



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 10:11 AM UTC

PDB ID : 2OXQ / pdb_00002oxq
Title : Structure of the Ubch5 :CHIP U-box complex
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Deposited on : 2007-02-20
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

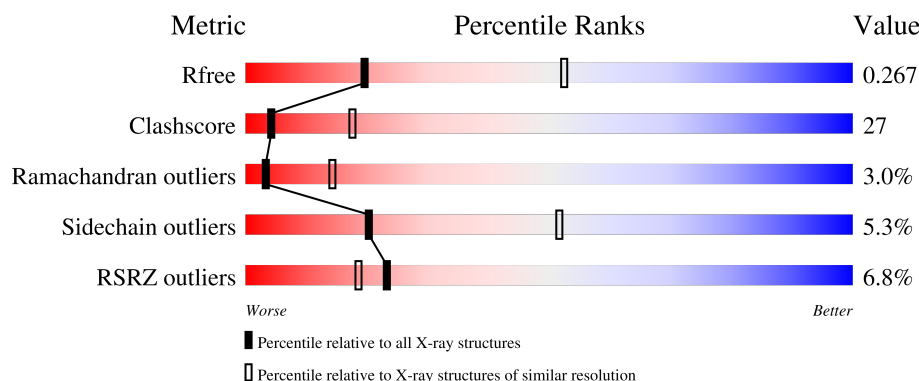
1 Overall quality at a glance

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	152	9% 50% 39% 6% ••
1	B	152	6% 59% 32% 7% •
2	C	80	4% 46% 35% 9% 10%
2	D	80	5% 40% 46% • 10%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	CL	D	7	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3525 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 Ubiquitin-conjugating enzyme E2D 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	149	Total	C	N	O	S	0	0	0
			1182	759	198	218	7			
1	B	149	Total	C	N	O	S	0	0	0
			1182	759	198	218	7			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP Q6PC58
A	-3	ALA	-	expression tag	UNP Q6PC58
A	-2	MET	-	expression tag	UNP Q6PC58
A	-1	GLY	-	expression tag	UNP Q6PC58
A	0	SER	-	expression tag	UNP Q6PC58
B	-4	GLY	-	expression tag	UNP Q6PC58
B	-3	ALA	-	expression tag	UNP Q6PC58
B	-2	MET	-	expression tag	UNP Q6PC58
B	-1	GLY	-	expression tag	UNP Q6PC58
B	0	SER	-	expression tag	UNP Q6PC58

- Molecule 2 is a protein called STIP1 homology and U-Box containing protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	72	Total	C	N	O	S	0	0	0
			577	365	94	114	4			
2	D	72	Total	C	N	O	S	0	0	0
			577	365	94	114	4			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	199	GLY	-	expression tag	UNP Q7ZTZ6
C	200	ALA	-	expression tag	UNP Q7ZTZ6

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Chain	Residue	Modelled	Actual	Comment	Reference
C	201	MET	-	expression tag	UNP Q7ZTZ6
C	202	GLY	-	expression tag	UNP Q7ZTZ6
C	203	SER	-	expression tag	UNP Q7ZTZ6
C	204	LYS	-	expression tag	UNP Q7ZTZ6
C	205	LYS	-	expression tag	UNP Q7ZTZ6
C	206	ARG	-	expression tag	UNP Q7ZTZ6
D	199	GLY	-	expression tag	UNP Q7ZTZ6
D	200	ALA	-	expression tag	UNP Q7ZTZ6
D	201	MET	-	expression tag	UNP Q7ZTZ6
D	202	GLY	-	expression tag	UNP Q7ZTZ6
D	203	SER	-	expression tag	UNP Q7ZTZ6
D	204	LYS	-	expression tag	UNP Q7ZTZ6
D	205	LYS	-	expression tag	UNP Q7ZTZ6
D	206	ARG	-	expression tag	UNP Q7ZTZ6

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	C	1	Total Cl 1 1	0	0
3	D	2	Total Cl 2 2	0	0

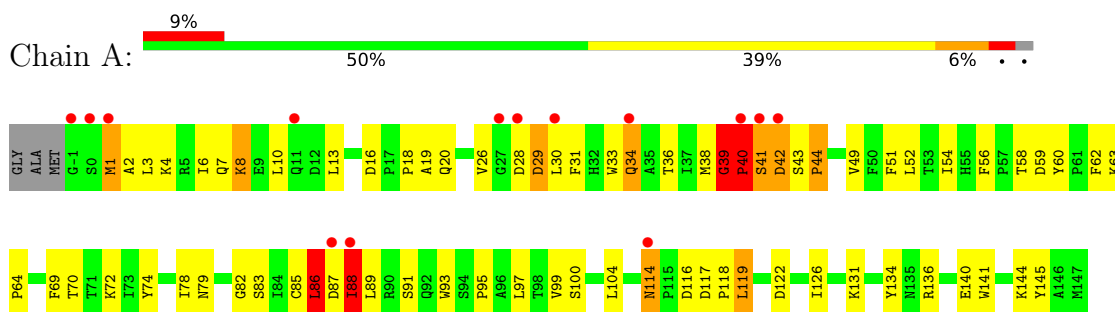
- Molecule 4 is water.

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

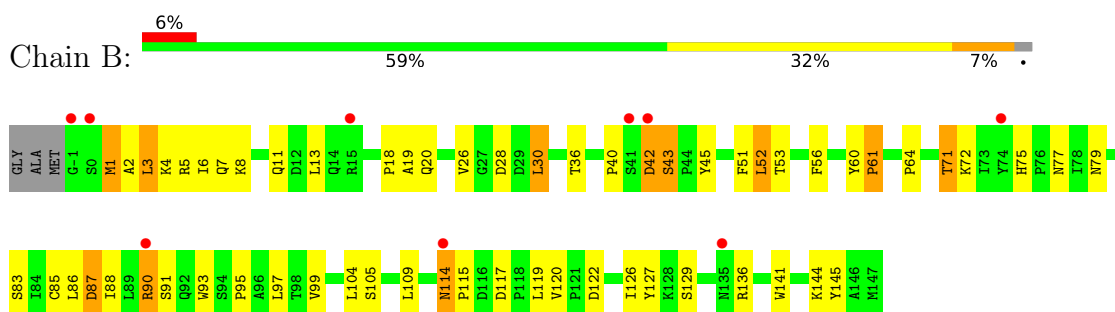
3 Residue-property plots

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.

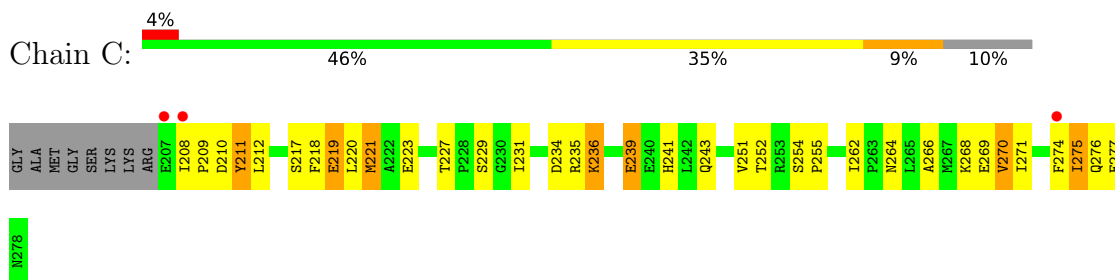
• Molecule 1: Ubiquitin-conjugating enzyme E2D 1



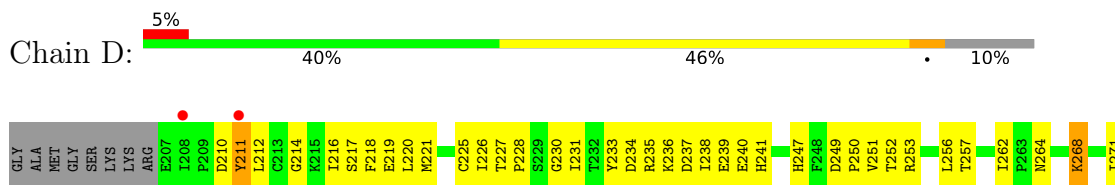
• Molecule 1: Ubiquitin-conjugating enzyme E2D 1



• Molecule 2: STIP1 homology and U-Box containing protein 1



• Molecule 2: STIP1 homology and U-Box containing protein 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.04Å 93.40Å 144.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	53.23 – 2.90 53.23 – 2.90	Depositor EDS
% Data completeness (in resolution range)	93.2 (53.23-2.90) 93.2 (53.23-2.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.93 (at 2.91Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.240 , 0.272 0.237 , 0.267	Depositor DCC
R_{free} test set	1165 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	65.4	Xtriage
Anisotropy	0.534	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3525	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.42% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section:
CL

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	0.49	0/1218	1.05	9/1659 (0.5%)
1	B	0.46	0/1218	1.02	5/1659 (0.3%)
2	C	0.44	0/589	0.94	2/798 (0.3%)
2	D	0.43	0/589	0.93	1/798 (0.1%)
All	All	0.46	0/3614	1.01	17/4914 (0.3%)

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	3
1	B	0	1
All	All	0	4

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	28	ASP	N-CA-C	-10.25	100.08	112.59
1	A	70	THR	N-CA-C	-8.66	103.22	113.88
1	A	119	LEU	N-CA-C	-7.56	103.42	114.39
1	A	114	ASN	CA-C-N	7.56	127.23	119.82
1	A	114	ASN	C-N-CA	7.56	127.23	119.82
1	B	114	ASN	CA-C-N	6.73	126.36	119.56
1	B	114	ASN	C-N-CA	6.73	126.36	119.56
1	A	1	MET	N-CA-C	-6.41	104.39	111.82
2	C	270	VAL	N-CA-C	-6.01	104.65	110.72
1	B	120	VAL	CA-C-N	5.89	127.21	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	120	VAL	C-N-CA	5.89	127.21	119.84
1	A	16	ASP	CA-C-N	5.89	126.44	120.38
1	A	16	ASP	C-N-CA	5.89	126.44	120.38
1	A	34	GLN	N-CA-C	-5.70	100.60	109.72
1	A	44	PRO	N-CA-C	-5.35	109.11	114.68
2	C	221	MET	N-CA-C	5.25	117.28	109.25
2	D	234	ASP	N-CA-C	-5.09	100.43	108.73

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	39	GLY	Peptide
1	A	40	PRO	Peptide
1	A	86	LEU	Peptide
1	B	40	PRO	Peptide

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	1182	0	1168	60	0
1	B	1182	0	1168	63	0
2	C	577	0	564	35	0
2	D	577	0	564	45	0
3	C	1	0	0	0	0
3	D	2	0	0	2	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
All	All	3525	0	3464	190	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 27.

All (190) 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:40:PRO:HB2	1:A:43:SER:OG	1.61	0.99
1:A:7:GLN:HG2	1:A:30:LEU:HD12	1.54	0.88
2:D:236:LYS:O	2:D:240:GLU:HG3	1.79	0.82
1:B:86:LEU:HD13	1:B:109:LEU:HD22	1.62	0.82
1:B:4:LYS:NZ	2:D:219:GLU:HG2	1.97	0.79
1:A:87:ASP:O	1:A:88:ILE:HB	1.80	0.79
1:B:77:ASN:HD21	1:B:114:ASN:H	1.30	0.77
1:B:51:PHE:H	1:B:71:THR:HG23	1.50	0.76
1:A:3:LEU:O	1:A:7:GLN:HG3	1.84	0.75
1:B:72:LYS:HE2	1:B:141:TRP:CZ3	2.22	0.75
1:B:75:HIS:HE1	1:B:77:ASN:ND2	1.84	0.75
1:B:75:HIS:HE1	1:B:77:ASN:HD22	1.36	0.74
1:B:51:PHE:H	1:B:71:THR:CG2	2.00	0.74
2:D:276:GLN:HE21	2:D:276:GLN:HA	1.52	0.73
1:B:79:ASN:HB3	1:B:119:LEU:HD22	1.71	0.72
2:C:268:LYS:O	2:C:268:LYS:HD3	1.90	0.72
1:A:4:LYS:O	1:A:8:LYS:HD2	1.91	0.71
1:A:26:VAL:HG22	1:A:34:GLN:HB2	1.71	0.71
1:A:39:GLY:O	1:A:41:SER:N	2.22	0.71
1:A:41:SER:O	1:A:42:ASP:CB	2.36	0.71
2:C:262:ILE:HD13	2:D:228:PRO:HA	1.71	0.70
1:B:79:ASN:ND2	1:B:83:SER:HB2	2.06	0.70
1:B:4:LYS:HZ2	2:D:219:GLU:HG2	1.57	0.70
1:A:72:LYS:HE3	1:A:141:TRP:CZ3	2.28	0.69
1:A:40:PRO:O	1:A:43:SER:N	2.21	0.69
1:B:86:LEU:HG	1:B:88:ILE:HG22	1.75	0.68
2:C:231:ILE:HD13	2:C:251:VAL:HB	1.75	0.68
1:B:72:LYS:HE2	1:B:141:TRP:CH2	2.30	0.67
1:B:85:CYS:HB2	1:B:119:LEU:HD12	1.77	0.67
2:D:247:HIS:HD2	2:D:257:THR:HG22	1.60	0.67
1:A:40:PRO:O	1:A:43:SER:OG	2.09	0.66
1:A:144:LYS:HD3	1:A:145:TYR:CZ	2.32	0.64
2:D:276:GLN:HA	2:D:276:GLN:NE2	2.12	0.64
2:D:262:ILE:HD12	2:D:262:ILE:N	2.13	0.63
2:D:277:GLU:O	2:D:278:ASN:ND2	2.32	0.63
1:A:87:ASP:O	1:A:88:ILE:CB	2.46	0.62
2:C:208:ILE:HG21	2:C:220:LEU:HD13	1.81	0.62
1:A:4:LYS:HG2	1:A:8:LYS:HE2	1.82	0.61
1:B:95:PRO:HG2	2:D:241:HIS:HD2	1.64	0.61
2:D:268:LYS:O	2:D:268:LYS:HD3	2.00	0.61
1:B:79:ASN:HD22	1:B:119:LEU:HD21	1.66	0.61
2:C:274:PHE:C	2:C:276:GLN:H	2.07	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:217:SER:O	2:D:218:PHE:HB2	1.99	0.61
2:D:247:HIS:CD2	2:D:257:THR:HG22	2.36	0.61
1:A:43:SER:HB2	1:A:44:PRO:HD2	1.84	0.60
1:A:26:VAL:CG2	1:A:34:GLN:HB2	2.31	0.60
1:B:51:PHE:C	1:B:52:LEU:HD23	2.27	0.60
1:B:79:ASN:HD21	1:B:83:SER:HB2	1.66	0.59
1:B:13:LEU:C	1:B:13:LEU:HD13	2.27	0.59
1:A:38:MET:O	1:A:40:PRO:HD3	2.03	0.58
2:D:225:CYS:HA	2:D:264:ASN:H	1.69	0.58
2:C:211:TYR:CD1	2:C:212:LEU:HG	2.39	0.57
2:D:268:LYS:HD3	2:D:268:LYS:C	2.29	0.57
1:A:1:MET:HE1	1:A:62:PHE:HE2	1.70	0.57
2:D:219:GLU:HG3	2:D:220:LEU:N	2.20	0.57
1:B:13:LEU:HD22	1:B:18:PRO:HD3	1.85	0.57
1:A:95:PRO:HG2	2:C:241:HIS:HD2	1.71	0.56
1:A:19:ALA:O	1:A:20:GLN:HB2	2.06	0.56
2:C:269:GLU:OE2	2:D:211:TYR:HB3	2.05	0.56
2:D:214:GLY:HA3	3:D:7:CL:CL	2.42	0.56
1:A:59:ASP:HB2	1:A:63:LYS:HB2	1.88	0.56
1:B:72:LYS:HD3	1:B:145:TYR:CE2	2.40	0.56
1:A:41:SER:O	1:A:42:ASP:HB2	2.06	0.55
2:D:276:GLN:HE21	2:D:276:GLN:CA	2.16	0.55
1:B:42:ASP:O	1:B:43:SER:C	2.49	0.55
1:A:29:ASP:C	1:A:31:PHE:H	2.14	0.55
1:B:7:GLN:HG2	1:B:30:LEU:HD22	1.88	0.54
2:D:264:ASN:C	2:D:264:ASN:HD22	2.16	0.54
1:A:7:GLN:HG2	1:A:30:LEU:CD1	2.34	0.54
1:A:41:SER:O	1:A:42:ASP:HB3	2.08	0.53
1:B:52:LEU:HD23	1:B:52:LEU:N	2.22	0.53
1:A:28:ASP:O	1:A:29:ASP:C	2.51	0.53
2:C:235:ARG:O	2:C:239:GLU:HB2	2.08	0.53
1:B:18:PRO:HB3	1:B:104:LEU:CD2	2.39	0.53
2:C:268:LYS:HD3	2:C:268:LYS:C	2.34	0.53
2:C:221:MET:O	2:C:271:ILE:HD13	2.08	0.53
2:C:275:ILE:O	2:C:275:ILE:HG22	2.09	0.53
2:D:253:ARG:HG3	2:D:253:ARG:HH11	1.73	0.52
1:B:86:LEU:HD21	1:B:105:SER:CB	2.38	0.52
1:A:4:LYS:CG	1:A:8:LYS:HE2	2.39	0.52
1:B:8:LYS:O	1:B:11:GLN:HB2	2.10	0.52
2:C:234:ASP:CG	2:C:236:LYS:HE2	2.35	0.52
2:C:254:SER:HB2	2:C:255:PRO:HD2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:217:SER:C	2:C:219:GLU:H	2.17	0.51
1:B:3:LEU:O	1:B:7:GLN:HG3	2.11	0.51
1:B:5:ARG:NH2	2:D:216:ILE:O	2.44	0.51
1:B:60:TYR:CD1	1:B:61:PRO:HA	2.46	0.51
1:B:19:ALA:O	1:B:20:GLN:HB2	2.11	0.50
1:B:64:PRO:HB3	1:B:93:TRP:CG	2.46	0.50
1:B:114:ASN:ND2	1:B:117:ASP:HB3	2.26	0.50
2:D:217:SER:HB3	2:D:237:ASP:OD2	2.12	0.49
1:B:2:ALA:O	1:B:6:ILE:HG13	2.11	0.49
2:D:227:THR:HB	2:D:228:PRO:HD2	1.94	0.49
1:B:13:LEU:CD2	1:B:18:PRO:HD3	2.43	0.49
2:D:210:ASP:C	2:D:212:LEU:H	2.20	0.49
1:A:64:PRO:HB3	1:A:93:TRP:CG	2.48	0.49
1:A:86:LEU:O	1:A:89:LEU:HB2	2.12	0.49
2:D:235:ARG:HG2	2:D:239:GLU:OE1	2.13	0.49
1:B:13:LEU:HD13	1:B:13:LEU:O	2.13	0.49
2:D:226:ILE:HD11	2:D:230:GLY:HA2	1.93	0.49
1:A:18:PRO:HB3	1:A:104:LEU:HD21	1.95	0.49
1:B:56:PHE:HZ	1:B:99:VAL:HG21	1.78	0.49
1:A:38:MET:SD	1:A:49:VAL:HG22	2.53	0.48
1:A:36:THR:HG22	1:A:51:PHE:HD1	1.77	0.48
1:B:42:ASP:O	1:B:43:SER:O	2.31	0.48
1:A:29:ASP:C	1:A:31:PHE:N	2.69	0.48
1:B:56:PHE:CZ	1:B:99:VAL:HG21	2.49	0.48
2:C:264:ASN:C	2:C:264:ASN:HD22	2.19	0.48
2:C:274:PHE:C	2:C:276:GLN:N	2.69	0.48
1:A:10:LEU:HB2	1:A:33:TRP:CZ2	2.49	0.47
1:B:95:PRO:HG2	2:D:241:HIS:CD2	2.47	0.47
1:A:78:ILE:HA	1:A:83:SER:O	2.14	0.47
2:D:241:HIS:CE1	2:D:250:PRO:HG3	2.50	0.47
1:B:90:ARG:HG3	1:B:90:ARG:HH11	1.79	0.47
1:B:122:ASP:O	1:B:126:ILE:HG12	2.14	0.47
2:D:221:MET:O	2:D:271:ILE:HD13	2.14	0.47
1:A:2:ALA:O	1:A:6:ILE:HG13	2.14	0.47
2:D:233:TYR:CB	2:D:238:ILE:HB	2.45	0.47
2:C:239:GLU:O	2:C:243:GLN:HG3	2.15	0.47
1:B:88:ILE:CG1	1:B:97:LEU:HD13	2.45	0.46
1:A:33:TRP:HB2	1:A:54:ILE:HB	1.98	0.46
2:C:217:SER:O	2:C:218:PHE:HB2	2.15	0.46
1:A:131:LYS:O	1:A:134:TYR:HB3	2.15	0.46
1:B:86:LEU:HD21	1:B:105:SER:HB2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:LYS:HD3	1:A:145:TYR:CE2	2.50	0.46
1:A:26:VAL:O	1:A:26:VAL:HG12	2.15	0.46
2:D:216:ILE:HB	3:D:7:CL:CL	2.53	0.46
1:B:4:LYS:HE3	2:D:219:GLU:OE1	2.15	0.46
1:B:1:MET:HE3	2:D:237:ASP:OD1	2.15	0.45
1:A:1:MET:HE1	1:A:62:PHE:CE2	2.49	0.45
1:A:116:ASP:C	1:A:118:PRO:HD3	2.41	0.45
1:A:136:ARG:O	1:A:140:GLU:HG3	2.17	0.45
1:B:4:LYS:HZ1	2:D:219:GLU:HG2	1.74	0.45
1:B:18:PRO:HB3	1:B:104:LEU:HD21	1.98	0.45
2:D:231:ILE:HD13	2:D:251:VAL:HB	1.99	0.45
2:C:231:ILE:CD1	2:C:251:VAL:HB	2.46	0.44
1:A:79:ASN:OD1	1:A:83:SER:HB2	2.17	0.44
1:A:85:CYS:HB2	1:A:119:LEU:HG	1.98	0.44
1:B:88:ILE:HD11	1:B:97:LEU:HD13	2.00	0.44
2:C:209:PRO:HD3	2:C:274:PHE:CZ	2.51	0.44
2:C:266:ALA:O	2:C:270:VAL:HG23	2.17	0.44
1:A:13:LEU:HD23	1:A:100:SER:HB2	1.98	0.44
1:A:88:ILE:HG13	1:A:97:LEU:HD13	2.00	0.44
2:D:277:GLU:O	2:D:278:ASN:CG	2.61	0.44
1:A:52:LEU:HD12	1:A:52:LEU:N	2.33	0.43
1:B:8:LYS:HG2	2:D:218:PHE:HB3	2.00	0.43
1:B:115:PRO:HB3	1:B:127:TYR:HD2	1.81	0.43
2:C:227:THR:C	2:C:229:SER:H	2.26	0.43
1:A:122:ASP:O	1:A:126:ILE:HG12	2.19	0.43
2:C:234:ASP:O	2:C:235:ARG:C	2.61	0.43
2:C:231:ILE:HD11	2:C:252:THR:HG23	1.99	0.43
1:A:36:THR:HG22	1:A:51:PHE:CD1	2.54	0.43
1:B:86:LEU:HD21	1:B:105:SER:HB3	2.00	0.43
2:D:249:ASP:O	2:D:253:ARG:N	2.48	0.43
1:B:90:ARG:CG	1:B:91:SER:N	2.80	0.43
2:C:211:TYR:CE1	2:C:212:LEU:HG	2.54	0.43
2:D:262:ILE:N	2:D:262:ILE:CD1	2.82	0.43
1:A:59:ASP:O	1:A:60:TYR:C	2.60	0.42
1:B:144:LYS:HE3	1:B:144:LYS:HB2	1.77	0.42
2:C:210:ASP:C	2:C:212:LEU:H	2.28	0.42
1:A:40:PRO:HB2	1:A:43:SER:CB	2.49	0.42
1:A:56:PHE:CZ	1:A:99:VAL:HG21	2.54	0.42
1:B:88:ILE:HD11	1:B:97:LEU:HD22	2.01	0.42
2:D:252:THR:O	2:D:253:ARG:HB2	2.20	0.42
2:C:212:LEU:O	2:C:220:LEU:HD12	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:86:LEU:O	1:B:87:ASP:C	2.62	0.42
2:C:210:ASP:C	2:C:212:LEU:N	2.78	0.42
1:A:72:LYS:HE3	1:A:141:TRP:CH2	2.55	0.41
1:B:43:SER:HB2	1:B:45:TYR:HD2	1.85	0.41
1:B:79:ASN:OD1	1:B:79:ASN:C	2.63	0.41
1:B:90:ARG:HG2	1:B:91:SER:N	2.35	0.41
2:D:247:HIS:HB3	2:D:256:LEU:O	2.20	0.41
1:B:13:LEU:C	1:B:13:LEU:CD1	2.93	0.41
1:B:79:ASN:HD22	1:B:119:LEU:CD2	2.32	0.41
2:C:277:GLU:CD	2:C:277:GLU:N	2.78	0.41
1:A:4:LYS:HE3	1:A:4:LYS:HB2	1.85	0.41
1:B:75:HIS:CE1	1:B:77:ASN:HD22	2.26	0.41
1:B:86:LEU:O	1:B:88:ILE:N	2.54	0.41
1:A:38:MET:O	1:A:40:PRO:CD	2.67	0.41
2:C:210:ASP:O	2:C:212:LEU:N	2.54	0.41
2:C:262:ILE:CD1	2:D:228:PRO:HA	2.45	0.41
1:A:34:GLN:HE22	1:A:51:PHE:HB3	1.85	0.41
1:A:74:TYR:HB2	1:A:141:TRP:CD2	2.56	0.41
2:C:231:ILE:CD1	2:C:252:THR:HG23	2.51	0.41
2:D:235:ARG:O	2:D:239:GLU:HG3	2.21	0.41
2:D:236:LYS:HG3	2:D:237:ASP:N	2.36	0.40
2:C:266:ALA:O	2:C:269:GLU:HG2	2.20	0.40
1:A:69:PHE:O	1:A:82:GLY:HA3	2.21	0.40
1:A:114:ASN:ND2	1:A:117:ASP:HB3	2.37	0.40
2:C:223:GLU:OE2	2:C:235:ARG:HD2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	147/152 (97%)	134 (91%)	7 (5%)	6 (4%)	2 9

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	147/152 (97%)	133 (90%)	10 (7%)	4 (3%)	4	16
2	C	70/80 (88%)	61 (87%)	7 (10%)	2 (3%)	3	15
2	D	70/80 (88%)	63 (90%)	6 (9%)	1 (1%)	9	30
All	All	434/464 (94%)	391 (90%)	30 (7%)	13 (3%)	3	14

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	40	PRO
1	A	88	ILE
1	B	87	ASP
1	A	42	ASP
1	B	43	SER
2	D	211	TYR
1	A	29	ASP
1	A	41	SER
1	B	129	SER
2	C	211	TYR
1	B	26	VAL
1	A	39	GLY
2	C	275	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	133/134 (99%)	128 (96%)	5 (4%)	29	64
1	B	133/134 (99%)	122 (92%)	11 (8%)	10	32
2	C	66/71 (93%)	63 (96%)	3 (4%)	24	58
2	D	66/71 (93%)	64 (97%)	2 (3%)	36	70
All	All	398/410 (97%)	377 (95%)	21 (5%)	20	52

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LYS
1	A	58	THR
1	A	86	LEU
1	A	88	ILE
1	A	91	SER
1	B	1	MET
1	B	3	LEU
1	B	30	LEU
1	B	36	THR
1	B	42	ASP
1	B	52	LEU
1	B	53	THR
1	B	61	PRO
1	B	71	THR
1	B	90	ARG
1	B	136	ARG
2	C	219	GLU
2	C	236	LYS
2	C	239	GLU
2	D	268	LYS
2	D	276	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	ASN
1	A	114	ASN
1	B	34	GLN
1	B	77	ASN
1	B	81	ASN
1	B	114	ASN
1	B	143	GLN
2	C	241	HIS
2	C	264	ASN
2	D	241	HIS
2	D	247	HIS
2	D	264	ASN
2	D	276	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	149/152 (98%)	0.32	14 (9%) 14 12	40, 55, 85, 99	0
1	B	149/152 (98%)	0.42	9 (6%) 27 21	43, 60, 94, 102	0
2	C	72/80 (90%)	0.29	3 (4%) 40 32	41, 63, 96, 113	0
2	D	72/80 (90%)	0.37	4 (5%) 30 23	51, 68, 95, 105	0
All	All	442/464 (95%)	0.36	30 (6%) 23 18	40, 60, 94, 113	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	-1	GLY	6.9
1	A	40	PRO	3.7
1	B	-1	GLY	3.6
1	A	88	ILE	3.5
1	A	87	ASP	3.2
1	A	0	SER	3.2
2	D	275	ILE	3.1
1	A	42	ASP	3.0
1	B	0	SER	3.0
1	A	28	ASP	3.0
1	A	114	ASN	2.9
1	B	41	SER	2.8
1	B	15	ARG	2.7
1	A	11	GLN	2.7
1	A	30	LEU	2.6
2	C	208	ILE	2.5
1	B	42	ASP	2.5
1	A	34	GLN	2.5
2	C	207	GLU	2.4
1	A	41	SER	2.4
1	A	27	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	114	ASN	2.3
2	D	211	TYR	2.3
1	B	90	ARG	2.2
1	B	74	TYR	2.2
2	D	208	ILE	2.1
2	D	278	ASN	2.1
1	A	1	MET	2.0
1	B	135	ASN	2.0
2	C	274	PHE	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](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	CL	D	6	1/1	0.88	0.14	97,97,97,97	0
3	CL	C	1	1/1	0.97	0.16	60,60,60,60	0
3	CL	D	7	1/1	0.98	0.07	65,65,65,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.