



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:25 PM UTC

PDB ID : 2J0N / pdb_00002j0n
Title : A proteolytically truncated form of Shigella Flexneri IpaD
Authors : Johnson, S.; Roversi, P.; Lea, S.M.
Deposited on : 2006-08-03
Resolution : 2.10 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
Xtriage (Phenix)	:	2.0
EDS	:	3.0
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

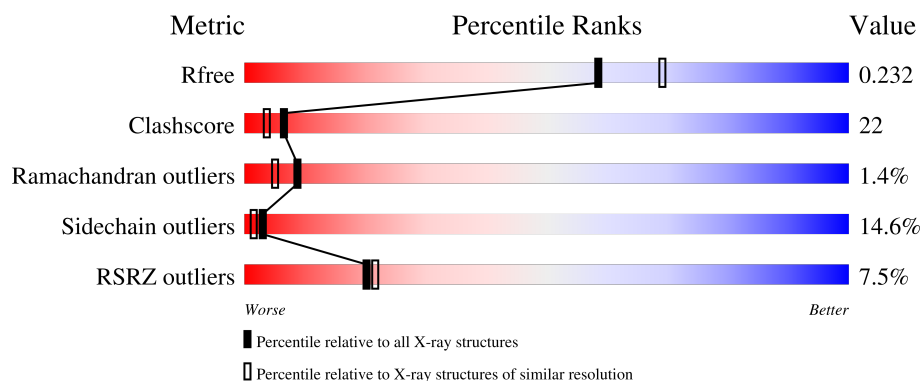
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	200	
1	B	200	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2986 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 INVASIN IPAD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	171	Total	C	N	O	S	0	0	0
			1329	835	221	269	4			
1	B	187	Total	C	N	O	S	0	0	0
			1452	914	241	293	4			

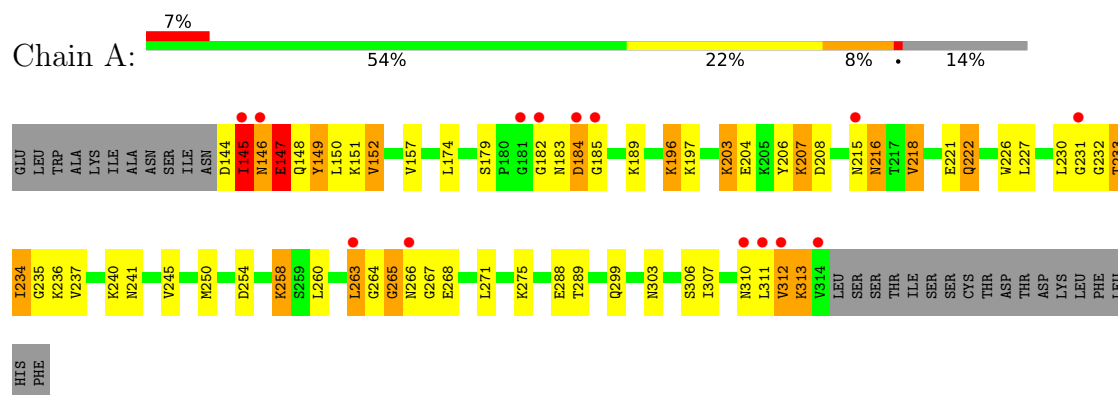
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	111	Total	O	0	0
			111	111		
2	B	94	Total	O	0	0
			94	94		

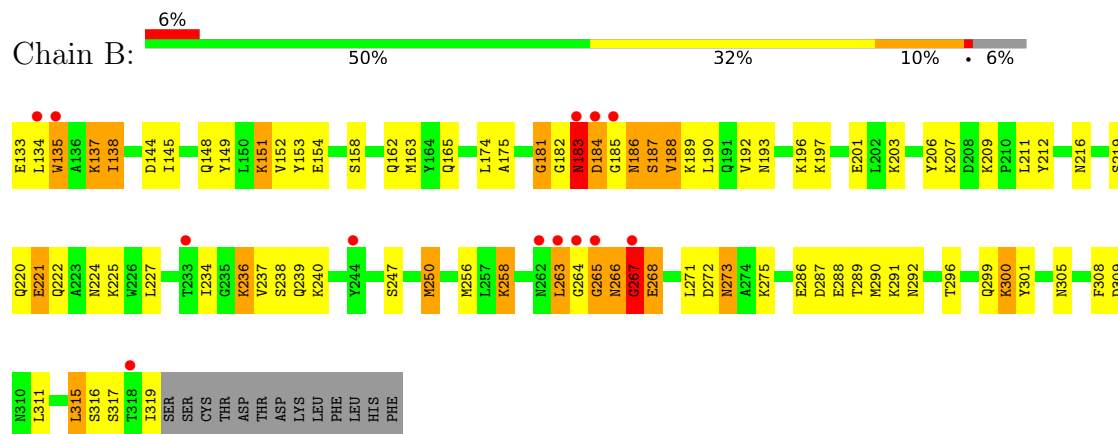
3 Residue-property plots [i](#)

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: INVASIN IPAD



• Molecule 1: INVASIN IPAD



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	77.79Å 91.36Å 54.91Å 90.00° 96.35° 90.00°	Depositor
Resolution (Å)	15.00 – 2.10 15.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	(Not available) (15.00-2.10) 99.5 (15.00-2.10)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 1.89Å)	Xtriage
Refinement program	TNT 5.6.1	Depositor
R, R_{free}	0.189 , (Not available) 0.199 , 0.232	Depositor DCC
R_{free} test set	1525 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	0.265	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 71.3	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	0.95	EDS
Total number of atoms	2986	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.49% 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

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	0.56	0/1351	1.04	13/1828 (0.7%)
1	B	0.59	2/1476 (0.1%)	1.23	17/1999 (0.9%)
All	All	0.58	2/2827 (0.1%)	1.15	30/3827 (0.8%)

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	1
1	B	0	6
All	All	0	7

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	268	GLU	N-CA	-6.78	1.38	1.46
1	B	266	ASN	N-CA	-6.05	1.37	1.46

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	263	LEU	N-CA-C	23.15	136.59	111.36
1	B	267	GLY	CA-C-N	11.55	139.80	122.47
1	B	267	GLY	C-N-CA	11.55	139.80	122.47
1	B	135	TRP	N-CA-C	-10.35	101.41	114.56
1	A	313	LYS	N-CA-C	-9.32	101.69	113.43
1	B	183	ASN	N-CA-C	8.11	123.16	108.17
1	A	313	LYS	CB-CA-C	7.87	122.88	109.65
1	B	266	ASN	N-CA-C	-7.65	101.09	110.88
1	A	263	LEU	N-CA-C	-7.50	95.38	108.20
1	A	149	TYR	N-CA-C	-7.27	103.38	112.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	179	SER	N-CA-C	7.04	114.56	108.22
1	B	183	ASN	CA-C-N	6.90	134.12	121.70
1	B	183	ASN	C-N-CA	6.90	134.12	121.70
1	B	301	TYR	N-CA-C	-6.80	103.99	112.90
1	A	215	ASN	N-CA-C	-6.72	98.44	109.40
1	A	174	LEU	N-CA-C	6.29	118.66	111.11
1	B	316	SER	N-CA-C	-6.25	105.48	113.23
1	B	188	VAL	N-CA-C	6.01	116.53	108.11
1	B	211	LEU	N-CA-C	-6.00	104.67	112.23
1	B	273	ASN	N-CA-C	5.90	118.19	111.11
1	B	184	ASP	N-CA-C	-5.90	94.48	111.00
1	A	147	GLU	N-CA-C	-5.84	106.00	113.01
1	A	218	VAL	N-CA-C	5.78	115.87	108.82
1	A	266	ASN	N-CA-C	-5.62	101.99	110.70
1	A	145	ILE	N-CA-C	-5.59	100.17	108.45
1	B	187	SER	N-CA-C	5.46	117.36	109.07
1	B	174	LEU	N-CA-C	5.44	117.91	111.33
1	A	310	ASN	N-CA-C	5.41	116.99	111.14
1	A	312	VAL	CB-CA-C	5.27	119.93	111.29
1	B	183	ASN	CA-C-O	-5.13	115.66	121.66

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	313	LYS	Mainchain
1	B	183	ASN	Peptide,Mainchain
1	B	265	GLY	Peptide,Mainchain
1	B	267	GLY	Peptide,Mainchain

5.2 Too-close contacts ⓘ

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	1329	0	1303	60	0
1	B	1452	0	1431	72	1
2	A	111	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	94	0	0	7	0
All	All	2986	0	2734	120	1

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 22.

All (120) 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:B:163:MET:HE1	1:B:290:MET:HE3	1.37	1.03
1:B:192:VAL:HG21	1:B:265:GLY:O	1.58	1.03
1:B:192:VAL:CG2	1:B:267:GLY:H	1.76	0.97
1:A:233:THR:HG22	1:B:165:GLN:HE22	1.40	0.87
1:A:233:THR:HG22	1:B:165:GLN:NE2	1.90	0.86
1:A:147:GLU:HB3	1:A:148:GLN:NE2	1.90	0.85
1:B:192:VAL:CG2	1:B:267:GLY:N	2.47	0.77
1:B:227:LEU:HD22	1:B:237:VAL:HG23	1.68	0.75
1:B:209:LYS:HD2	2:B:2044:HOH:O	1.87	0.75
1:B:197:LYS:O	1:B:201:GLU:HG3	1.86	0.75
1:B:134:LEU:HB2	2:B:2002:HOH:O	1.87	0.73
1:A:307:ILE:O	1:A:311:LEU:HD23	1.90	0.71
1:B:256:MET:HE3	1:B:286:GLU:HB2	1.72	0.71
1:B:192:VAL:HG22	1:B:267:GLY:H	1.53	0.70
1:B:163:MET:CE	1:B:290:MET:HE3	2.20	0.70
1:A:230:LEU:HB3	1:A:234:ILE:HD11	1.75	0.68
1:A:203:LYS:HE3	1:A:254:ASP:OD1	1.93	0.68
1:A:258:LYS:NZ	2:A:2087:HOH:O	2.27	0.68
1:B:236:LYS:HE3	1:B:247:SER:HB2	1.76	0.68
1:B:133:GLU:HG2	1:B:134:LEU:N	2.09	0.67
1:B:189:LYS:HD3	1:B:268:GLU:CD	2.19	0.66
1:A:289:THR:HG21	1:B:165:GLN:HG2	1.76	0.66
1:B:133:GLU:HG3	2:B:2003:HOH:O	1.96	0.65
1:A:299:GLN:O	1:B:305:ASN:ND2	2.30	0.65
1:A:288:GLU:HG3	1:B:291:LYS:HE2	1.79	0.64
1:B:292:ASN:O	1:B:296:THR:HG23	1.98	0.63
1:B:192:VAL:HG22	1:B:267:GLY:N	2.13	0.62
1:A:231:GLY:O	1:A:233:THR:N	2.27	0.62
1:B:189:LYS:HD3	1:B:268:GLU:OE2	1.99	0.62
1:A:204:GLU:O	1:A:207:LYS:HD2	2.00	0.60
1:A:184:ASP:CG	1:A:185:GLY:N	2.59	0.60
1:A:307:ILE:HG22	1:A:311:LEU:CD2	2.30	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:THR:HG22	1:B:165:GLN:CD	2.25	0.60
1:B:236:LYS:CE	1:B:247:SER:HB2	2.33	0.58
1:A:234:ILE:HB	2:A:2106:HOH:O	2.03	0.58
1:A:230:LEU:HB3	1:A:234:ILE:CD1	2.35	0.56
1:A:307:ILE:HG22	1:A:311:LEU:HD21	1.86	0.56
1:B:182:GLY:O	1:B:184:ASP:HA	2.05	0.56
1:B:206:TYR:HB2	1:B:250:MET:CE	2.36	0.56
1:A:207:LYS:HG2	1:A:208:ASP:N	2.23	0.54
1:A:234:ILE:HG12	1:A:235:GLY:N	2.21	0.54
1:A:241:ASN:ND2	2:A:2075:HOH:O	2.40	0.54
1:A:288:GLU:HG2	2:A:2100:HOH:O	2.06	0.54
1:B:192:VAL:HG23	1:B:267:GLY:N	2.22	0.54
1:A:189:LYS:HD2	1:A:268:GLU:CD	2.33	0.53
1:A:254:ASP:O	1:A:258:LYS:HG2	2.08	0.53
1:B:288:GLU:HG3	2:B:2087:HOH:O	2.08	0.53
1:B:311:LEU:HD22	1:B:315:LEU:CD2	2.39	0.53
1:A:233:THR:HG23	1:B:162:GLN:HG2	1.89	0.53
1:B:272:ASP:OD2	1:B:275:LYS:HD3	2.09	0.53
1:B:163:MET:SD	1:B:290:MET:HE1	2.48	0.52
1:A:236:LYS:NZ	2:A:2071:HOH:O	2.29	0.52
1:B:151:LYS:HE2	2:B:2011:HOH:O	2.09	0.51
1:B:192:VAL:HG22	1:B:267:GLY:HA2	1.91	0.51
1:B:222:GLN:OE1	1:B:222:GLN:HA	2.11	0.51
1:A:307:ILE:HG23	1:A:311:LEU:HD22	1.92	0.51
1:B:192:VAL:HG22	1:B:267:GLY:CA	2.40	0.51
1:A:240:LYS:HD2	1:A:245:VAL:HG11	1.91	0.51
1:A:307:ILE:CG2	1:A:311:LEU:HD22	2.39	0.51
1:B:133:GLU:HG2	2:B:2002:HOH:O	2.10	0.50
1:B:158:SER:O	1:B:162:GLN:HG3	2.11	0.50
1:A:271:LEU:HD22	1:A:275:LYS:HG2	1.92	0.50
1:B:181:GLY:HA3	1:B:187:SER:OG	2.12	0.50
1:B:263:LEU:N	1:B:264:GLY:HA2	2.24	0.50
1:B:220:GLN:O	1:B:224:ASN:OD1	2.30	0.50
1:A:312:VAL:HG12	1:A:312:VAL:O	2.12	0.49
1:B:219:SER:OG	1:B:221:GLU:HG2	2.13	0.49
1:A:303:ASN:ND2	1:B:309:ASP:OD2	2.46	0.49
1:B:152:VAL:HG13	1:B:212:TYR:HD2	1.77	0.49
1:B:190:LEU:HD23	1:B:192:VAL:HG12	1.95	0.48
1:B:236:LYS:HB3	1:B:236:LYS:HE2	1.61	0.47
1:B:145:ILE:O	1:B:149:TYR:HB3	2.15	0.47
1:A:206:TYR:HB2	1:A:250:MET:CE	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:VAL:HG21	1:A:222:GLN:HG2	1.96	0.47
1:B:137:LYS:HD2	1:B:137:LYS:HA	1.49	0.47
1:A:265:GLY:C	1:A:267:GLY:H	2.23	0.47
1:A:307:ILE:CG2	1:A:311:LEU:CD2	2.93	0.46
1:A:146:ASN:HB3	1:A:148:GLN:H	1.81	0.46
1:A:184:ASP:OD1	1:A:185:GLY:N	2.48	0.46
1:A:182:GLY:HA3	1:A:183:ASN:HA	1.53	0.46
1:B:300:LYS:HE2	2:B:2093:HOH:O	2.15	0.46
1:B:258:LYS:HA	1:B:258:LYS:HD2	1.47	0.46
1:B:311:LEU:HD22	1:B:315:LEU:HD23	1.97	0.46
1:B:144:ASP:O	1:B:148:GLN:HG3	2.17	0.45
1:A:189:LYS:HB2	1:A:189:LYS:HE2	1.72	0.44
1:A:196:LYS:HB3	1:A:196:LYS:HE3	1.28	0.44
1:A:233:THR:HG22	1:B:165:GLN:OE1	2.18	0.44
1:B:192:VAL:HG21	1:B:265:GLY:C	2.36	0.44
1:A:227:LEU:HD22	1:A:237:VAL:HG23	2.00	0.44
1:A:145:ILE:HD11	1:A:150:LEU:HD21	2.00	0.44
1:A:203:LYS:HA	1:A:250:MET:HE3	2.00	0.44
1:A:312:VAL:O	1:A:312:VAL:CG1	2.65	0.44
1:A:307:ILE:C	1:A:311:LEU:HD23	2.43	0.43
1:B:134:LEU:HD23	1:B:134:LEU:HA	1.82	0.43
1:A:204:GLU:HG3	1:A:207:LYS:NZ	2.34	0.43
1:A:260:LEU:O	1:A:263:LEU:O	2.37	0.43
1:B:153:TYR:CE2	1:B:300:LYS:HG2	2.53	0.43
1:B:196:LYS:HE2	1:B:196:LYS:HB3	1.82	0.43
1:A:146:ASN:H	1:A:149:TYR:HB3	1.84	0.42
1:A:275:LYS:HE3	1:B:175:ALA:O	2.20	0.42
1:B:315:LEU:HD22	1:B:315:LEU:N	2.33	0.42
1:A:265:GLY:C	1:A:267:GLY:N	2.78	0.42
1:B:138:ILE:HD12	1:B:138:ILE:HA	1.84	0.42
1:A:184:ASP:CG	1:A:185:GLY:H	2.27	0.42
1:A:307:ILE:O	1:A:311:LEU:CD2	2.65	0.42
1:B:163:MET:SD	1:B:290:MET:CE	3.07	0.42
1:B:184:ASP:OD2	1:B:185:GLY:HA2	2.19	0.42
1:A:152:VAL:HG11	1:A:226:TRP:CH2	2.55	0.42
1:A:289:THR:CG2	1:B:165:GLN:HG2	2.47	0.42
1:B:151:LYS:HD2	1:B:154:GLU:OE1	2.19	0.42
1:B:152:VAL:HG13	1:B:212:TYR:CD2	2.54	0.41
1:A:307:ILE:HD11	1:B:308:PHE:HE2	1.85	0.41
1:B:286:GLU:OE1	1:B:286:GLU:HA	2.19	0.41
1:B:151:LYS:HD2	1:B:151:LYS:HA	1.87	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:315:LEU:HD13	1:B:315:LEU:HA	1.52	0.41
1:A:206:TYR:HB2	1:A:250:MET:HE2	2.02	0.41
1:A:197:LYS:HE2	1:A:197:LYS:HB2	1.58	0.41
1:B:186:ASN:HD22	1:B:186:ASN:HA	1.64	0.40
1:B:193:ASN:H	1:B:267:GLY:HA3	1.86	0.40
1:A:221:GLU:H	1:A:221:GLU:CD	2.29	0.40

All (1) 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)
1:B:216:ASN:OD1	1:B:263:LEU:O[4_455]	2.08	0.12

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	169/200 (84%)	159 (94%)	6 (4%)	4 (2%)	4	2
1	B	185/200 (92%)	172 (93%)	12 (6%)	1 (0%)	24	22
All	All	354/400 (88%)	331 (94%)	18 (5%)	5 (1%)	9	5

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	232	GLY
1	A	216	ASN
1	A	264	GLY
1	B	181	GLY
1	A	265	GLY

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	147/174 (84%)	130 (88%)	17 (12%)	5	3
1	B	161/174 (92%)	133 (83%)	28 (17%)	2	1
All	All	308/348 (88%)	263 (85%)	45 (15%)	3	1

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	144	ASP
1	A	145	ILE
1	A	146	ASN
1	A	147	GLU
1	A	151	LYS
1	A	152	VAL
1	A	157	VAL
1	A	184	ASP
1	A	196	LYS
1	A	203	LYS
1	A	207	LYS
1	A	216	ASN
1	A	222	GLN
1	A	233	THR
1	A	234	ILE
1	A	258	LYS
1	A	306	SER
1	B	135	TRP
1	B	137	LYS
1	B	138	ILE
1	B	151	LYS
1	B	183	ASN
1	B	186	ASN
1	B	188	VAL
1	B	203	LYS
1	B	207	LYS
1	B	221	GLU

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Mol	Chain	Res	Type
1	B	225	LYS
1	B	234	ILE
1	B	236	LYS
1	B	238	SER
1	B	239	GLN
1	B	240	LYS
1	B	250	MET
1	B	258	LYS
1	B	266	ASN
1	B	271	LEU
1	B	273	ASN
1	B	287	ASP
1	B	289	THR
1	B	299	GLN
1	B	300	LYS
1	B	315	LEU
1	B	317	SER
1	B	319	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	148	GLN
1	A	162	GLN
1	A	183	ASN
1	A	216	ASN
1	A	295	GLN
1	B	146	ASN
1	B	186	ASN
1	B	241	ASN
1	B	262	ASN
1	B	273	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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](#)

There are no ligands in this entry.

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	171/200 (85%)	0.04	14 (8%) 17 18	17, 37, 84, 92	0
1	B	187/200 (93%)	0.29	13 (6%) 22 24	20, 44, 77, 98	0
All	All	358/400 (89%)	0.17	27 (7%) 20 21	17, 41, 82, 98	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	182	GLY	6.8
1	B	183	ASN	4.0
1	B	265	GLY	4.0
1	A	314	VAL	3.9
1	B	267	GLY	3.9
1	B	264	GLY	3.8
1	B	135	TRP	3.5
1	A	146	ASN	3.3
1	A	311	LEU	3.3
1	A	185	GLY	3.1
1	A	266	ASN	2.9
1	B	185	GLY	2.6
1	A	181	GLY	2.6
1	A	184	ASP	2.6
1	B	134	LEU	2.6
1	A	215	ASN	2.4
1	B	318	THR	2.4
1	B	184	ASP	2.3
1	A	312	VAL	2.3
1	B	263	LEU	2.3
1	B	244	TYR	2.3
1	B	233	THR	2.2
1	B	262	ASN	2.2
1	A	310	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	145	ILE	2.1
1	A	263	LEU	2.1
1	A	231	GLY	2.1

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

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.