



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:44 AM UTC

PDB ID : 2FJY / pdb_00002fjy
Title : Crystal Structure of B-form Bombyx mori Pheromone Binding Protein
Authors : Lautenschlager, C.; Leal, W.S.; Clardy, J.
Deposited on : 2006-01-03
Resolution : 2.30 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

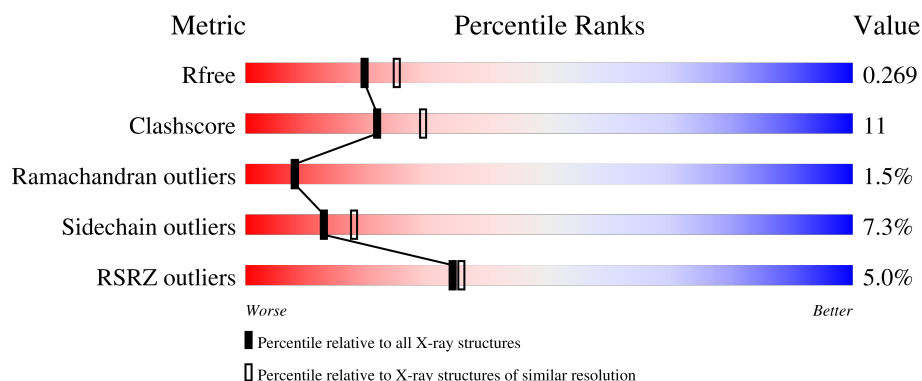
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	142	3% 54% 31% 11% ..
1	B	142	7% 51% 32% 13% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2261 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 Pheromone-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	136	Total	C	N	O	S	0	0	0
			1059	664	176	207	12			
1	B	142	Total	C	N	O	S	0	0	0
			1107	693	184	217	13			

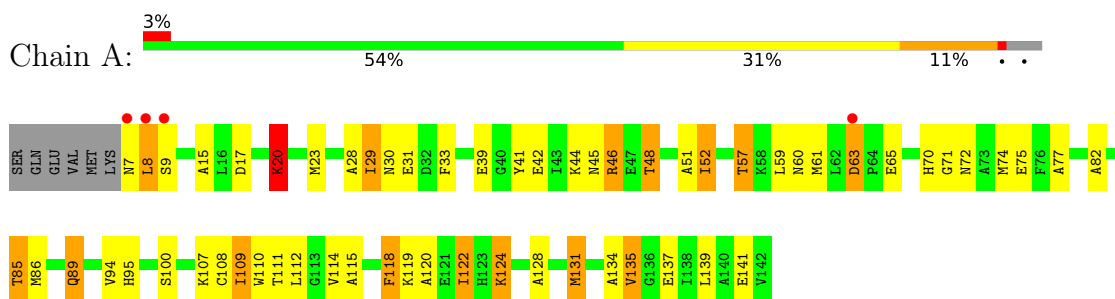
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	48	Total	O	0	0
			48	48		
2	B	47	Total	O	0	0
			47	47		

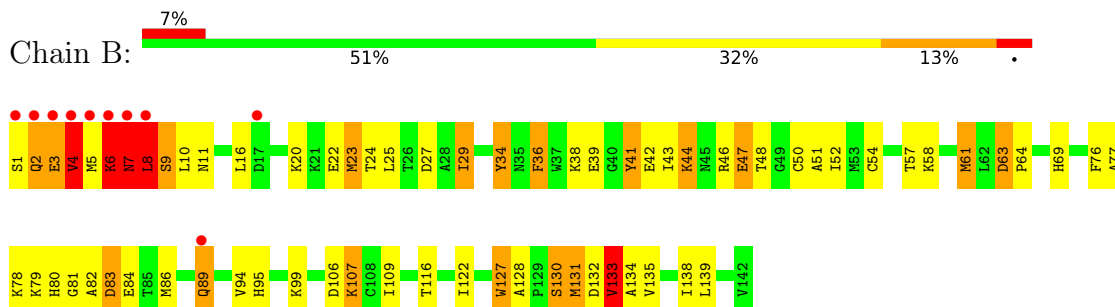
3 Residue-property plots [i](#)

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: Pheromone-binding protein



- Molecule 1: Pheromone-binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.24Å 70.79Å 75.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.00 – 2.30 43.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	89.9 (43.00-2.30) 89.9 (43.00-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.75 (at 2.29Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.220 , 0.270 0.216 , 0.269	Depositor DCC
R_{free} test set	605 reflections (4.50%)	wwPDB-VP
Wilson B-factor (Å ²)	20.8	Xtriage
Anisotropy	0.157	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 26.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2261	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.69 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1045e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	2.53	65/1079 (6.0%)	1.55	6/1454 (0.4%)
1	B	2.65	64/1127 (5.7%)	1.58	6/1517 (0.4%)
All	All	2.60	129/2206 (5.8%)	1.57	12/2971 (0.4%)

All (129) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	61	MET	SD-CE	-19.54	1.30	1.79
1	A	63	ASP	CB-CG	12.13	1.82	1.52
1	B	23	MET	SD-CE	-11.22	1.51	1.79
1	B	63	ASP	CB-CG	11.15	1.79	1.52
1	A	128	ALA	C-O	-10.81	1.14	1.24
1	B	131	MET	SD-CE	10.47	2.05	1.79
1	B	77	ALA	CA-CB	10.45	1.69	1.53
1	B	4	VAL	CA-CB	10.02	1.68	1.54
1	A	95	HIS	C-O	-9.57	1.12	1.24
1	B	89	GLN	CD-NE2	9.32	1.52	1.33
1	B	8	LEU	CA-C	9.27	1.65	1.52
1	B	128	ALA	CA-CB	9.20	1.65	1.53
1	A	89	GLN	CG-CD	8.95	1.74	1.52
1	B	107	LYS	CE-NZ	8.75	1.75	1.49
1	B	95	HIS	N-CA	8.41	1.56	1.46
1	A	23	MET	CG-SD	-8.03	1.60	1.80
1	B	89	GLN	CG-CD	7.96	1.72	1.52
1	A	86	MET	CA-C	7.95	1.62	1.52
1	A	115	ALA	CA-CB	-7.82	1.41	1.53
1	B	94	VAL	CA-C	-7.59	1.43	1.52
1	A	111	THR	N-CA	-7.59	1.37	1.46
1	A	41	TYR	CA-C	-7.50	1.43	1.52
1	B	51	ALA	C-O	-7.40	1.15	1.24
1	B	107	LYS	CD-CE	7.28	1.74	1.52
1	A	82	ALA	CA-CB	-7.27	1.40	1.53
1	B	134	ALA	CA-CB	-7.25	1.42	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	135	VAL	CA-C	-7.23	1.43	1.52
1	A	77	ALA	N-CA	-7.18	1.37	1.46
1	A	39	GLU	C-O	7.18	1.32	1.23
1	B	109	ILE	C-O	-7.16	1.16	1.24
1	A	48	THR	C-O	-7.12	1.15	1.24
1	A	57	THR	N-CA	7.03	1.55	1.46
1	B	5	MET	CA-C	6.96	1.61	1.52
1	A	118	PHE	C-O	6.90	1.32	1.24
1	B	34	TYR	CA-C	-6.86	1.44	1.52
1	A	131	MET	SD-CE	6.86	1.96	1.79
1	A	131	MET	C-O	-6.85	1.16	1.24
1	B	78	LYS	CA-C	-6.79	1.44	1.52
1	B	29	ILE	CA-CB	-6.78	1.46	1.54
1	A	119	LYS	N-CA	6.77	1.54	1.46
1	B	7	ASN	N-CA	6.74	1.54	1.46
1	A	51	ALA	C-O	-6.73	1.16	1.24
1	A	59	LEU	CA-C	-6.70	1.43	1.52
1	B	7	ASN	C-O	6.67	1.32	1.24
1	A	61	MET	SD-CE	-6.67	1.62	1.79
1	A	122	ILE	CA-C	6.54	1.61	1.52
1	A	44	LYS	CB-CG	6.52	1.72	1.52
1	A	8	LEU	CA-C	6.47	1.61	1.53
1	B	83	ASP	N-CA	6.43	1.53	1.46
1	A	39	GLU	CA-CB	6.40	1.63	1.53
1	B	130	SER	C-O	6.39	1.31	1.23
1	B	24	THR	CA-CB	-6.38	1.42	1.53
1	A	71	GLY	C-O	-6.37	1.15	1.23
1	B	64	PRO	C-O	-6.37	1.15	1.24
1	B	47	GLU	C-O	-6.26	1.16	1.24
1	A	33	PHE	CA-C	-6.25	1.44	1.52
1	A	135	VAL	CA-CB	-6.20	1.46	1.54
1	A	72	ASN	CA-CB	6.18	1.64	1.53
1	A	112	LEU	C-O	6.17	1.31	1.24
1	A	107	LYS	CA-C	6.12	1.61	1.52
1	A	46	ARG	CG-CD	6.12	1.70	1.52
1	A	124	LYS	CE-NZ	6.07	1.67	1.49
1	B	6	LYS	N-CA	6.07	1.54	1.46
1	A	45	ASN	CA-C	6.07	1.60	1.52
1	A	137	GLU	C-O	6.06	1.31	1.24
1	A	60	ASN	CA-C	-6.06	1.44	1.52
1	B	36	PHE	CA-C	-6.05	1.44	1.52
1	B	79	LYS	N-CA	6.04	1.54	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	89	GLN	CD-NE2	6.03	1.46	1.33
1	B	84	GLU	CA-C	-5.99	1.44	1.52
1	B	127	TRP	C-O	5.94	1.31	1.23
1	A	28	ALA	CA-C	5.94	1.60	1.52
1	B	95	HIS	C-O	-5.92	1.17	1.24
1	B	109	ILE	N-CA	5.90	1.53	1.46
1	B	11	ASN	CA-CB	5.86	1.61	1.53
1	B	122	ILE	CA-CB	5.86	1.62	1.54
1	A	15	ALA	N-CA	5.86	1.53	1.46
1	B	79	LYS	CD-CE	5.85	1.70	1.52
1	B	2	GLN	N-CA	5.82	1.52	1.45
1	A	141	GLU	C-O	-5.81	1.16	1.24
1	B	86	MET	CA-C	5.80	1.60	1.52
1	B	6	LYS	CA-C	5.78	1.60	1.52
1	A	46	ARG	CZ-NH2	5.78	1.41	1.33
1	B	25	LEU	N-CA	-5.77	1.39	1.46
1	B	9	SER	N-CA	5.74	1.53	1.46
1	A	31	GLU	CD-OE1	5.74	1.36	1.25
1	B	82	ALA	CA-CB	-5.74	1.42	1.53
1	A	110	TRP	C-O	-5.74	1.17	1.24
1	B	39	GLU	CG-CD	5.73	1.66	1.52
1	B	57	THR	C-O	-5.71	1.16	1.24
1	A	85	THR	C-O	-5.70	1.17	1.24
1	A	61	MET	CG-SD	5.70	1.95	1.80
1	A	124	LYS	C-O	-5.69	1.16	1.24
1	A	114	VAL	CA-C	-5.68	1.46	1.52
1	B	44	LYS	CB-CG	5.63	1.69	1.52
1	B	131	MET	CG-SD	5.62	1.94	1.80
1	B	2	GLN	CB-CG	5.61	1.69	1.52
1	A	28	ALA	CA-CB	-5.59	1.44	1.53
1	A	114	VAL	CA-CB	-5.58	1.47	1.54
1	A	94	VAL	CA-C	-5.57	1.46	1.52
1	A	111	THR	CA-C	-5.55	1.45	1.52
1	A	134	ALA	CA-CB	-5.52	1.44	1.53
1	B	89	GLN	CD-OE1	5.49	1.33	1.23
1	B	135	VAL	CA-C	-5.48	1.46	1.52
1	A	74	MET	CG-SD	5.46	1.94	1.80
1	B	69	HIS	C-O	-5.40	1.17	1.23
1	B	39	GLU	CA-CB	5.38	1.62	1.53
1	A	70	HIS	C-O	-5.34	1.17	1.23
1	B	41	TYR	CZ-OH	5.32	1.49	1.38
1	B	10	LEU	C-O	-5.31	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	46	ARG	CD-NE	5.29	1.53	1.46
1	A	139	LEU	C-O	-5.29	1.17	1.24
1	B	138	ILE	C-N	-5.28	1.26	1.34
1	A	109	ILE	C-O	-5.26	1.18	1.24
1	B	89	GLN	N-CA	-5.25	1.40	1.46
1	A	20	LYS	CA-C	5.23	1.59	1.52
1	A	29	ILE	C-O	-5.22	1.16	1.24
1	A	9	SER	CA-C	5.21	1.59	1.52
1	A	120	ALA	CA-CB	5.21	1.61	1.53
1	B	106	ASP	C-O	-5.17	1.17	1.24
1	A	108	CYS	CA-C	-5.16	1.46	1.52
1	A	131	MET	N-CA	5.15	1.52	1.46
1	B	10	LEU	CA-CB	5.15	1.60	1.53
1	B	2	GLN	CA-C	5.10	1.58	1.52
1	A	51	ALA	CA-CB	5.09	1.61	1.53
1	B	99	LYS	CA-C	-5.09	1.46	1.52
1	B	116	THR	CA-C	5.07	1.59	1.52
1	B	22	GLU	CD-OE2	5.05	1.34	1.25
1	B	78	LYS	CD-CE	5.04	1.67	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	133	VAL	CB-CA-C	-10.73	98.24	111.97
1	B	46	ARG	NE-CZ-NH2	6.69	125.22	119.20
1	A	46	ARG	NE-CZ-NH2	6.46	125.02	119.20
1	B	81	GLY	N-CA-C	6.08	123.75	115.30
1	A	100	SER	N-CA-C	5.80	119.61	112.54
1	A	134	ALA	N-CA-C	5.53	116.99	111.07
1	A	65	GLU	N-CA-C	-5.50	105.94	113.30
1	B	134	ALA	N-CA-C	5.35	117.11	111.28
1	A	52	ILE	CA-CB-CG1	5.33	119.46	110.40
1	A	65	GLU	CB-CA-C	5.30	119.14	110.17
1	B	2	GLN	N-CA-C	5.17	118.15	110.14
1	B	139	LEU	O-C-N	5.05	128.13	122.22

There are no chirality outliers.

There are no planarity outliers.

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	1059	0	1018	16	0
1	B	1107	0	1071	30	0
2	A	48	0	0	1	0
2	B	47	0	0	3	0
All	All	2261	0	2089	46	0

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 11.

All (46) 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:124:LYS:CE	1:A:124:LYS:NZ	1.67	1.57
1:A:63:ASP:CB	1:A:63:ASP:CG	1.82	1.53
1:B:63:ASP:CB	1:B:63:ASP:CG	1.79	1.52
1:B:107:LYS:CE	1:B:107:LYS:NZ	1.75	1.47
1:B:131:MET:SD	1:B:131:MET:CE	2.05	1.43
1:B:130:SER:OG	1:B:133:VAL:HG23	1.64	0.96
1:B:7:ASN:O	1:B:8:LEU:HB2	1.79	0.80
1:B:7:ASN:OD1	1:B:131:MET:HE3	1.84	0.76
1:A:29:ILE:C	1:A:29:ILE:HD12	2.21	0.65
1:B:7:ASN:O	1:B:8:LEU:CB	2.48	0.61
1:B:127:TRP:HB2	2:B:150:HOH:O	2.03	0.58
1:B:58:LYS:HD3	1:B:61:MET:HE2	1.86	0.58
1:B:3:GLU:HG2	1:B:4:VAL:H	1.70	0.57
1:B:20:LYS:HE2	1:B:27:ASP:HA	1.87	0.57
1:B:23:MET:HE1	1:B:50:CYS:O	2.06	0.56
1:B:58:LYS:HB3	1:B:61:MET:HE2	1.88	0.55
1:B:38:LYS:HG3	2:B:175:HOH:O	2.06	0.55
1:B:80:HIS:HB3	1:B:133:VAL:HG22	1.89	0.54
1:A:46:ARG:NH1	2:A:174:HOH:O	2.40	0.54
1:A:57:THR:O	1:A:57:THR:HG22	2.07	0.53
1:B:130:SER:OG	1:B:133:VAL:CG2	2.47	0.53
1:A:57:THR:O	1:A:57:THR:CG2	2.56	0.53
1:B:58:LYS:HD3	1:B:61:MET:CE	2.39	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:LYS:HG3	1:B:29:ILE:HD11	1.93	0.51
1:A:7:ASN:OD1	1:A:131:MET:HE3	2.10	0.50
1:A:118:PHE:O	1:A:122:ILE:HG23	2.11	0.50
1:B:3:GLU:HG2	1:B:4:VAL:N	2.27	0.49
1:B:83:ASP:OD1	1:B:83:ASP:C	2.55	0.49
1:B:36:PHE:HA	1:B:43:ILE:HD12	1.93	0.49
1:B:23:MET:HE1	1:B:50:CYS:C	2.39	0.48
1:A:85:THR:O	1:A:89:GLN:HG2	2.14	0.47
1:A:20:LYS:HE2	1:A:30:ASN:HD21	1.80	0.47
1:B:8:LEU:HD11	1:B:34:TYR:HE2	1.80	0.46
1:A:20:LYS:HG3	1:A:29:ILE:HD11	1.96	0.46
1:B:9:SER:HB3	1:B:132:ASP:OD1	2.16	0.46
1:A:46:ARG:HG2	1:A:109:ILE:HG12	1.98	0.44
1:B:38:LYS:HD2	1:B:41:TYR:CZ	2.52	0.44
1:A:131:MET:O	1:A:135:VAL:HG23	2.17	0.44
1:B:48:THR:O	1:B:52:ILE:HG13	2.18	0.44
1:A:29:ILE:HD12	1:A:30:ASN:N	2.34	0.43
1:B:8:LEU:HD11	2:B:170:HOH:O	2.20	0.42
1:A:48:THR:O	1:A:52:ILE:HG12	2.19	0.42
1:A:7:ASN:OD1	1:A:131:MET:CE	2.69	0.41
1:B:23:MET:CE	1:B:54:CYS:HB2	2.51	0.40
1:B:16:LEU:HD22	1:B:29:ILE:HD12	2.03	0.40
1:B:3:GLU:O	1:B:4:VAL:C	2.65	0.40

There are no symmetry-related clashes.

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	134/142 (94%)	129 (96%)	5 (4%)	0	100	100
1	B	140/142 (99%)	128 (91%)	8 (6%)	4 (3%)	3	2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	274/284 (96%)	257 (94%)	13 (5%)	4 (2%)	8 8

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	6	LYS
1	B	8	LEU
1	B	7	ASN
1	B	4	VAL

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	114/120 (95%)	109 (96%)	5 (4%)	25 38
1	B	120/120 (100%)	108 (90%)	12 (10%)	7 9
All	All	234/240 (98%)	217 (93%)	17 (7%)	13 18

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LEU
1	A	17	ASP
1	A	20	LYS
1	A	42	GLU
1	A	75	GLU
1	B	1	SER
1	B	2	GLN
1	B	3	GLU
1	B	4	VAL
1	B	6	LYS
1	B	8	LEU
1	B	42	GLU
1	B	44	LYS
1	B	47	GLU

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Mol	Chain	Res	Type
1	B	76	PHE
1	B	89	GLN
1	B	133	VAL

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	67	ASN
1	A	69	HIS
1	B	2	GLN
1	B	60	ASN
1	B	67	ASN

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](#)

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	136/142 (95%)	-0.12	4 (2%) 53 56	14, 23, 42, 108	0
1	B	142/142 (100%)	0.27	10 (7%) 22 24	14, 25, 67, 109	0
All	All	278/284 (97%)	0.08	14 (5%) 34 35	14, 24, 43, 109	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	SER	9.9
1	B	4	VAL	8.2
1	B	2	GLN	8.0
1	A	8	LEU	4.5
1	B	8	LEU	4.4
1	B	3	GLU	4.2
1	A	7	ASN	3.6
1	B	5	MET	3.3
1	B	7	ASN	3.2
1	B	6	LYS	3.0
1	A	9	SER	2.8
1	B	89	GLN	2.4
1	A	63	ASP	2.3
1	B	17	ASP	2.0

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.