



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:08 AM UTC

PDB ID : 1X9Y / pdb_00001x9y
Title : The prostaphopain B structure
Authors : Filipek, R.; Szczepanowski, R.; Sabat, A.; Potempa, J.; Bochtler, M.
Deposited on : 2004-08-24
Resolution : 2.50 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

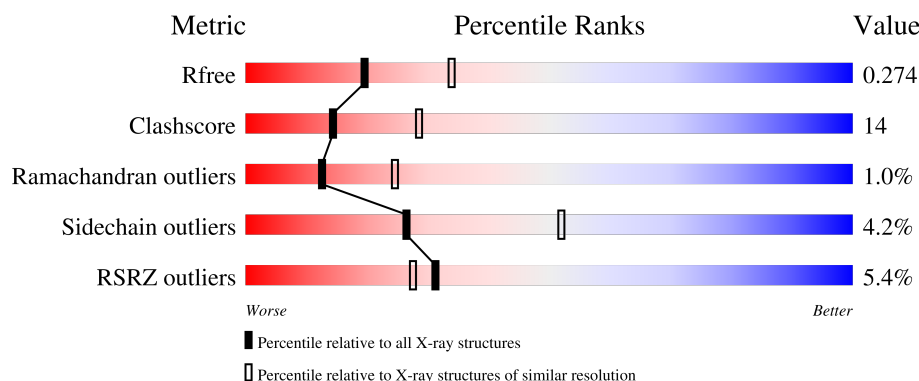
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

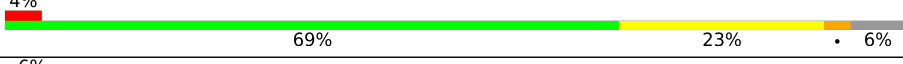
The reported resolution of this entry is 2.50 Å.

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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	367	
1	B	367	
1	C	367	
1	D	367	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11355 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 cysteine proteinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	346	Total	C	N	O	S	0	0	0
			2778	1756	463	551	8			
1	B	346	Total	C	N	O	S	0	0	0
			2778	1756	463	551	8			
1	C	346	Total	C	N	O	S	0	0	0
			2778	1756	463	551	8			
1	D	346	Total	C	N	O	S	0	0	0
			2778	1756	463	551	8			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	35	GLY	-	cloning artifact	UNP Q70UQ8
A	36	SER	-	cloning artifact	UNP Q70UQ8
A	394	LEU	-	cloning artifact	UNP Q70UQ8
A	395	GLU	-	cloning artifact	UNP Q70UQ8
A	396	VAL	-	cloning artifact	UNP Q70UQ8
A	397	PRO	-	cloning artifact	UNP Q70UQ8
A	398	ILE	-	cloning artifact	UNP Q70UQ8
A	399	HIS	-	cloning artifact	UNP Q70UQ8
A	400	ARG	-	cloning artifact	UNP Q70UQ8
A	401	ASP	-	cloning artifact	UNP Q70UQ8
B	35	GLY	-	cloning artifact	UNP Q70UQ8
B	36	SER	-	cloning artifact	UNP Q70UQ8
B	394	LEU	-	cloning artifact	UNP Q70UQ8
B	395	GLU	-	cloning artifact	UNP Q70UQ8
B	396	VAL	-	cloning artifact	UNP Q70UQ8
B	397	PRO	-	cloning artifact	UNP Q70UQ8
B	398	ILE	-	cloning artifact	UNP Q70UQ8
B	399	HIS	-	cloning artifact	UNP Q70UQ8
B	400	ARG	-	cloning artifact	UNP Q70UQ8
B	401	ASP	-	cloning artifact	UNP Q70UQ8
C	35	GLY	-	cloning artifact	UNP Q70UQ8

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Chain	Residue	Modelled	Actual	Comment	Reference
C	36	SER	-	cloning artifact	UNP Q70UQ8
C	394	LEU	-	cloning artifact	UNP Q70UQ8
C	395	GLU	-	cloning artifact	UNP Q70UQ8
C	396	VAL	-	cloning artifact	UNP Q70UQ8
C	397	PRO	-	cloning artifact	UNP Q70UQ8
C	398	ILE	-	cloning artifact	UNP Q70UQ8
C	399	HIS	-	cloning artifact	UNP Q70UQ8
C	400	ARG	-	cloning artifact	UNP Q70UQ8
C	401	ASP	-	cloning artifact	UNP Q70UQ8
D	35	GLY	-	cloning artifact	UNP Q70UQ8
D	36	SER	-	cloning artifact	UNP Q70UQ8
D	394	LEU	-	cloning artifact	UNP Q70UQ8
D	395	GLU	-	cloning artifact	UNP Q70UQ8
D	396	VAL	-	cloning artifact	UNP Q70UQ8
D	397	PRO	-	cloning artifact	UNP Q70UQ8
D	398	ILE	-	cloning artifact	UNP Q70UQ8
D	399	HIS	-	cloning artifact	UNP Q70UQ8
D	400	ARG	-	cloning artifact	UNP Q70UQ8
D	401	ASP	-	cloning artifact	UNP Q70UQ8

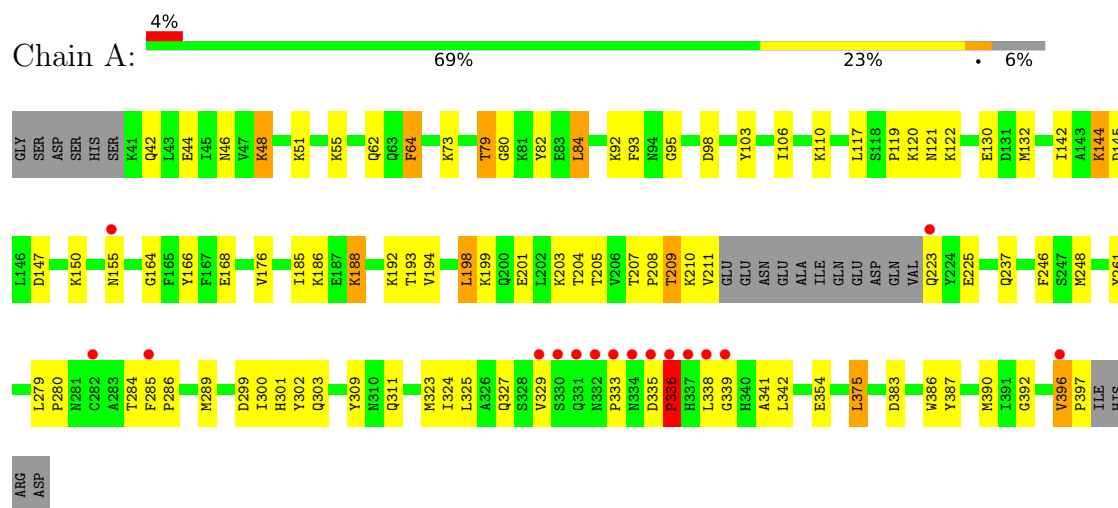
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	68	Total O 68 68	0	0
2	B	58	Total O 58 58	0	0
2	C	49	Total O 49 49	0	0
2	D	68	Total O 68 68	0	0

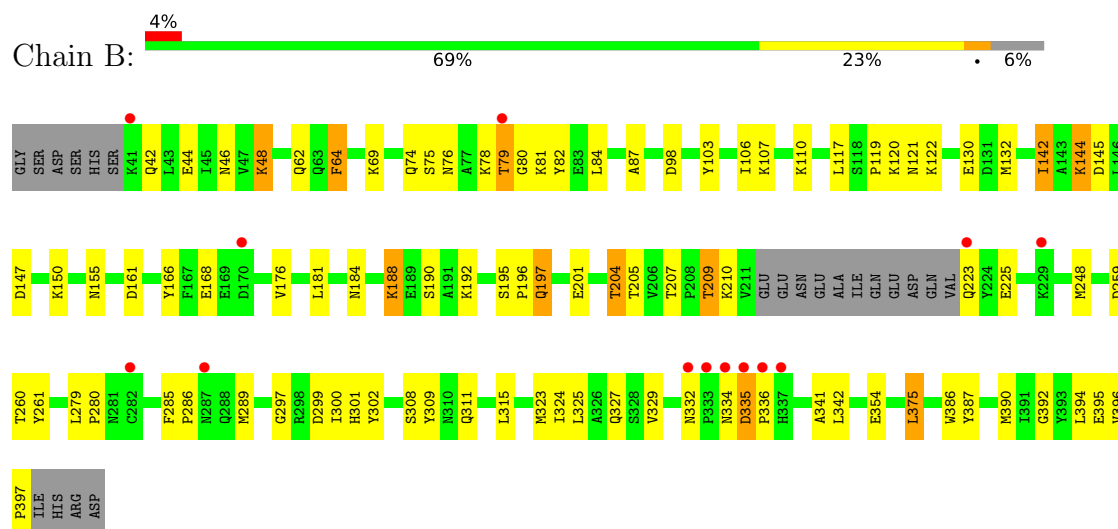
3 Residue-property plots [i](#)

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: cysteine proteinase

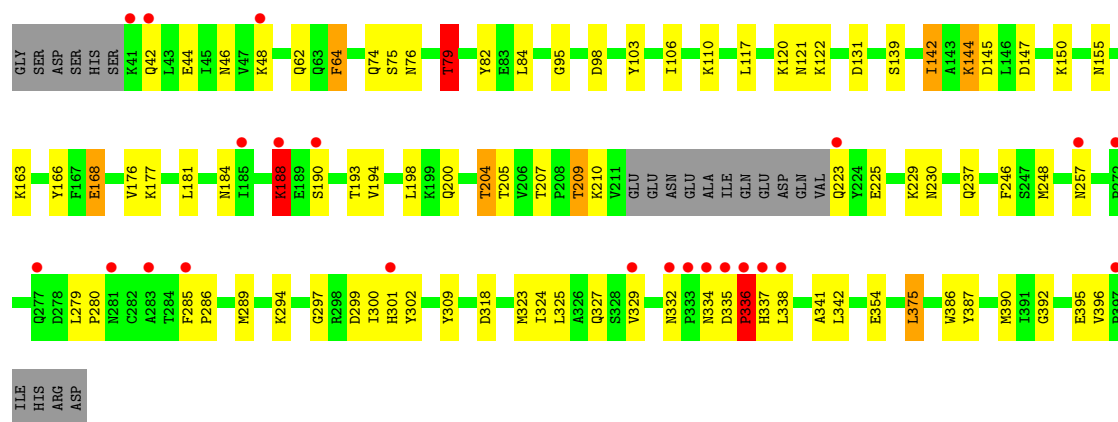


• Molecule 1: cysteine proteinase

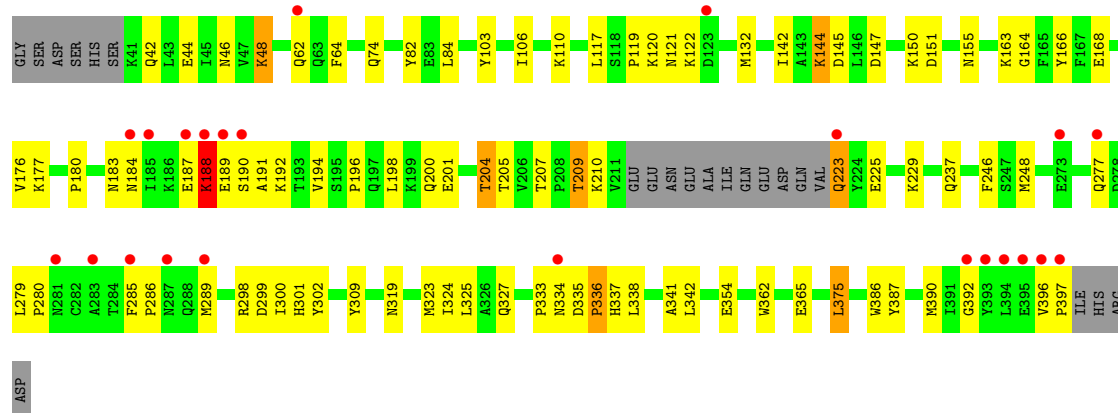


• Molecule 1: cysteine proteinase





● Molecule 1: cysteine proteinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	84.62Å 104.92Å 173.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.86 – 2.50 19.86 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.1 (19.86-2.50) 99.0 (19.86-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 2.50Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.244 , 0.276 0.242 , 0.274	Depositor DCC
R_{free} test set	2730 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	36.5	Xtriage
Anisotropy	0.721	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 36.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11355	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.07% 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	0.46	0/2833	0.89	5/3831 (0.1%)
1	B	0.45	0/2833	0.91	4/3831 (0.1%)
1	C	0.44	0/2833	0.89	3/3831 (0.1%)
1	D	0.45	0/2833	0.89	5/3831 (0.1%)
All	All	0.45	0/11332	0.90	17/15324 (0.1%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	396	VAL	N-CA-C	8.64	116.70	108.15
1	A	396	VAL	N-CA-C	6.34	114.80	107.89
1	B	196	PRO	N-CA-C	-6.15	106.17	113.86
1	B	64	PHE	N-CA-C	6.11	118.44	111.11
1	C	64	PHE	N-CA-C	5.99	118.30	111.11
1	A	120	LYS	N-CA-C	5.89	117.50	111.14
1	D	188	LYS	N-CA-C	5.77	117.61	108.67
1	A	164	GLY	N-CA-C	5.75	117.74	110.38
1	A	64	PHE	N-CA-C	5.63	118.14	111.33
1	D	64	PHE	N-CA-C	5.62	117.85	111.11
1	D	120	LYS	N-CA-C	5.48	117.06	111.14
1	C	120	LYS	N-CA-C	5.48	117.06	111.14
1	D	164	GLY	N-CA-C	5.46	117.37	110.38
1	D	191	ALA	N-CA-C	5.42	117.56	109.59
1	B	120	LYS	N-CA-C	5.21	116.77	111.14
1	B	396	VAL	N-CA-C	-5.21	103.84	108.95
1	A	198	LEU	N-CA-C	-5.05	105.68	111.14

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	2778	0	2724	83	0
1	B	2778	0	2724	88	0
1	C	2778	0	2724	84	0
1	D	2778	0	2724	77	0
2	A	68	0	0	10	0
2	B	58	0	0	6	0
2	C	49	0	0	6	0
2	D	68	0	0	7	0
All	All	11355	0	10896	310	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 14.

All (310) 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:195:SER:HB3	1:B:197:GLN:HG3	1.31	1.09
1:B:308:SER:HB3	1:C:257:ASN:CG	1.86	0.99
1:C:122:LYS:HE2	1:C:329:VAL:HG13	1.46	0.96
1:A:204:THR:HG22	1:A:205:THR:H	1.33	0.94
1:B:332:ASN:HD22	1:B:335:ASP:HB3	1.34	0.92
1:B:79:THR:HG22	1:B:80:GLY:H	1.37	0.89
1:D:204:THR:HG22	1:D:205:THR:H	1.38	0.89
1:C:336:PRO:HB3	1:C:387:TYR:CE1	2.07	0.88
1:A:336:PRO:HD2	1:D:397:PRO:HB3	1.57	0.87
1:D:207:THR:OG1	1:D:210:LYS:HB2	1.76	0.86
1:B:204:THR:HG22	1:B:205:THR:H	1.39	0.85
1:C:294:LYS:HD3	2:C:435:HOH:O	1.75	0.85
1:B:207:THR:OG1	1:B:210:LYS:HB2	1.75	0.85
1:D:289:MET:SD	1:D:323:MET:HE1	2.16	0.84
1:C:207:THR:OG1	1:C:210:LYS:HB2	1.77	0.84
1:A:289:MET:SD	1:A:323:MET:HE1	2.18	0.83
1:A:207:THR:OG1	1:A:210:LYS:HB2	1.77	0.83
1:B:315:LEU:HD21	1:C:229:LYS:HE3	1.64	0.80
1:C:204:THR:HG22	1:C:205:THR:H	1.46	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:289:MET:SD	1:C:323:MET:HE1	2.21	0.80
1:A:302:TYR:HB2	1:A:390:MET:HE1	1.63	0.80
1:C:144:LYS:HE3	2:C:434:HOH:O	1.83	0.78
1:B:289:MET:SD	1:B:323:MET:HE1	2.24	0.77
1:C:302:TYR:HB2	1:C:390:MET:HE1	1.65	0.77
1:B:302:TYR:HB2	1:B:390:MET:HE1	1.66	0.77
1:D:302:TYR:HB2	1:D:390:MET:HE1	1.66	0.76
1:A:188:LYS:N	1:A:188:LYS:HE3	2.02	0.75
1:D:48:LYS:HE3	1:D:201:GLU:O	1.86	0.75
1:B:46:ASN:HD21	1:B:155:ASN:ND2	1.85	0.75
1:B:308:SER:CB	1:C:257:ASN:OD1	2.34	0.75
1:C:131:ASP:HA	1:D:277:GLN:HE22	1.52	0.74
1:A:79:THR:HG22	1:A:80:GLY:N	1.99	0.74
1:B:308:SER:CB	1:C:257:ASN:CG	2.61	0.73
1:A:122:LYS:NZ	1:A:333:PRO:HB3	2.03	0.73
1:D:46:ASN:HD21	1:D:155:ASN:ND2	1.86	0.73
1:A:79:THR:HG22	1:A:80:GLY:H	1.54	0.73
1:C:46:ASN:HD21	1:C:155:ASN:ND2	1.86	0.72
1:D:188:LYS:HD2	1:D:188:LYS:N	2.03	0.72
1:B:79:THR:HG22	1:B:80:GLY:N	2.06	0.71
1:A:144:LYS:HE2	1:A:144:LYS:HA	1.72	0.71
1:D:144:LYS:HA	1:D:144:LYS:HE2	1.73	0.71
1:B:144:LYS:HA	1:B:144:LYS:HE2	1.72	0.70
1:D:196:PRO:O	1:D:200:GLN:HG2	1.91	0.70
1:B:375:LEU:HB3	2:B:423:HOH:O	1.91	0.70
1:D:177:LYS:HE2	2:D:412:HOH:O	1.89	0.70
1:D:286:PRO:N	1:D:289:MET:HE3	2.06	0.69
1:A:44:GLU:HG2	1:A:207:THR:HG22	1.73	0.69
1:C:144:LYS:HE2	1:C:144:LYS:HA	1.74	0.69
1:A:329:VAL:HG23	1:A:383:ASP:O	1.92	0.69
1:B:48:LYS:HB2	2:B:422:HOH:O	1.93	0.69
1:A:46:ASN:HD21	1:A:155:ASN:ND2	1.91	0.69
1:B:332:ASN:HD22	1:B:335:ASP:CB	2.06	0.69
1:C:181:LEU:HB2	1:C:184:ASN:ND2	2.08	0.68
1:B:207:THR:HG1	1:B:210:LYS:HB2	1.56	0.68
1:B:308:SER:HB3	1:C:257:ASN:OD1	1.92	0.68
1:C:44:GLU:HG2	1:C:207:THR:HG22	1.76	0.68
1:C:163:LYS:HE3	1:C:188:LYS:O	1.93	0.67
1:B:188:LYS:HE3	1:B:188:LYS:N	2.08	0.67
1:B:315:LEU:CD2	1:C:229:LYS:HE3	2.24	0.67
1:C:286:PRO:N	1:C:289:MET:HE3	2.09	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:324:ILE:HG12	1:C:386:TRP:CZ2	2.30	0.67
1:D:44:GLU:HG2	1:D:207:THR:HG22	1.77	0.67
1:D:324:ILE:HG12	1:D:386:TRP:CZ2	2.29	0.66
1:A:286:PRO:N	1:A:289:MET:HE3	2.10	0.66
1:B:286:PRO:N	1:B:289:MET:HE3	2.11	0.66
1:B:44:GLU:HG2	1:B:207:THR:HG22	1.78	0.66
1:A:324:ILE:HG12	1:A:386:TRP:CZ2	2.30	0.66
1:D:207:THR:HG1	1:D:210:LYS:HB2	1.59	0.66
1:C:300:ILE:HD13	1:C:392:GLY:HA2	1.78	0.65
1:B:302:TYR:HD1	1:B:390:MET:CE	2.09	0.65
1:A:300:ILE:HD13	1:A:392:GLY:HA2	1.78	0.65
1:C:188:LYS:HE3	1:C:188:LYS:N	2.12	0.64
1:B:324:ILE:HG12	1:B:386:TRP:CZ2	2.33	0.64
1:A:194:VAL:HG13	1:A:198:LEU:HD23	1.77	0.64
1:D:300:ILE:HD13	1:D:392:GLY:HA2	1.80	0.64
1:D:163:LYS:HE3	1:D:189:GLU:HB2	1.79	0.63
1:A:166:TYR:OH	1:A:192:LYS:NZ	2.32	0.63
1:A:302:TYR:HD1	1:A:390:MET:CE	2.12	0.62
1:B:300:ILE:HD13	1:B:392:GLY:HA2	1.79	0.62
1:A:208:PRO:HD2	2:A:426:HOH:O	1.99	0.62
1:C:334:ASN:C	1:C:336:PRO:HD3	2.25	0.62
1:A:48:LYS:HE3	1:A:201:GLU:O	2.00	0.62
1:A:95:GLY:HA2	1:A:198:LEU:HD13	1.82	0.61
1:B:46:ASN:HD21	1:B:155:ASN:HD22	1.46	0.61
1:B:78:LYS:O	1:B:107:LYS:NZ	2.30	0.61
1:C:302:TYR:HD1	1:C:390:MET:CE	2.14	0.61
1:B:74:GLN:NE2	1:B:225:GLU:HG3	2.16	0.61
1:B:161:ASP:OD2	1:B:192:LYS:NZ	2.34	0.60
1:C:46:ASN:HD21	1:C:155:ASN:HD22	1.49	0.59
1:D:74:GLN:NE2	1:D:225:GLU:HG3	2.17	0.59
1:C:84:LEU:HD22	1:C:103:TYR:HB3	1.84	0.59
1:A:79:THR:CG2	1:A:80:GLY:H	2.15	0.59
1:C:299:ASP:OD2	1:C:301:HIS:NE2	2.31	0.59
1:D:302:TYR:HD1	1:D:390:MET:CE	2.16	0.59
1:A:79:THR:CG2	1:A:80:GLY:N	2.66	0.59
1:C:188:LYS:H	1:C:188:LYS:CD	2.15	0.59
1:D:248:MET:HE2	1:D:289:MET:HG2	1.85	0.59
1:D:299:ASP:OD2	1:D:301:HIS:NE2	2.33	0.59
1:D:46:ASN:HD21	1:D:155:ASN:HD22	1.50	0.58
1:D:335:ASP:N	1:D:336:PRO:HD3	2.19	0.58
1:A:84:LEU:HD22	1:A:103:TYR:HB3	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:130:GLU:HG2	2:B:434:HOH:O	2.04	0.58
1:C:188:LYS:H	1:C:188:LYS:HD2	1.69	0.57
1:A:302:TYR:CB	1:A:390:MET:HE1	2.33	0.57
1:D:84:LEU:HD22	1:D:103:TYR:HB3	1.84	0.57
1:A:204:THR:HG22	1:A:205:THR:N	2.12	0.57
1:B:302:TYR:CB	1:B:390:MET:HE1	2.33	0.57
1:B:84:LEU:HD22	1:B:103:TYR:HB3	1.86	0.56
1:B:335:ASP:CG	1:B:335:ASP:O	2.48	0.56
1:C:42:GLN:OE1	1:C:209:THR:HG21	2.06	0.56
1:D:42:GLN:OE1	1:D:209:THR:HG21	2.06	0.56
1:B:285:PHE:C	1:B:289:MET:HE3	2.30	0.55
1:C:294:LYS:HB2	2:C:438:HOH:O	2.06	0.55
1:A:42:GLN:OE1	1:A:209:THR:HG21	2.05	0.55
1:D:285:PHE:C	1:D:289:MET:HE3	2.31	0.55
1:B:195:SER:HB3	1:B:197:GLN:CG	2.22	0.55
1:B:46:ASN:ND2	1:B:155:ASN:HD22	2.04	0.54
1:D:184:ASN:HB3	2:D:424:HOH:O	2.07	0.54
1:D:302:TYR:CB	1:D:390:MET:HE1	2.35	0.54
1:B:308:SER:OG	1:C:230:ASN:OD1	2.21	0.54
1:C:285:PHE:C	1:C:289:MET:HE3	2.32	0.54
1:C:302:TYR:CB	1:C:390:MET:HE1	2.34	0.54
1:B:311:GLN:HG3	1:C:229:LYS:HG3	1.89	0.54
1:A:248:MET:HE2	1:A:289:MET:HG2	1.89	0.54
1:B:42:GLN:OE1	1:B:209:THR:HG21	2.08	0.54
1:A:46:ASN:HD21	1:A:155:ASN:HD22	1.54	0.54
1:D:289:MET:HE1	1:D:338:LEU:HD22	1.90	0.54
1:C:248:MET:HE2	1:C:289:MET:HG2	1.89	0.54
1:A:122:LYS:HE2	1:A:329:VAL:HG13	1.90	0.53
1:B:48:LYS:HD2	2:B:422:HOH:O	2.09	0.53
1:C:177:LYS:HE2	2:C:419:HOH:O	2.08	0.53
1:D:46:ASN:ND2	1:D:155:ASN:HD22	2.06	0.53
1:D:334:ASN:C	1:D:336:PRO:HD3	2.33	0.53
1:B:181:LEU:HB2	1:B:184:ASN:ND2	2.23	0.53
1:A:303:GLN:HG3	1:D:319:ASN:HB2	1.91	0.53
1:A:207:THR:HG1	1:A:210:LYS:HB2	1.72	0.53
1:A:285:PHE:C	1:A:289:MET:HE3	2.34	0.53
1:D:189:GLU:HG3	1:D:192:LYS:NZ	2.24	0.53
1:A:311:GLN:HG3	1:D:229:LYS:HG3	1.90	0.52
1:A:211:VAL:HG12	2:A:418:HOH:O	2.09	0.52
1:B:308:SER:CA	1:C:257:ASN:OD1	2.58	0.52
1:B:327:GLN:OE1	1:B:336:PRO:HA	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:46:ASN:ND2	1:C:155:ASN:HD22	2.06	0.52
1:C:95:GLY:HA2	1:C:198:LEU:HD13	1.91	0.52
1:A:130:GLU:OE2	1:B:260:THR:OG1	2.21	0.52
1:C:297:GLY:HA2	1:C:395:GLU:CD	2.34	0.52
1:D:325:LEU:HD22	1:D:338:LEU:HD13	1.92	0.51
1:D:327:GLN:HG2	1:D:387:TYR:CD2	2.45	0.51
1:A:73:LYS:NZ	1:A:225:GLU:HB3	2.26	0.51
1:B:327:GLN:HG2	1:B:387:TYR:CD2	2.46	0.51
1:A:336:PRO:HD2	1:D:397:PRO:CB	2.36	0.51
1:A:185:ILE:HG23	1:A:186:LYS:N	2.25	0.50
1:B:302:TYR:CD1	1:B:390:MET:CE	2.94	0.50
1:C:327:GLN:HG2	1:C:387:TYR:CD2	2.46	0.50
1:B:248:MET:HE2	1:B:289:MET:HG2	1.93	0.50
1:A:55:LYS:HE2	2:A:442:HOH:O	2.11	0.50
1:D:189:GLU:HG3	1:D:192:LYS:HZ3	1.77	0.50
1:B:79:THR:CG2	1:B:80:GLY:H	2.16	0.50
1:B:332:ASN:ND2	1:B:335:ASP:HB3	2.14	0.50
1:A:46:ASN:ND2	1:A:155:ASN:HD22	2.10	0.49
1:B:204:THR:HG22	1:B:205:THR:N	2.18	0.49
2:B:426:HOH:O	1:C:257:ASN:HB2	2.11	0.49
1:B:301:HIS:ND1	1:C:318:ASP:OD1	2.46	0.49
1:A:327:GLN:HG2	1:A:387:TYR:CD2	2.47	0.49
1:C:74:GLN:NE2	1:C:225:GLU:HG3	2.28	0.49
1:B:74:GLN:NE2	1:B:225:GLU:CG	2.75	0.49
1:C:302:TYR:HA	1:C:390:MET:HE2	1.95	0.49
1:B:74:GLN:HE21	1:B:225:GLU:HG3	1.77	0.49
1:D:145:ASP:CG	1:D:176:VAL:HG13	2.37	0.49
1:D:309:TYR:CD1	1:D:354:GLU:HB3	2.48	0.49
1:B:147:ASP:O	1:B:150:LYS:HG2	2.13	0.48
1:C:335:ASP:O	1:C:337:HIS:N	2.46	0.48
1:A:284:THR:O	1:A:338:LEU:HD23	2.13	0.48
1:D:302:TYR:HA	1:D:390:MET:HE2	1.95	0.48
1:D:223:GLN:NE2	2:D:460:HOH:O	2.33	0.48
1:A:302:TYR:HA	1:A:390:MET:HE2	1.94	0.48
1:B:308:SER:HA	1:C:257:ASN:OD1	2.12	0.48
1:D:204:THR:CG2	1:D:205:THR:H	2.15	0.48
1:D:122:LYS:NZ	1:D:333:PRO:HG3	2.29	0.48
1:D:147:ASP:O	1:D:150:LYS:HG2	2.14	0.48
1:A:211:VAL:HG11	2:A:423:HOH:O	2.14	0.48
1:A:51:LYS:HD2	1:B:259:ASP:OD1	2.14	0.47
1:C:309:TYR:CD1	1:C:354:GLU:HB3	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:329:VAL:HG12	1:A:329:VAL:O	2.15	0.47
1:D:204:THR:HG22	1:D:205:THR:N	2.18	0.47
1:D:337:HIS:HB3	2:D:451:HOH:O	2.13	0.47
1:B:334:ASN:C	1:B:336:PRO:HD3	2.40	0.47
1:C:145:ASP:CG	1:C:176:VAL:HG13	2.38	0.47
1:D:248:MET:CE	1:D:289:MET:HG2	2.45	0.47
1:D:299:ASP:CG	1:D:301:HIS:HE2	2.21	0.47
1:A:145:ASP:CG	1:A:176:VAL:HG13	2.40	0.47
1:B:308:SER:HB3	1:C:257:ASN:ND2	2.29	0.47
1:D:279:LEU:N	1:D:280:PRO:CD	2.78	0.47
1:C:279:LEU:N	1:C:280:PRO:CD	2.78	0.47
1:B:309:TYR:CD1	1:B:354:GLU:HB3	2.49	0.46
1:C:207:THR:HG1	1:C:210:LYS:HB2	1.79	0.46
1:D:365:GLU:HB3	2:D:459:HOH:O	2.15	0.46
1:A:144:LYS:HA	1:A:144:LYS:CE	2.43	0.46
1:A:261:TYR:HA	2:A:462:HOH:O	2.15	0.46
1:A:341:ALA:C	1:A:342:LEU:HD12	2.41	0.46
1:D:119:PRO:HG3	1:D:132:MET:SD	2.55	0.46
1:A:98:ASP:HB3	2:A:421:HOH:O	2.16	0.46
1:A:130:GLU:HB3	2:A:429:HOH:O	2.15	0.46
1:A:302:TYR:CD1	1:A:390:MET:CE	2.95	0.46
1:A:93:PHE:HD1	2:A:432:HOH:O	1.97	0.46
1:A:289:MET:SD	1:A:323:MET:CE	2.98	0.46
1:B:302:TYR:HA	1:B:390:MET:HE2	1.98	0.46
1:B:341:ALA:C	1:B:342:LEU:HD12	2.41	0.46
1:A:166:TYR:CZ	1:A:192:LYS:NZ	2.81	0.46
1:A:325:LEU:HD11	1:A:390:MET:HE3	1.98	0.46
1:A:194:VAL:HB	1:A:199:LYS:HE3	1.97	0.46
1:C:188:LYS:HE3	1:C:188:LYS:H	1.78	0.46
1:B:325:LEU:HD11	1:B:390:MET:HE3	1.97	0.46
1:B:145:ASP:CG	1:B:176:VAL:HG13	2.41	0.45
1:D:183:ASN:O	1:D:187:GLU:HG3	2.16	0.45
1:B:279:LEU:N	1:B:280:PRO:CD	2.79	0.45
1:C:279:LEU:HB3	1:C:280:PRO:HD3	1.98	0.45
1:C:302:TYR:CD1	1:C:390:MET:CE	2.98	0.45
1:C:341:ALA:C	1:C:342:LEU:HD12	2.42	0.45
1:A:342:LEU:HD12	1:A:342:LEU:N	2.31	0.45
1:C:248:MET:CE	1:C:289:MET:HG2	2.46	0.45
1:A:110:LYS:HA	1:A:150:LYS:HZ3	1.82	0.45
1:A:309:TYR:CD1	1:A:354:GLU:HB3	2.51	0.45
1:B:81:LYS:O	1:B:107:LYS:HA	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:289:MET:SD	1:C:323:MET:CE	3.00	0.45
1:C:299:ASP:CG	1:C:301:HIS:HE2	2.23	0.45
1:A:147:ASP:O	1:A:150:LYS:HG2	2.16	0.45
1:B:297:GLY:HA2	1:B:395:GLU:OE1	2.16	0.45
1:C:144:LYS:HA	1:C:144:LYS:CE	2.44	0.45
1:D:48:LYS:HD2	1:D:201:GLU:HB3	1.98	0.45
1:D:298:ARG:NE	2:D:415:HOH:O	2.49	0.45
1:C:79:THR:HG22	1:C:82:TYR:OH	2.17	0.45
1:A:396:VAL:HA	1:A:397:PRO:HD3	1.72	0.45
1:C:147:ASP:O	1:C:150:LYS:HG2	2.17	0.45
1:D:279:LEU:HB3	1:D:280:PRO:HD3	1.99	0.45
1:A:48:LYS:NZ	1:A:203:LYS:HE2	2.31	0.44
1:A:299:ASP:OD2	1:A:301:HIS:NE2	2.43	0.44
1:C:110:LYS:HA	1:C:150:LYS:HZ3	1.81	0.44
1:B:342:LEU:HD12	1:B:342:LEU:N	2.32	0.44
1:D:74:GLN:NE2	1:D:225:GLU:CG	2.81	0.44
1:A:279:LEU:HB3	1:A:280:PRO:HD3	1.99	0.44
1:D:289:MET:SD	1:D:323:MET:CE	2.99	0.44
1:A:73:LYS:HZ1	1:A:225:GLU:HB3	1.82	0.44
1:B:64:PHE:CG	1:B:84:LEU:HG	2.53	0.44
1:A:279:LEU:N	1:A:280:PRO:CD	2.80	0.44
1:A:335:ASP:N	1:A:336:PRO:HD3	2.33	0.44
1:A:130:GLU:HG3	1:B:261:TYR:CZ	2.53	0.43
1:B:122:LYS:HE2	1:B:329:VAL:HG13	1.99	0.43
1:C:188:LYS:H	1:C:188:LYS:CE	2.31	0.43
1:B:279:LEU:HB3	1:B:280:PRO:HD3	1.99	0.43
1:C:139:SER:HB2	2:C:447:HOH:O	2.18	0.43
1:D:144:LYS:HA	1:D:144:LYS:CE	2.46	0.43
1:D:325:LEU:HD11	1:D:390:MET:HE3	2.00	0.43
1:A:336:PRO:CD	1:D:397:PRO:HB3	2.38	0.43
1:C:342:LEU:HD12	1:C:342:LEU:N	2.33	0.43
1:D:151:ASP:HB2	2:D:461:HOH:O	2.18	0.43
1:C:121:ASN:C	1:C:375:LEU:HD22	2.43	0.43
1:A:82:TYR:HA	1:A:106:ILE:O	2.19	0.43
1:C:110:LYS:HA	1:C:150:LYS:NZ	2.34	0.43
1:B:69:LYS:HA	1:B:69:LYS:HD2	1.90	0.43
1:D:342:LEU:HD12	1:D:342:LEU:N	2.34	0.43
1:B:48:LYS:HE3	1:B:201:GLU:O	2.19	0.43
1:C:332:ASN:HB2	1:C:335:ASP:HB3	2.01	0.43
1:D:341:ALA:C	1:D:342:LEU:HD12	2.44	0.43
1:B:82:TYR:HA	1:B:106:ILE:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:168:GLU:OE2	1:C:194:VAL:HB	2.18	0.43
1:A:339:GLY:HA2	2:A:459:HOH:O	2.18	0.42
1:D:237:GLN:HB2	1:D:246:PHE:CD2	2.55	0.42
1:D:121:ASN:C	1:D:375:LEU:HD22	2.44	0.42
1:D:335:ASP:N	1:D:336:PRO:CD	2.82	0.42
1:A:237:GLN:HB2	1:A:246:PHE:CD2	2.55	0.42
1:A:248:MET:CE	1:A:289:MET:HG2	2.49	0.42
1:C:64:PHE:CG	1:C:84:LEU:HG	2.54	0.42
1:D:194:VAL:HG13	1:D:198:LEU:HD23	2.00	0.42
1:A:92:LYS:HB3	2:A:431:HOH:O	2.19	0.42
1:B:110:LYS:HA	1:B:150:LYS:HZ3	1.84	0.42
1:D:180:PRO:HD2	1:D:362:TRP:NE1	2.35	0.42
1:B:98:ASP:HB3	2:B:404:HOH:O	2.19	0.42
1:C:75:SER:O	1:C:76:ASN:C	2.63	0.42
1:B:119:PRO:HG3	1:B:132:MET:SD	2.60	0.42
1:C:82:TYR:HA	1:C:106:ILE:O	2.20	0.42
1:D:180:PRO:HD2	1:D:362:TRP:CD1	2.54	0.42
1:A:64:PHE:CG	1:A:84:LEU:HG	2.55	0.42
1:B:121:ASN:C	1:B:375:LEU:HD22	2.44	0.42
1:A:42:GLN:CD	1:A:209:THR:HG21	2.45	0.41
1:A:121:ASN:C	1:A:375:LEU:HD22	2.45	0.41
1:B:87:ALA:HB2	1:B:103:TYR:CE2	2.55	0.41
1:B:144:LYS:HA	1:B:144:LYS:CE	2.45	0.41
1:C:42:GLN:CD	1:C:209:THR:HG21	2.45	0.41
1:A:119:PRO:HG3	1:A:132:MET:SD	2.61	0.41
1:B:299:ASP:OD2	1:B:301:HIS:NE2	2.46	0.41
1:C:325:LEU:HD11	1:C:390:MET:HE3	2.02	0.41
1:B:289:MET:SD	1:B:323:MET:CE	3.03	0.41
1:C:302:TYR:HD1	1:C:390:MET:HE3	1.85	0.41
1:D:302:TYR:CD1	1:D:390:MET:CE	3.00	0.41
1:D:110:LYS:HA	1:D:150:LYS:HZ3	1.86	0.41
1:A:166:TYR:CD1	1:A:166:TYR:N	2.89	0.41
1:B:142:ILE:HG22	1:B:176:VAL:HG12	2.03	0.41
1:C:98:ASP:HB3	2:C:402:HOH:O	2.20	0.41
1:C:142:ILE:HG22	1:C:176:VAL:HG12	2.03	0.41
1:C:166:TYR:CD1	1:C:166:TYR:N	2.89	0.41
1:D:42:GLN:CD	1:D:209:THR:HG21	2.46	0.41
1:B:74:GLN:HE22	1:B:225:GLU:CG	2.34	0.40
1:D:82:TYR:HA	1:D:106:ILE:O	2.21	0.40
1:D:166:TYR:CD1	1:D:166:TYR:N	2.89	0.40
1:B:308:SER:CB	1:C:257:ASN:ND2	2.84	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:300:ILE:HD13	1:D:392:GLY:CA	2.49	0.40
1:B:75:SER:O	1:B:76:ASN:HB2	2.21	0.40
1:B:166:TYR:CD1	1:B:166:TYR:N	2.89	0.40
1:C:237:GLN:HB2	1:C:246:PHE:CD2	2.56	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	342/367 (93%)	323 (94%)	16 (5%)	3 (1%)	14	27
1	B	342/367 (93%)	321 (94%)	18 (5%)	3 (1%)	14	27
1	C	342/367 (93%)	321 (94%)	15 (4%)	6 (2%)	6	12
1	D	342/367 (93%)	322 (94%)	18 (5%)	2 (1%)	21	38
All	All	1368/1468 (93%)	1287 (94%)	67 (5%)	14 (1%)	12	24

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	188	LYS
1	A	336	PRO
1	B	190	SER
1	C	190	SER
1	C	336	PRO
1	B	79	THR
1	A	79	THR
1	C	79	THR
1	C	338	LEU
1	D	190	SER
1	A	142	ILE

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Mol	Chain	Res	Type
1	B	142	ILE
1	C	142	ILE
1	D	142	ILE

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	309/328 (94%)	297 (96%)	12 (4%)	28	55
1	B	309/328 (94%)	295 (96%)	14 (4%)	24	49
1	C	309/328 (94%)	295 (96%)	14 (4%)	24	49
1	D	309/328 (94%)	297 (96%)	12 (4%)	28	55
All	All	1236/1312 (94%)	1184 (96%)	52 (4%)	26	52

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	LYS
1	A	62	GLN
1	A	84	LEU
1	A	117	LEU
1	A	144	LYS
1	A	168	GLU
1	A	188	LYS
1	A	193	THR
1	A	209	THR
1	A	223	GLN
1	A	336	PRO
1	A	375	LEU
1	B	48	LYS
1	B	62	GLN
1	B	117	LEU
1	B	144	LYS
1	B	168	GLU
1	B	188	LYS

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Mol	Chain	Res	Type
1	B	197	GLN
1	B	204	THR
1	B	209	THR
1	B	223	GLN
1	B	335	ASP
1	B	375	LEU
1	B	394	LEU
1	B	397	PRO
1	C	48	LYS
1	C	62	GLN
1	C	79	THR
1	C	117	LEU
1	C	144	LYS
1	C	168	GLU
1	C	188	LYS
1	C	193	THR
1	C	200	GLN
1	C	204	THR
1	C	209	THR
1	C	223	GLN
1	C	336	PRO
1	C	375	LEU
1	D	48	LYS
1	D	62	GLN
1	D	117	LEU
1	D	144	LYS
1	D	168	GLU
1	D	188	LYS
1	D	204	THR
1	D	209	THR
1	D	223	GLN
1	D	336	PRO
1	D	375	LEU
1	D	396	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	GLN
1	A	126	ASN
1	A	155	ASN
1	A	223	GLN

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Mol	Chain	Res	Type
1	A	257	ASN
1	A	303	GLN
1	A	311	GLN
1	A	327	GLN
1	B	74	GLN
1	B	76	ASN
1	B	126	ASN
1	B	153	ASN
1	B	155	ASN
1	B	197	GLN
1	B	223	GLN
1	B	230	ASN
1	B	257	ASN
1	B	303	GLN
1	B	311	GLN
1	B	332	ASN
1	B	337	HIS
1	C	74	GLN
1	C	126	ASN
1	C	155	ASN
1	C	223	GLN
1	C	303	GLN
1	C	311	GLN
1	C	327	GLN
1	C	331	GLN
1	C	337	HIS
1	D	74	GLN
1	D	76	ASN
1	D	126	ASN
1	D	155	ASN
1	D	183	ASN
1	D	223	GLN
1	D	257	ASN
1	D	277	GLN
1	D	281	ASN
1	D	303	GLN
1	D	311	GLN
1	D	327	GLN
1	D	331	GLN
1	D	337	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](#)

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	346/367 (94%)	0.25	16 (4%) 37 33	19, 33, 64, 108	0
1	B	346/367 (94%)	0.25	13 (3%) 44 39	19, 34, 63, 100	0
1	C	346/367 (94%)	0.29	23 (6%) 24 21	20, 35, 65, 100	0
1	D	346/367 (94%)	0.28	23 (6%) 24 21	19, 35, 64, 94	0
All	All	1384/1468 (94%)	0.27	75 (5%) 31 27	19, 34, 64, 108	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	396	VAL	9.0
1	B	333	PRO	6.9
1	D	397	PRO	6.3
1	B	336	PRO	5.9
1	A	335	ASP	5.7
1	A	332	ASN	5.5
1	A	333	PRO	5.3
1	B	334	ASN	5.2
1	A	338	LEU	4.4
1	A	223	GLN	4.4
1	D	190	SER	4.4
1	A	337	HIS	4.3
1	D	392	GLY	4.3
1	A	334	ASN	3.9
1	D	189	GLU	3.9
1	A	336	PRO	3.9
1	C	336	PRO	3.7
1	D	289	MET	3.6
1	D	184	ASN	3.6
1	D	223	GLN	3.6
1	C	333	PRO	3.6

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Mol	Chain	Res	Type	RSRZ
1	D	188	LYS	3.4
1	C	334	ASN	3.3
1	B	335	ASP	3.2
1	D	273	GLU	3.1
1	C	332	ASN	3.1
1	A	155	ASN	3.0
1	C	257	ASN	3.0
1	D	185	ILE	3.0
1	C	185	ILE	2.9
1	C	283	ALA	2.9
1	C	223	GLN	2.9
1	C	190	SER	2.8
1	D	277	GLN	2.8
1	C	335	ASP	2.7
1	A	285	PHE	2.7
1	A	330	SER	2.7
1	D	393	TYR	2.7
1	C	337	HIS	2.7
1	B	79	THR	2.6
1	B	229	LYS	2.6
1	B	282	CYS	2.6
1	D	283	ALA	2.6
1	B	223	GLN	2.6
1	C	42	GLN	2.6
1	C	285	PHE	2.6
1	B	332	ASN	2.5
1	C	281	ASN	2.5
1	D	287	ASN	2.5
1	D	285	PHE	2.5
1	C	188	LYS	2.5
1	B	41	LYS	2.5
1	D	281	ASN	2.4
1	C	272	PRO	2.3
1	D	394	LEU	2.3
1	A	339	GLY	2.3
1	C	338	LEU	2.3
1	D	123	ASP	2.3
1	B	287	ASN	2.3
1	C	277	GLN	2.3
1	C	329	VAL	2.3
1	A	396	VAL	2.2
1	C	48	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	187	GLU	2.2
1	D	395	GLU	2.2
1	D	334	ASN	2.1
1	A	282	CYS	2.1
1	C	397	PRO	2.1
1	A	329	VAL	2.1
1	C	301	HIS	2.1
1	A	331	GLN	2.0
1	D	62	GLN	2.0
1	C	41	LYS	2.0
1	B	337	HIS	2.0
1	B	170	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.