:O	C:73:THR:HG22	0.407
1	C:466:SER:OG	C:467:SEP:N	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	850	796	45	9

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	758	711	35	12

Detailed list of outliers are tabulated below.

Model ID	Chain	Residue ID	Residue type
1	A	6	THR
1	A	18	LEU
1	A	26	ASP
1	A	58	THR
1	A	79	LEU
1	A	143	ASP
1	A	156	GLU
1	A	214	ILE
1	A	335	LEU
1	A	353	HIS
1	C	373	THR

Model ID	Chain	Residue ID	Residue type
1	C	479	VAL

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

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 byRCSB 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.