



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:48 PM UTC

PDB ID : 6CP2 / pdb_00006cp2
Title : SidC in complex with UbcH7 Ub
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Deposited on : 2018-03-13
Resolution : 2.90 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

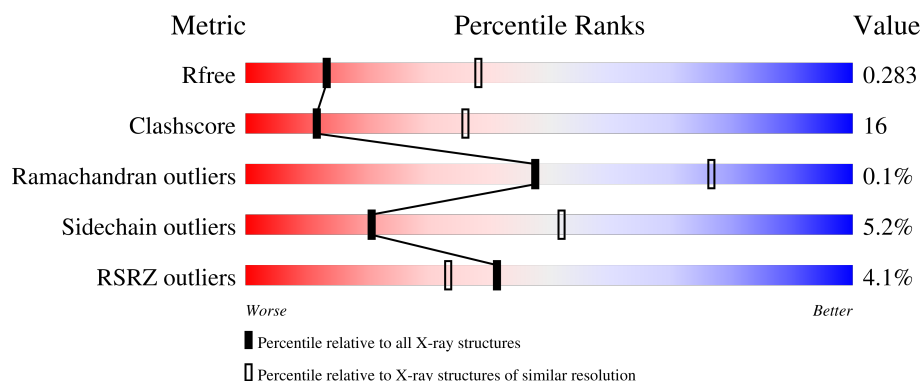
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	542	 4% 68% 22% 7%
2	B	154	 3% 64% 31% 2% 2%
3	C	78	 3% 63% 29% 6%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5957 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 SidC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	502	Total	C	N	O	S	0	0	0
			4065	2569	692	800	4			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	46	ALA	CYS	engineered mutation	UNP Q6RCR4

- Molecule 2 is a protein called Ubiquitin-conjugating enzyme E2 L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	153	Total	C	N	O	S	0	0	0
			1252	803	217	227	5			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	86	LYS	CYS	engineered mutation	UNP P68036

- Molecule 3 is a protein called Polyubiquitin-B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	78	Total	C	N	O	S	0	0	0
			611	383	107	120	1			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	SER	GLY	conflict	UNP P0CG47

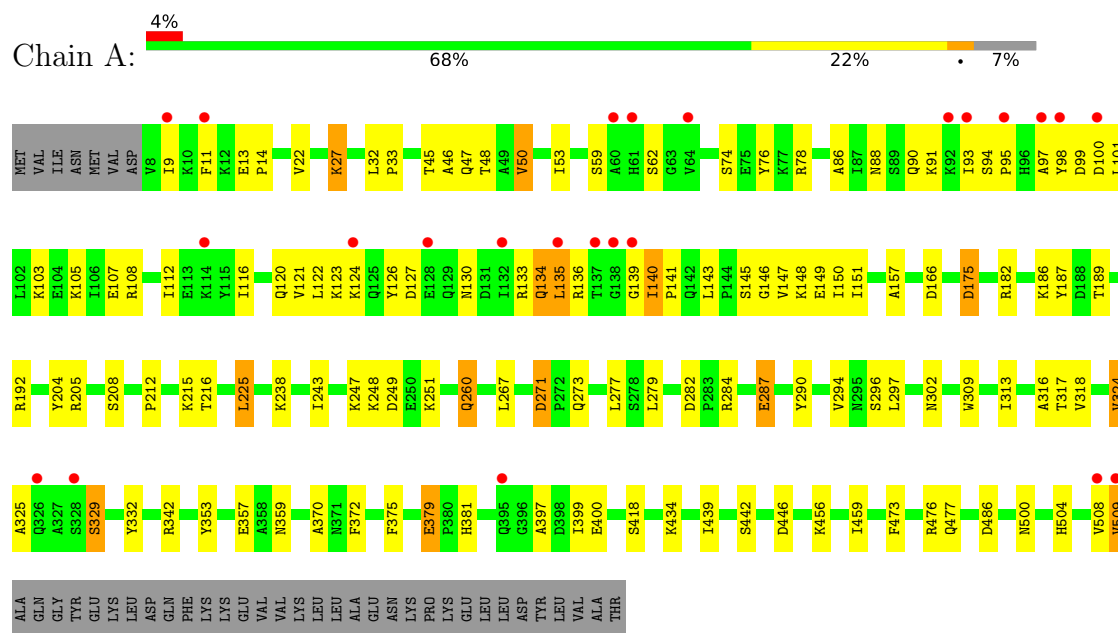
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	25	Total 25	O 25	0	0
4	B	3	Total 3	O 3	0	0
4	C	1	Total 1	O 1	0	0

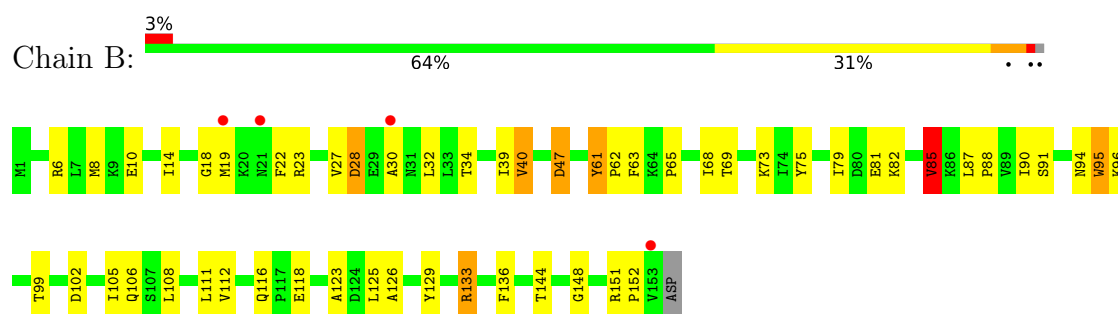
3 Residue-property plots

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.

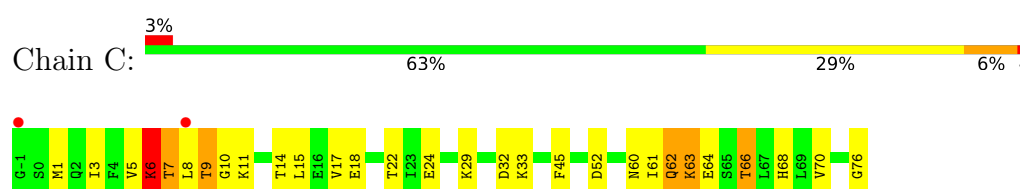
• Molecule 1: SidC



• Molecule 2: Ubiquitin-conjugating enzyme E2 L3



• Molecule 3: Polyubiquitin-B



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	101.53Å 101.53Å 352.33Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.90 50.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.7 (50.00-2.90) 99.7 (50.00-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.91 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.210 , 0.288 0.217 , 0.283	Depositor DCC
R_{free} test set	1261 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	62.5	Xtriage
Anisotropy	0.288	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 51.6	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	0.93	EDS
Total number of atoms	5957	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.58% 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	0.81	0/4148	1.06	22/5601 (0.4%)
2	B	0.80	0/1283	1.14	10/1732 (0.6%)
3	C	0.73	0/617	1.11	5/829 (0.6%)
All	All	0.80	0/6048	1.08	37/8162 (0.5%)

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	329	SER	N-CA-C	9.96	117.19	108.22
1	A	124	LYS	N-CA-C	8.77	124.47	113.43
3	C	63	LYS	N-CA-C	8.00	121.02	111.02
1	A	370	ALA	N-CA-C	7.73	120.55	110.43
3	C	9	THR	N-CA-C	7.57	122.59	112.30
2	B	81	GLU	N-CA-C	-7.51	103.08	111.71
2	B	61	TYR	C-N-CD	7.43	136.94	120.60
1	A	317	THR	N-CA-C	7.32	122.27	113.12
1	A	379	GLU	C-N-CD	7.27	136.60	120.60
1	A	50	VAL	CB-CA-C	-6.99	103.02	111.97
3	C	7	THR	N-CA-C	-6.88	101.45	110.53
1	A	271	ASP	CA-C-N	6.69	128.20	119.84
1	A	271	ASP	C-N-CA	6.69	128.20	119.84
1	A	399	ILE	N-CA-C	6.40	117.97	111.00
1	A	297	LEU	N-CA-C	6.34	118.27	111.36
3	C	11	LYS	N-CA-C	6.15	119.19	109.96
2	B	30	ALA	N-CA-C	6.03	117.86	111.28
1	A	135	LEU	N-CA-C	6.01	118.92	109.96
3	C	6	LYS	N-CA-C	5.63	118.14	109.07
1	A	318	VAL	N-CA-C	5.56	116.58	108.58
2	B	40	VAL	CA-C-N	-5.54	114.55	120.03
2	B	40	VAL	C-N-CA	-5.54	114.55	120.03
1	A	90	GLN	N-CA-C	5.48	118.06	108.52
1	A	140	ILE	CA-C-N	-5.36	114.40	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	140	ILE	C-N-CA	-5.36	114.40	119.76
2	B	95	TRP	N-CA-C	5.34	118.10	110.24
1	A	50	VAL	N-CA-C	5.32	115.53	110.42
1	A	189	THR	CA-C-N	-5.29	113.17	119.05
1	A	189	THR	C-N-CA	-5.29	113.17	119.05
1	A	175	ASP	CA-C-N	-5.21	114.37	119.99
1	A	175	ASP	C-N-CA	-5.21	114.37	119.99
2	B	85	VAL	CB-CA-C	-5.18	102.96	110.63
1	A	59	SER	CB-CA-C	5.13	119.10	109.86
1	A	139	GLY	N-CA-C	-5.07	101.16	113.18
2	B	96	LYS	CA-C-N	-5.02	114.44	119.56
2	B	96	LYS	C-N-CA	-5.02	114.44	119.56
2	B	28	ASP	N-CA-C	-5.02	102.22	109.59

There are no chirality outliers.

There are no planarity outliers.

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	4065	0	4005	128	0
2	B	1252	0	1271	53	0
3	C	611	0	637	14	0
4	A	25	0	0	9	0
4	B	3	0	0	2	0
4	C	1	0	0	0	0
All	All	5957	0	5913	188	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 16.

All (188) 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:48:THR:HB	1:A:140:ILE:HD11	1.23	1.19

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:LEU:O	1:A:127:ASP:HB2	1.48	1.14
1:A:53:ILE:HG12	1:A:135:LEU:CD2	1.91	1.00
2:B:65:PRO:HB2	2:B:90:ILE:HG23	1.50	0.94
1:A:53:ILE:HG23	1:A:135:LEU:HD13	1.51	0.93
1:A:74:SER:HB3	1:A:78:ARG:HH12	1.33	0.90
1:A:48:THR:CB	1:A:140:ILE:HD11	2.01	0.89
1:A:101:LEU:CD2	1:A:105:LYS:HG3	2.07	0.84
1:A:53:ILE:HG12	1:A:135:LEU:HD22	1.59	0.83
1:A:74:SER:CB	1:A:78:ARG:HH12	1.92	0.82
3:C:8:LEU:O	3:C:8:LEU:HD12	1.80	0.81
1:A:74:SER:C	1:A:78:ARG:NH1	2.40	0.79
1:A:101:LEU:HD23	1:A:105:LYS:HG3	1.61	0.79
1:A:53:ILE:HG12	1:A:135:LEU:CD1	2.11	0.79
1:A:53:ILE:HG23	1:A:135:LEU:CD1	2.12	0.79
1:A:101:LEU:HD23	1:A:101:LEU:O	1.82	0.79
1:A:101:LEU:HD21	1:A:105:LYS:HE3	1.67	0.76
1:A:135:LEU:HD23	1:A:141:PRO:HG3	1.66	0.76
1:A:53:ILE:CG1	1:A:135:LEU:CD2	2.63	0.76
1:A:135:LEU:CD2	1:A:141:PRO:HG3	2.15	0.76
2:B:22:PHE:CG	2:B:105:ILE:HD13	2.21	0.76
2:B:22:PHE:CD1	2:B:105:ILE:HD13	2.21	0.74
1:A:91:LYS:HA	1:A:94:SER:HB3	1.69	0.73
1:A:146:GLY:O	1:A:150:ILE:HG13	1.89	0.72
2:B:22:PHE:CD2	2:B:105:ILE:CD1	2.73	0.72
2:B:22:PHE:CG	2:B:105:ILE:CD1	2.73	0.72
1:A:212:PRO:HD3	4:A:606:HOH:O	1.90	0.71
1:A:225:LEU:C	1:A:225:LEU:HD23	2.15	0.71
1:A:13:GLU:OE2	1:A:108:ARG:NH1	2.25	0.70
1:A:53:ILE:CG1	1:A:135:LEU:HD22	2.22	0.70
1:A:508:VAL:O	1:A:509:VAL:HG13	1.91	0.69
1:A:309:TRP:O	1:A:313:ILE:HG13	1.93	0.68
1:A:192:ARG:CZ	2:B:73:LYS:HE2	2.23	0.68
1:A:103:LYS:O	1:A:107:GLU:HG3	1.95	0.67
1:A:45:THR:HG22	4:A:609:HOH:O	1.94	0.67
1:A:273:GLN:HE22	3:C:68:HIS:HB2	1.60	0.66
1:A:47:GLN:O	1:A:50:VAL:HG23	1.96	0.66
1:A:287:GLU:HG2	4:A:622:HOH:O	1.96	0.65
1:A:134:GLN:HE21	1:A:134:GLN:HA	1.60	0.65
1:A:88:ASN:HB3	1:A:91:LYS:HZ1	1.62	0.65
1:A:88:ASN:O	1:A:91:LYS:HG3	1.96	0.64
1:A:74:SER:C	1:A:78:ARG:HH11	2.05	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:47:ASP:O	2:B:151:ARG:HD2	1.97	0.64
1:A:504:HIS:HE1	3:C:32:ASP:O	1.81	0.64
3:C:29:LYS:O	3:C:33:LYS:HG2	1.97	0.64
1:A:134:GLN:HE21	1:A:134:GLN:CA	2.10	0.64
2:B:27:VAL:HG22	2:B:28:ASP:N	2.12	0.64
3:C:6:LYS:HD2	3:C:66:THR:HG22	1.80	0.64
1:A:47:GLN:O	1:A:48:THR:C	2.41	0.64
1:A:192:ARG:NH1	2:B:73:LYS:HE2	2.13	0.63
1:A:53:ILE:CG2	1:A:135:LEU:HD13	2.27	0.63
2:B:129:TYR:O	2:B:133:ARG:HG2	1.98	0.63
1:A:500:ASN:ND2	4:A:601:HOH:O	2.24	0.63
2:B:22:PHE:CD2	2:B:105:ILE:HD12	2.33	0.62
1:A:53:ILE:HG12	1:A:135:LEU:HD21	1.79	0.62
1:A:53:ILE:CG1	1:A:135:LEU:HD21	2.30	0.62
2:B:75:TYR:CD1	2:B:125:LEU:HD21	2.34	0.61
1:A:53:ILE:HG12	1:A:135:LEU:HD13	1.83	0.61
1:A:101:LEU:HD23	1:A:101:LEU:C	2.26	0.61
1:A:88:ASN:C	1:A:91:LYS:HZ2	2.09	0.60
3:C:7:THR:OG1	3:C:10:GLY:N	2.35	0.60
3:C:8:LEU:HG	3:C:9:THR:HG23	1.83	0.60
1:A:329:SER:HB2	1:A:400:GLU:OE2	2.01	0.60
2:B:22:PHE:HB2	2:B:105:ILE:HG21	1.82	0.60
2:B:65:PRO:CB	2:B:90:ILE:HG23	2.29	0.59
1:A:74:SER:O	1:A:78:ARG:NH1	2.34	0.59
1:A:99:ASP:O	1:A:100:ASP:HB2	2.02	0.59
1:A:212:PRO:CD	4:A:606:HOH:O	2.48	0.59
1:A:98:TYR:O	1:A:99:ASP:HB2	2.03	0.58
1:A:147:VAL:O	1:A:151:ILE:HG13	2.03	0.58
1:A:260:GLN:HG3	1:A:279:LEU:HG	1.86	0.58
1:A:282:ASP:OD2	1:A:284:ARG:NH2	2.36	0.58
1:A:74:SER:HB3	1:A:78:ARG:NH1	2.13	0.58
1:A:182:ARG:NH2	1:A:446:ASP:OD1	2.37	0.57
1:A:247:LYS:O	1:A:251:LYS:HG3	2.03	0.57
1:A:508:VAL:C	1:A:509:VAL:HG22	2.28	0.57
1:A:145:SER:O	1:A:149:GLU:HG3	2.04	0.57
1:A:32:LEU:HD12	1:A:33:PRO:HD2	1.87	0.57
2:B:61:TYR:CD1	2:B:62:PRO:HA	2.40	0.57
1:A:375:PHE:CZ	1:A:381:HIS:CE1	2.93	0.56
2:B:91:SER:HB3	2:B:94:ASN:HD22	1.71	0.56
1:A:442:SER:HB3	4:A:612:HOH:O	2.05	0.54
1:A:225:LEU:HD11	1:A:267:LEU:HD23	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:90:ILE:HG22	2:B:90:ILE:O	2.08	0.54
1:A:143:LEU:O	1:A:148:LYS:HE3	2.08	0.53
2:B:10:GLU:OE2	2:B:61:TYR:OH	2.23	0.53
2:B:39:ILE:O	2:B:39:ILE:HG13	2.08	0.53
1:A:93:ILE:O	1:A:93:ILE:HG12	2.08	0.53
2:B:91:SER:HB3	2:B:94:ASN:ND2	2.23	0.53
1:A:277:LEU:C	1:A:277:LEU:HD23	2.34	0.53
1:A:504:HIS:H	1:A:504:HIS:CD2	2.27	0.53
1:A:9:ILE:HD11	1:A:86:ALA:HB1	1.90	0.52
1:A:95:PRO:HG2	1:A:97:ALA:O	2.08	0.52
2:B:23:ARG:NH1	2:B:152:PRO:HB2	2.23	0.52
2:B:82:LYS:HD2	4:B:203:HOH:O	2.09	0.52
1:A:98:TYR:O	1:A:98:TYR:HD1	1.92	0.52
1:A:46:ALA:HB3	3:C:76:GLY:HA3	1.90	0.52
1:A:74:SER:C	1:A:78:ARG:HH12	2.18	0.51
1:A:287:GLU:CG	4:A:622:HOH:O	2.57	0.51
2:B:47:ASP:N	2:B:47:ASP:OD1	2.41	0.51
1:A:116:ILE:O	1:A:120:GLN:HG2	2.11	0.51
2:B:27:VAL:CG2	2:B:28:ASP:N	2.73	0.51
3:C:1:MET:SD	3:C:62:GLN:HA	2.51	0.51
1:A:135:LEU:C	1:A:136:ARG:HG3	2.36	0.50
2:B:144:THR:O	2:B:148:GLY:HA3	2.11	0.50
1:A:473:PHE:CZ	1:A:477:GLN:HG3	2.46	0.50
2:B:65:PRO:HB2	2:B:90:ILE:CG2	2.33	0.50
1:A:130:ASN:OD1	1:A:133:ARG:CD	2.59	0.50
1:A:332:TYR:OH	1:A:397:ALA:O	2.30	0.50
1:A:135:LEU:HD23	1:A:141:PRO:CG	2.37	0.50
2:B:102:ASP:O	2:B:106:GLN:HG2	2.12	0.50
3:C:63:LYS:O	3:C:64:GLU:HB2	2.12	0.49
3:C:1:MET:N	3:C:17:VAL:O	2.42	0.49
1:A:225:LEU:C	1:A:225:LEU:CD2	2.85	0.49
1:A:88:ASN:O	1:A:91:LYS:NZ	2.30	0.49
1:A:101:LEU:CD2	1:A:101:LEU:C	2.85	0.49
1:A:130:ASN:OD1	1:A:133:ARG:HD3	2.13	0.49
2:B:75:TYR:CE1	2:B:125:LEU:HD21	2.47	0.49
2:B:40:VAL:HG13	2:B:151:ARG:NH1	2.28	0.48
1:A:88:ASN:CB	1:A:91:LYS:HZ1	2.26	0.48
1:A:48:THR:HB	1:A:140:ILE:CD1	2.16	0.48
1:A:53:ILE:CD1	1:A:135:LEU:HD22	2.43	0.48
1:A:204:TYR:O	1:A:208:SER:OG	2.27	0.48
2:B:14:ILE:HG23	2:B:18:GLY:HA2	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:GLN:O	1:A:134:GLN:NE2	2.46	0.47
2:B:79:ILE:HG12	2:B:85:VAL:HG13	1.96	0.47
1:A:101:LEU:HD21	1:A:105:LYS:CE	2.43	0.47
2:B:87:LEU:HD12	2:B:111:LEU:HD22	1.97	0.47
1:A:456:LYS:HA	4:A:601:HOH:O	2.14	0.47
1:A:260:GLN:CG	1:A:279:LEU:HG	2.44	0.47
2:B:32:LEU:HD12	2:B:32:LEU:N	2.30	0.47
2:B:90:ILE:O	2:B:91:SER:C	2.56	0.46
1:A:359:ASN:ND2	1:A:372:PHE:H	2.14	0.46
1:A:53:ILE:HG12	1:A:135:LEU:CG	2.44	0.46
1:A:9:ILE:HG23	1:A:11:PHE:CZ	2.51	0.46
1:A:290:TYR:O	1:A:294:VAL:HG23	2.15	0.46
2:B:82:LYS:CD	4:B:203:HOH:O	2.64	0.45
1:A:14:PRO:HD2	1:A:76:TYR:CE2	2.51	0.45
1:A:324:VAL:O	1:A:325:ALA:C	2.57	0.45
1:A:98:TYR:O	1:A:98:TYR:CD1	2.70	0.45
1:A:205:ARG:NE	1:A:379:GLU:OE2	2.39	0.45
1:A:243:ILE:O	1:A:302:ASN:HA	2.17	0.45
2:B:22:PHE:CE2	2:B:105:ILE:HD11	2.52	0.44
1:A:434:LYS:HG2	4:A:616:HOH:O	2.18	0.44
3:C:8:LEU:HD12	3:C:8:LEU:C	2.35	0.44
1:A:282:ASP:OD1	1:A:282:ASP:C	2.59	0.44
2:B:22:PHE:CE2	2:B:105:ILE:CD1	3.00	0.44
3:C:45:PHE:CE2	3:C:61:ILE:HG12	2.53	0.44
1:A:101:LEU:HD21	1:A:105:LYS:HG3	1.98	0.44
3:C:22:THR:HG22	3:C:52:ASP:O	2.18	0.44
1:A:123:LYS:HD3	1:A:123:LYS:HA	1.72	0.43
1:A:353:TYR:OH	1:A:357:GLU:OE2	2.35	0.43
1:A:94:SER:HA	1:A:95:PRO:HD3	1.84	0.43
2:B:88:PRO:O	2:B:94:ASN:ND2	2.51	0.43
1:A:175:ASP:OD1	1:A:175:ASP:N	2.51	0.43
2:B:6:ARG:NH1	2:B:62:PRO:HG3	2.33	0.43
2:B:28:ASP:O	2:B:32:LEU:HD12	2.19	0.43
2:B:123:ALA:HA	2:B:126:ALA:HB3	1.99	0.43
2:B:62:PRO:O	2:B:95:TRP:NE1	2.46	0.43
1:A:88:ASN:CA	1:A:91:LYS:NZ	2.82	0.42
2:B:129:TYR:CE1	2:B:133:ARG:NH1	2.87	0.42
1:A:225:LEU:HD23	1:A:225:LEU:O	2.19	0.42
2:B:32:LEU:N	2:B:32:LEU:CD1	2.82	0.42
1:A:76:TYR:HD1	1:A:112:ILE:HD13	1.85	0.42
1:A:135:LEU:HD23	1:A:141:PRO:CB	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:79:ILE:HG12	2:B:85:VAL:CG1	2.49	0.42
2:B:136:PHE:C	2:B:136:PHE:CD1	2.97	0.42
1:A:271:ASP:CG	1:A:273:GLN:HG2	2.45	0.41
1:A:316:ALA:HA	2:B:63:PHE:CE1	2.55	0.41
1:A:22:VAL:HA	1:A:27:LYS:O	2.20	0.41
1:A:439:ILE:H	1:A:439:ILE:HG13	1.77	0.41
2:B:23:ARG:CZ	2:B:152:PRO:HG2	2.49	0.41
1:A:121:VAL:CG1	1:A:126:TYR:CD2	3.03	0.41
1:A:248:LYS:HG3	1:A:249:ASP:N	2.36	0.41
2:B:61:TYR:HA	2:B:62:PRO:HA	1.78	0.41
2:B:108:LEU:O	2:B:112:VAL:HG23	2.20	0.41
1:A:76:TYR:CD1	1:A:112:ILE:HD13	2.56	0.41
1:A:130:ASN:OD1	1:A:133:ARG:HB3	2.21	0.41
2:B:22:PHE:CD1	2:B:105:ILE:CD1	2.97	0.41
1:A:192:ARG:NH1	2:B:73:LYS:HG2	2.36	0.41
1:A:238:LYS:HD2	1:A:238:LYS:HA	1.93	0.41
1:A:135:LEU:O	1:A:136:ARG:CB	2.69	0.40
1:A:186:LYS:HE2	1:A:187:TYR:CZ	2.56	0.40
2:B:6:ARG:NH2	2:B:99:THR:O	2.54	0.40
1:A:157:ALA:HB3	1:A:459:ILE:HG13	2.02	0.40
1:A:166:ASP:C	1:A:166:ASP:OD1	2.64	0.40
1:A:53:ILE:HG13	1:A:135:LEU:HD21	2.03	0.40
1:A:150:ILE:HG12	1:A:476:ARG:HD2	2.04	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	500/542 (92%)	480 (96%)	20 (4%)	0	100	100
2	B	151/154 (98%)	143 (95%)	8 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	76/78 (97%)	70 (92%)	5 (7%)	1 (1%)	9	32
All	All	727/774 (94%)	693 (95%)	33 (4%)	1 (0%)	48	77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	24	GLU

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	454/490 (93%)	440 (97%)	14 (3%)	35	69
2	B	137/138 (99%)	127 (93%)	10 (7%)	13	38
3	C	69/69 (100%)	59 (86%)	10 (14%)	3	10
All	All	660/697 (95%)	626 (95%)	34 (5%)	21	52

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LYS
1	A	62	SER
1	A	134	GLN
1	A	215	LYS
1	A	216	THR
1	A	225	LEU
1	A	260	GLN
1	A	287	GLU
1	A	296	SER
1	A	324	VAL
1	A	342	ARG
1	A	418	SER
1	A	486	ASP
1	A	509	VAL

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Mol	Chain	Res	Type
2	B	8	MET
2	B	19	MET
2	B	34	THR
2	B	47	ASP
2	B	68	ILE
2	B	69	THR
2	B	85	VAL
2	B	116	GLN
2	B	118	GLU
2	B	133	ARG
3	C	3	ILE
3	C	5	VAL
3	C	6	LYS
3	C	14	THR
3	C	15	LEU
3	C	18	GLU
3	C	60	ASN
3	C	62	GLN
3	C	66	THR
3	C	70	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	72	GLN
1	A	129	GLN
1	A	134	GLN
1	A	273	GLN
1	A	302	ASN
1	A	365	ASN
1	A	381	HIS
1	A	462	HIS
1	A	469	HIS
1	A	477	GLN
1	A	500	ASN
1	A	503	HIS
1	A	504	HIS
2	B	24	ASN
2	B	76	HIS
2	B	78	ASN
2	B	84	GLN

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Mol	Chain	Res	Type
2	B	94	ASN
2	B	116	GLN
3	C	25	ASN
3	C	40	GLN
3	C	60	ASN
3	C	62	GLN
3	C	68	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	502/542 (92%)	0.10	24 (4%) 35 28	32, 61, 125, 159	3 (0%)
2	B	153/154 (99%)	0.24	4 (2%) 57 48	40, 65, 112, 126	0
3	C	78/78 (100%)	0.32	2 (2%) 57 48	47, 86, 121, 128	0
All	All	733/774 (94%)	0.15	30 (4%) 41 33	32, 64, 122, 159	3 (0%)

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	114	LYS	6.9
1	A	395	GLN	6.1
1	A	98	TYR	4.1
1	A	93	ILE	4.0
1	A	509	VAL	4.0
1	A	135	LEU	3.8
1	A	97	ALA	3.3
2	B	30	ALA	3.2
1	A	9	ILE	3.2
1	A	128	GLU	3.2
2	B	19	MET	3.2
1	A	326	GLN	3.1
1	A	139	GLY	2.9
1	A	508	VAL	2.9
1	A	328	SER	2.8
2	B	153	VAL	2.7
3	C	8	LEU	2.7
1	A	137	THR	2.5
1	A	92	LYS	2.4
1	A	100	ASP	2.3
1	A	64	VAL	2.3
2	B	21	ASN	2.3
1	A	132	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	11	PHE	2.2
1	A	61	HIS	2.2
1	A	95	PRO	2.0
1	A	138	GLY	2.0
3	C	-1	GLY	2.0
1	A	60	ALA	2.0
1	A	124	LYS	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.