



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 03:16 AM UTC

PDB ID : 9B96 / pdb_00009b96
Title : Crystal structure of the human PAD2 protein bound to inhibitor
Authors : Byrnes, L.J.; Vajdos, F.
Deposited on : 2024-04-01
Resolution : 1.64 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

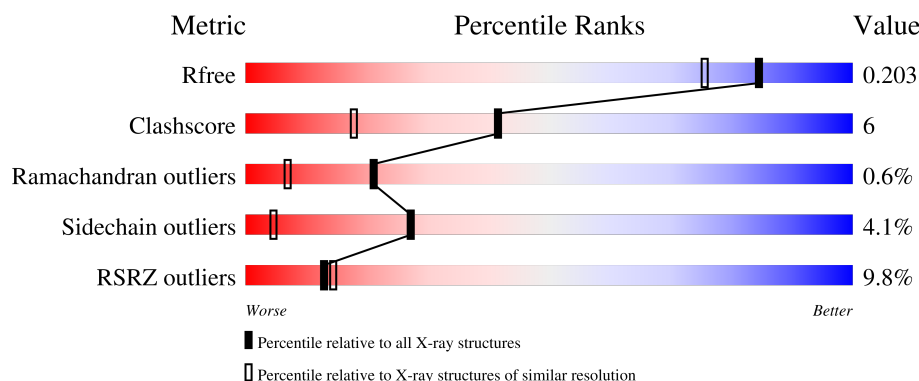
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.64 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1141 (1.64-1.64)
Clashscore	190562	1171 (1.64-1.64)
Ramachandran outliers	187476	1151 (1.64-1.64)
Sidechain outliers	187428	1150 (1.64-1.64)
RSRZ outliers	180081	1141 (1.64-1.64)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	690	9% 77% 13% 8%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5613 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Protein-arginine deiminase type-2.

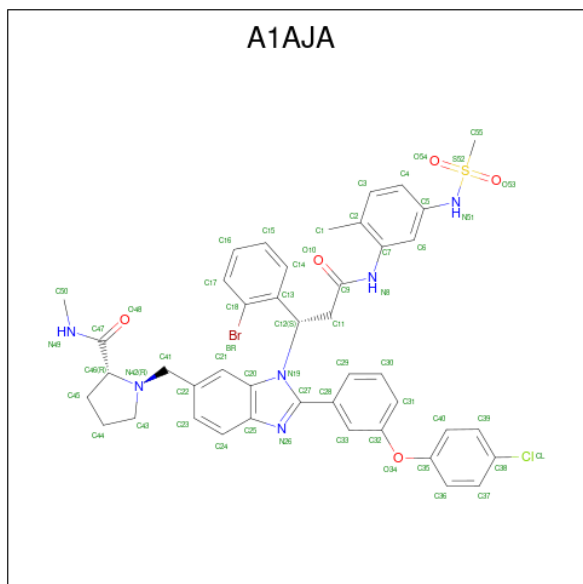
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	633	Total	C	N	O	S	0	7	0
			5073	3261	840	940	32			

There are 25 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	initiating methionine	UNP Q9Y2J8
A	-19	GLY	-	expression tag	UNP Q9Y2J8
A	-18	HIS	-	expression tag	UNP Q9Y2J8
A	-17	HIS	-	expression tag	UNP Q9Y2J8
A	-16	HIS	-	expression tag	UNP Q9Y2J8
A	-15	HIS	-	expression tag	UNP Q9Y2J8
A	-14	HIS	-	expression tag	UNP Q9Y2J8
A	-13	HIS	-	expression tag	UNP Q9Y2J8
A	-12	HIS	-	expression tag	UNP Q9Y2J8
A	-11	HIS	-	expression tag	UNP Q9Y2J8
A	-10	HIS	-	expression tag	UNP Q9Y2J8
A	-9	HIS	-	expression tag	UNP Q9Y2J8
A	-8	SER	-	expression tag	UNP Q9Y2J8
A	-7	SER	-	expression tag	UNP Q9Y2J8
A	-6	GLY	-	expression tag	UNP Q9Y2J8
A	-5	HIS	-	expression tag	UNP Q9Y2J8
A	-4	ILE	-	expression tag	UNP Q9Y2J8
A	-3	GLU	-	expression tag	UNP Q9Y2J8
A	-2	GLY	-	expression tag	UNP Q9Y2J8
A	-1	ARG	-	expression tag	UNP Q9Y2J8
A	0	HIS	-	expression tag	UNP Q9Y2J8
A	666	SER	-	expression tag	UNP Q9Y2J8
A	667	ARG	-	expression tag	UNP Q9Y2J8
A	668	ARG	-	expression tag	UNP Q9Y2J8
A	669	SER	-	expression tag	UNP Q9Y2J8

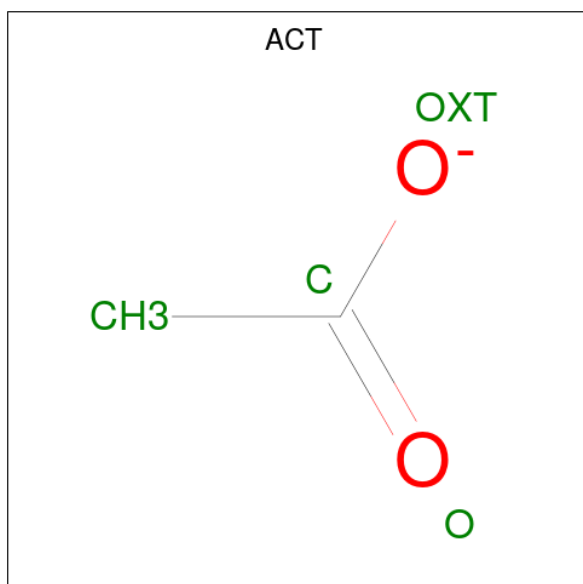
- Molecule 2 is 1-((2P)-1-((1R)-1-(2-bromophenyl)-3-[5-(methanesulfonamido)-2-methylanili

no]-3-oxopropyl}-2-[3-(4-chlorophenoxy)phenyl]-1H-1,3-benzimidazol-6-yl)methyl)-N-methyl-D-prolinamide (CCD ID: A1AJA) (formula: $C_{43}H_{42}BrClN_6O_5S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
			Total	Br	C	Cl	N	O	S		
2	A	1	57	1	43	1	6	5	1	0	0

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	4	2	2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total	Ca	0	0
			4	4		

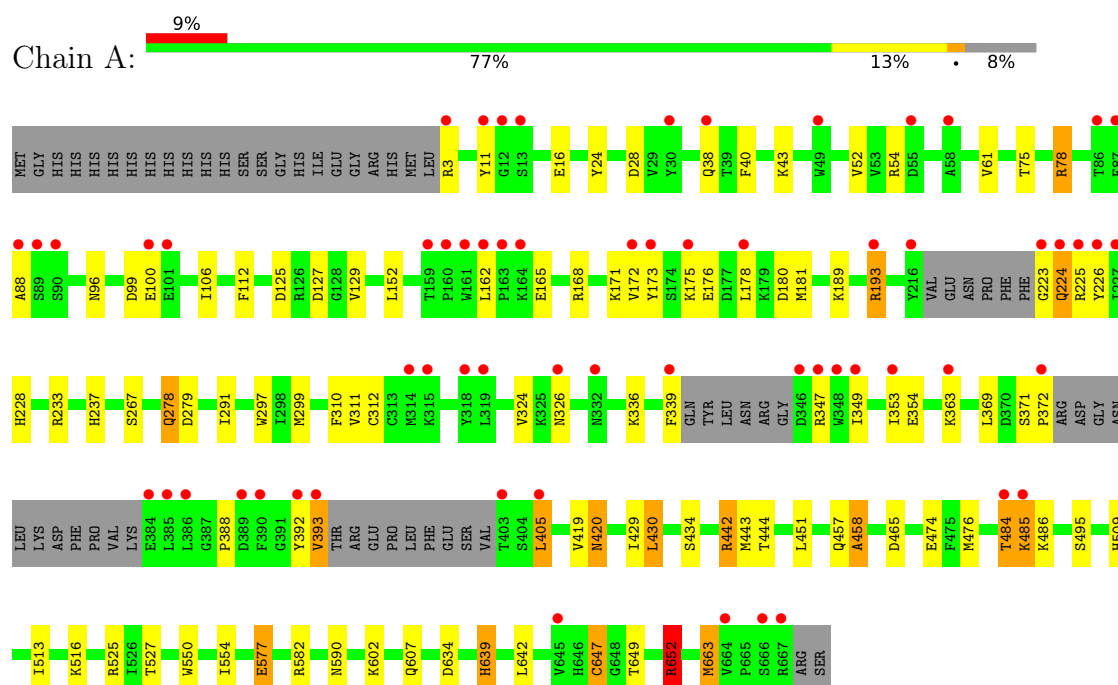
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	471	Total	O	0	0
			471	471		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Protein-arginine deiminase type-2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	202.26Å 50.73Å 75.56Å 90.00° 105.08° 90.00°	Depositor
Resolution (Å)	97.65 – 1.64 97.65 – 1.64	Depositor EDS
% Data completeness (in resolution range)	67.1 (97.65-1.64) 67.1 (97.65-1.64)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 1.64Å)	Xtriage
Refinement program	REFMAC 5.8.0349	Depositor
R, R_{free}	0.172 , 0.211 (Not available) , 0.203	Depositor DCC
R_{free} test set	2970 reflections (3.26%)	wwPDB-VP
Wilson B-factor (Å ²)	22.2	Xtriage
Anisotropy	0.077	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5613	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.37% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, ACT, A1AJA

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.64	4/5207 (0.1%)	0.98	5/7051 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	10

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	639	HIS	CE1-NE2	7.18	1.39	1.32
1	A	228	HIS	CE1-NE2	5.89	1.38	1.32
1	A	237	HIS	CE1-NE2	5.53	1.38	1.32
1	A	509	HIS	CE1-NE2	5.45	1.38	1.32

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	634	ASP	CA-CB-CG	6.55	119.15	112.60
1	A	484	THR	CB-CA-C	6.15	123.64	111.22
1	A	28	ASP	CA-CB-CG	5.61	118.21	112.60
1	A	484	THR	CA-C-N	5.28	131.21	121.70
1	A	484	THR	C-N-CA	5.28	131.21	121.70

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	168	ARG	Sidechain
1	A	193	ARG	Sidechain
1	A	233	ARG	Sidechain
1	A	3	ARG	Sidechain
1	A	347	ARG	Sidechain
1	A	442	ARG	Sidechain
1	A	525	ARG	Sidechain
1	A	582	ARG	Sidechain
1	A	652	ARG	Sidechain
1	A	78	ARG	Sidechain

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5073	0	5050	65	0
2	A	57	0	0	0	0
3	A	8	0	6	0	0
4	A	4	0	0	0	0
5	A	471	0	0	14	2
All	All	5613	0	5056	65	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (65) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:485:LYS:HA	5:A:1187:HOH:O	1.31	1.31
1:A:484:THR:HB	1:A:485:LYS:HB3	1.32	1.11
1:A:663:MET:HE1	5:A:1249:HOH:O	1.65	0.97
1:A:339:PHE:C	5:A:938:HOH:O	2.08	0.95
1:A:484:THR:CB	1:A:485:LYS:HB3	2.07	0.84
1:A:474[A]:GLU:OE2	1:A:590:ASN:ND2	2.13	0.81
1:A:88:ALA:O	5:A:801:HOH:O	1.98	0.80
1:A:652:ARG:HD3	5:A:1078:HOH:O	1.83	0.79
1:A:484:THR:HA	1:A:485:LYS:HB2	1.67	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:513:ILE:HD13	1:A:527:THR:HG22	1.73	0.69
1:A:484:THR:HB	1:A:485:LYS:CB	2.19	0.65
1:A:99:ASP:HB3	5:A:1225:HOH:O	1.98	0.64
1:A:577:GLU:H	1:A:577:GLU:CD	2.06	0.62
1:A:484:THR:CA	1:A:485:LYS:CB	2.78	0.61
1:A:647:CYS:HB2	5:A:1130:HOH:O	2.01	0.60
1:A:484:THR:HA	1:A:485:LYS:CB	2.31	0.59
1:A:312[B]:CYS:SG	1:A:349:ILE:HB	2.43	0.59
1:A:392:TYR:O	1:A:393:VAL:C	2.44	0.59
1:A:162:LEU:HD11	1:A:173:TYR:CE2	2.37	0.58
1:A:550:TRP:CZ3	1:A:554:ILE:HD11	2.40	0.57
1:A:223:GLY:O	1:A:224:GLN:CB	2.55	0.55
1:A:639:HIS:HD2	5:A:1093:HOH:O	1.90	0.55
1:A:152:LEU:HD11	1:A:363:LYS:HG3	1.90	0.54
1:A:550:TRP:CH2	1:A:554:ILE:HD11	2.43	0.54
1:A:54:ARG:HG2	1:A:75:THR:OG1	2.06	0.54
1:A:38:GLN:HG3	1:A:100:GLU:OE1	2.09	0.53
1:A:172:VAL:CG2	1:A:181:MET:HE1	2.40	0.52
1:A:311:VAL:HG22	1:A:649:THR:HG22	1.92	0.52
1:A:550:TRP:CZ3	1:A:554:ILE:CD1	2.93	0.52
1:A:430:LEU:HD13	1:A:476:MET:HE1	1.92	0.52
1:A:495:SER:HB3	5:A:1255:HOH:O	2.10	0.50
1:A:388:PRO:O	5:A:802:HOH:O	2.20	0.50
1:A:602:LYS:HE3	5:A:1029:HOH:O	2.12	0.50
1:A:178:LEU:HD11	1:A:226:TYR:HE1	1.75	0.50
1:A:52:VAL:HG23	1:A:61:VAL:HG21	1.95	0.49
1:A:106:ILE:CD1	5:A:1225:HOH:O	2.62	0.48
1:A:43:LYS:HE3	5:A:936:HOH:O	2.13	0.47
1:A:299:MET:HE3	1:A:354:GLU:CD	2.38	0.47
1:A:178:LEU:HD11	1:A:226:TYR:CE1	2.50	0.46
1:A:419:VAL:O	1:A:420:ASN:ND2	2.49	0.46
1:A:474[A]:GLU:CD	1:A:590:ASN:HD22	2.22	0.46
1:A:513:ILE:HD13	1:A:527:THR:CG2	2.45	0.46
1:A:297:TRP:CE2	1:A:369:LEU:HG	2.52	0.45
1:A:16:GLU:HG2	1:A:112:PHE:HB2	1.99	0.45
1:A:40:PHE:HA	1:A:96:ASN:O	2.17	0.44
1:A:405:LEU:HD22	1:A:443:MET:HB2	1.99	0.44
1:A:176:GLU:HG2	5:A:845:HOH:O	2.18	0.44
1:A:457:GLN:O	1:A:458:ALA:C	2.60	0.44
1:A:180:ASP:HA	1:A:363:LYS:HE2	2.00	0.44
1:A:223:GLY:HA2	1:A:225:ARG:HH22	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:516[A]:LYS:HD2	1:A:607:GLN:HG3	2.00	0.43
1:A:278:GLN:O	1:A:279:ASP:HB2	2.18	0.43
1:A:516[B]:LYS:HZ2	1:A:516[B]:LYS:HB2	1.84	0.43
1:A:429:ILE:HG21	1:A:451:LEU:HD13	2.00	0.42
1:A:223:GLY:CA	1:A:225:ARG:HH22	2.32	0.42
1:A:267:SER:OG	1:A:291:ILE:CD1	2.68	0.42
1:A:162:LEU:HD11	1:A:173:TYR:HE2	1.80	0.42
1:A:310:PHE:CE2	1:A:336:LYS:HD2	2.54	0.42
1:A:172:VAL:HG22	1:A:181:MET:HE1	2.01	0.42
1:A:516[B]:LYS:HB2	1:A:516[B]:LYS:NZ	2.35	0.41
1:A:24:TYR:CE2	1:A:78:ARG:HD2	2.55	0.41
1:A:125:ASP:C	1:A:125:ASP:OD1	2.62	0.41
1:A:127:ASP:OD1	1:A:127:ASP:C	2.63	0.41
1:A:311:VAL:HG21	1:A:324:VAL:HG11	2.03	0.40
1:A:434:SER:HA	1:A:465:ASP:HB2	2.04	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:1198:HOH:O	5:A:1224:HOH:O[4_746]	2.13	0.07
5:A:960:HOH:O	5:A:1131:HOH:O[4_756]	2.17	0.03

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	630/690 (91%)	612 (97%)	14 (2%)	4 (1%)	21 7

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	224	GLN
1	A	485	LYS
1	A	11	TYR
1	A	458	ALA

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	565/612 (92%)	542 (96%)	23 (4%)	27 4

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	129	VAL
1	A	165	GLU
1	A	171	LYS
1	A	175	LYS
1	A	189	LYS
1	A	193	ARG
1	A	278	GLN
1	A	326	ASN
1	A	353	ILE
1	A	371	SER
1	A	372	PRO
1	A	393	VAL
1	A	405	LEU
1	A	420	ASN
1	A	430	LEU
1	A	442	ARG
1	A	444	THR
1	A	486	LYS
1	A	577	GLU
1	A	642	LEU
1	A	647	CYS
1	A	652	ARG
1	A	663	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	96	ASN
1	A	108	GLN
1	A	420	ASN
1	A	639	HIS
1	A	646	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	A1AJA	A	701	-	63,63,63	0.60	1 (1%)	89,91,91	0.81	3 (3%)
3	ACT	A	702	-	3,3,3	1.04	0	3,3,3	0.77	0
3	ACT	A	703	-	3,3,3	0.86	0	3,3,3	0.78	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	A1AJA	A	701	-	-	4/39/49/49	0/7/7/7

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	A1AJA	BR-C18	2.80	1.96	1.89

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	A1AJA	C13-C12-N19	3.82	116.00	111.90
2	A	701	A1AJA	C21-C20-N19	2.47	132.93	128.59
2	A	701	A1AJA	C47-C46-N42	2.04	115.00	112.38

There are no chirality outliers.

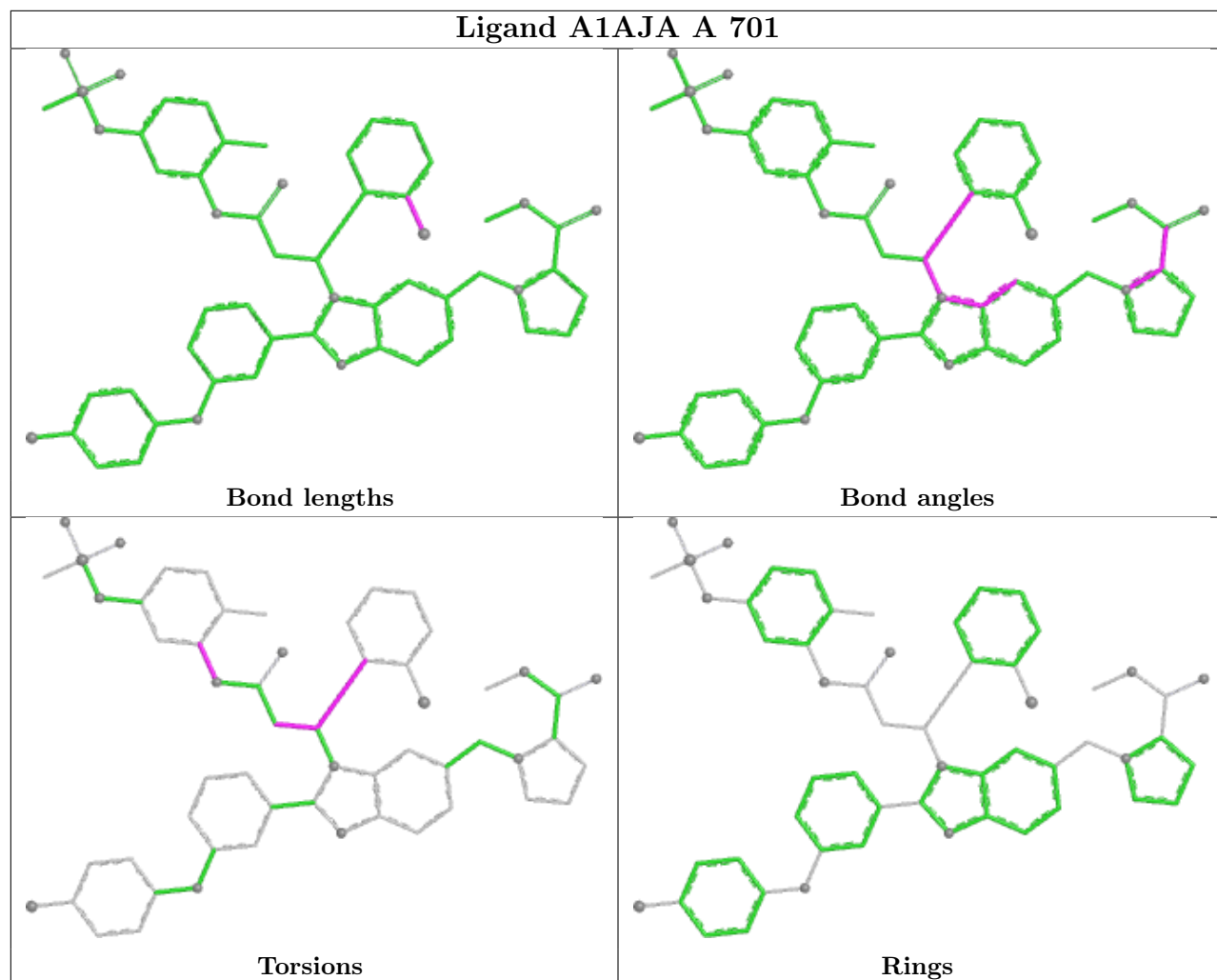
All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	A1AJA	N19-C12-C13-C14
2	A	701	A1AJA	C9-C11-C12-C13
2	A	701	A1AJA	C6-C7-N8-C9
2	A	701	A1AJA	C2-C7-N8-C9

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	633/690 (91%)	0.43	62 (9%) 13 14	9, 24, 56, 83	7 (1%)

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	393	VAL	9.3
1	A	161	TRP	7.6
1	A	162	LEU	6.3
1	A	392	TYR	5.6
1	A	372	PRO	4.9
1	A	88	ALA	4.7
1	A	90	SER	4.5
1	A	49	TRP	4.4
1	A	224	GLN	4.1
1	A	385	LEU	4.0
1	A	349	ILE	3.9
1	A	403	THR	3.6
1	A	225	ARG	3.6
1	A	163	PRO	3.6
1	A	13	SER	3.5
1	A	172	VAL	3.3
1	A	160	PRO	3.3
1	A	216	TYR	3.2
1	A	89	SER	3.2
1	A	666	SER	3.2
1	A	178	LEU	3.1
1	A	346	ASP	3.1
1	A	353	ILE	3.0
1	A	55	ASP	3.0
1	A	645	VAL	2.9
1	A	164	LYS	2.9
1	A	667	ARG	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	314	MET	2.7
1	A	318	TYR	2.7
1	A	159	THR	2.7
1	A	12	GLY	2.6
1	A	223	GLY	2.6
1	A	386	LEU	2.6
1	A	664	VAL	2.6
1	A	348	TRP	2.6
1	A	3	ARG	2.6
1	A	226	TYR	2.6
1	A	389	ASP	2.5
1	A	87	GLU	2.5
1	A	339	PHE	2.5
1	A	484	THR	2.5
1	A	173	TYR	2.5
1	A	363	LYS	2.5
1	A	175	LYS	2.4
1	A	390	PHE	2.4
1	A	86	THR	2.4
1	A	193	ARG	2.3
1	A	485	LYS	2.3
1	A	101	GLU	2.3
1	A	100	GLU	2.3
1	A	384	GLU	2.3
1	A	347	ARG	2.2
1	A	11	TYR	2.2
1	A	326	ASN	2.2
1	A	227	ILE	2.2
1	A	38	GLN	2.1
1	A	315	LYS	2.1
1	A	405	LEU	2.1
1	A	319	LEU	2.1
1	A	332	ASN	2.0
1	A	30	TYR	2.0
1	A	58	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

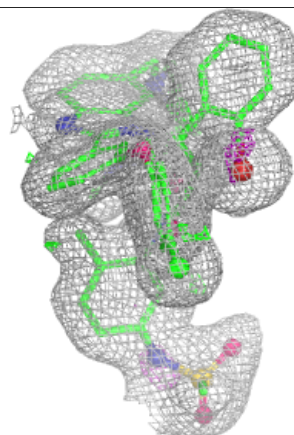
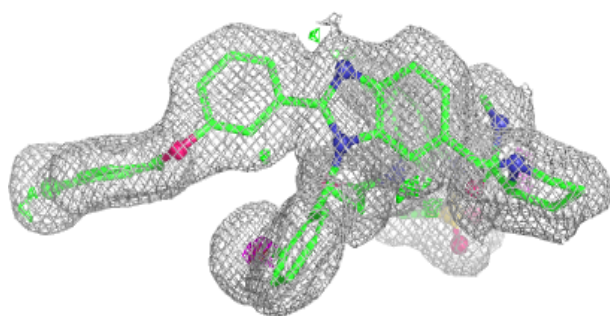
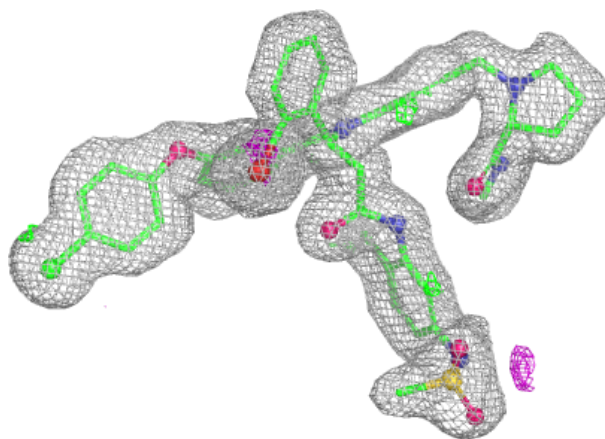
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ACT	A	702	4/4	0.88	0.16	48,49,52,54	0
4	CA	A	705	1/1	0.90	0.13	48,48,48,48	0
4	CA	A	704	1/1	0.91	0.12	52,52,52,52	0
3	ACT	A	703	4/4	0.91	0.13	40,41,42,53	0
4	CA	A	707	1/1	0.94	0.07	38,38,38,38	0
4	CA	A	706	1/1	0.97	0.08	49,49,49,49	0
2	A1AJA	A	701	57/57	0.98	0.06	14,20,34,37	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around A1AJA A 701:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.