



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:39 AM UTC

PDB ID : 8S7L / pdb_00008s7l
Title : Crystal structure of a double mutant of VirB8-like OrfG central and C-terminal domains of Streptococcus thermophilus ICEst3 (Gram positive conjugative type IV secretion system).
Authors : Favier, F.; Didierjean, C.; Maffo-Woulefack, R.; Douzi, B.; Leblond-Bourget, N.
Deposited on : 2024-03-03
Resolution : 2.60 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

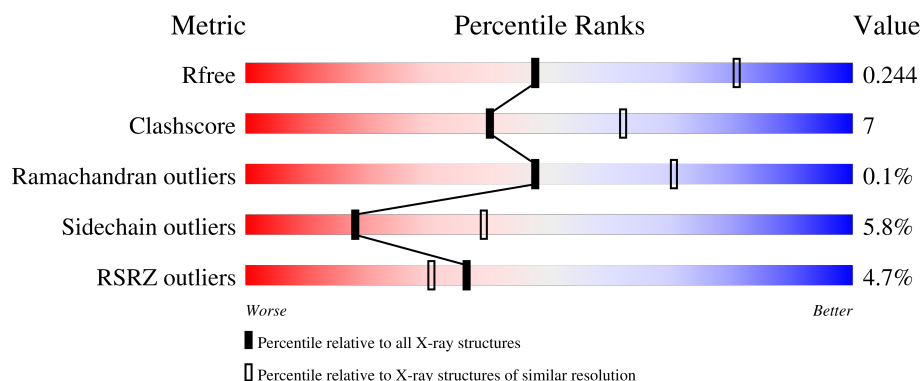
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	268	
1	B	268	
1	C	268	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 11398 atoms, of which 5486 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 Putative transfer protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	241	Total	C	H	N	O	S	0	0	0
			3823	1271	1840	316	393	3			
1	B	238	Total	C	H	N	O	S	0	0	0
			3777	1257	1819	312	386	3			
1	C	240	Total	C	H	N	O	S	0	0	0
			3798	1264	1827	314	390	3			

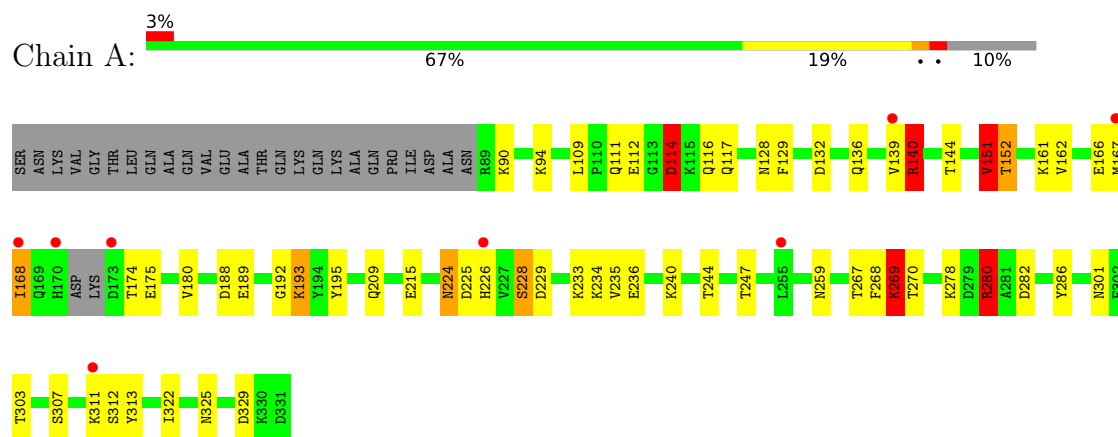
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	204	CYS	ALA	engineered mutation	UNP Q70CA4
A	264	CYS	ALA	engineered mutation	UNP Q70CA4
B	204	CYS	ALA	engineered mutation	UNP Q70CA4
B	264	CYS	ALA	engineered mutation	UNP Q70CA4
C	204	CYS	ALA	engineered mutation	UNP Q70CA4
C	264	CYS	ALA	engineered mutation	UNP Q70CA4

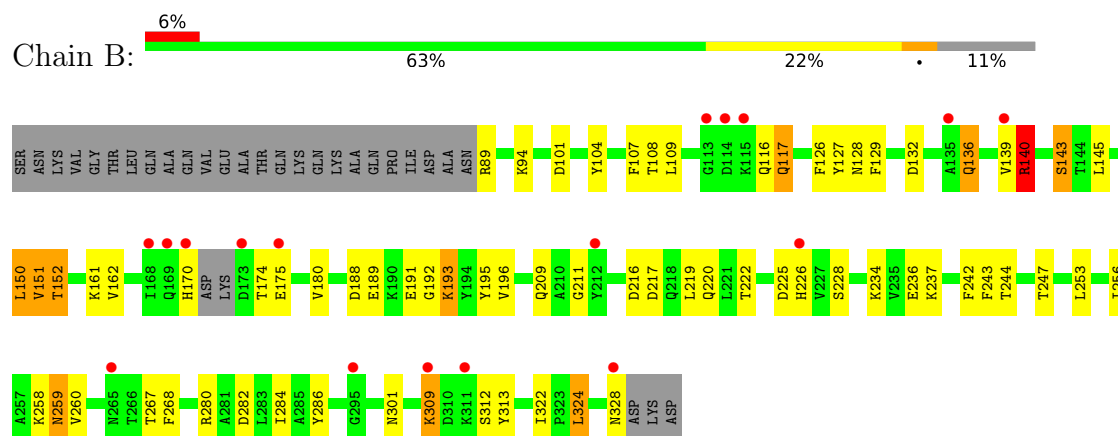
3 Residue-property plots [i](#)

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.

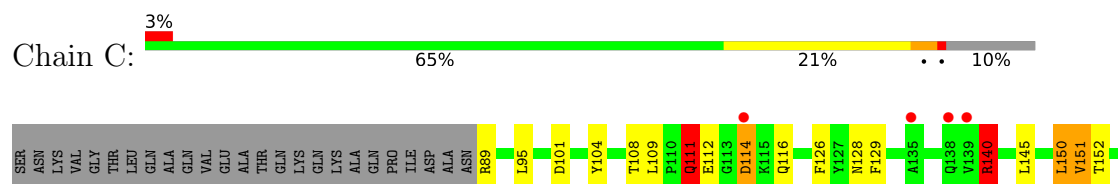
• Molecule 1: Putative transfer protein

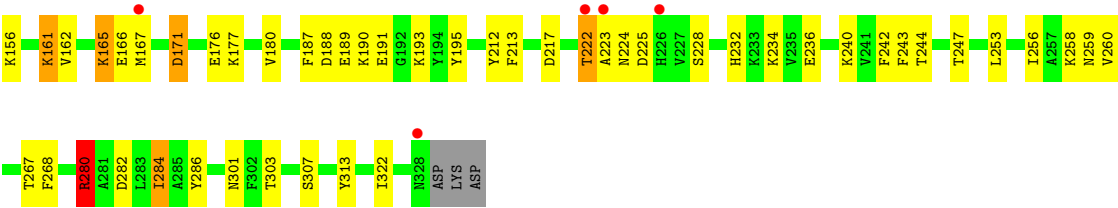


• Molecule 1: Putative transfer protein



• Molecule 1: Putative transfer protein





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	152.74Å 89.19Å 98.42Å 90.00° 112.31° 90.00°	Depositor
Resolution (Å)	91.05 – 2.60 91.05 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.6 (91.05-2.60) 98.6 (91.05-2.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 2.62Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.202 , 0.256 (Not available) , 0.244	Depositor DCC
R_{free} test set	1819 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	60.3	Xtriage
Anisotropy	0.300	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 61.3	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	0.95	EDS
Total number of atoms	11398	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.26% 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 (i)

5.1 Standard geometry (i)

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	0.84	0/2029	1.65	33/2749 (1.2%)
1	B	0.85	0/2004	1.65	33/2716 (1.2%)
1	C	0.76	0/2018	1.58	21/2737 (0.8%)
All	All	0.82	0/6051	1.63	87/8202 (1.1%)

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	3
1	C	0	4
All	All	0	10

There are no bond length outliers.

All (87) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	226	HIS	CB-CA-C	12.63	128.69	111.14
1	A	226	HIS	CB-CA-C	12.40	127.40	111.40
1	A	236	GLU	CB-CG-CD	8.59	127.20	112.60
1	C	111	GLN	CB-CA-C	-8.48	97.81	110.95
1	B	282	ASP	CA-CB-CG	8.35	120.95	112.60
1	B	313	TYR	N-CA-CB	-8.18	97.95	111.20
1	B	170	HIS	CA-CB-CG	8.12	121.92	113.80
1	B	161	LYS	CB-CA-C	7.86	124.09	109.71
1	A	234	LYS	N-CA-CB	-7.77	98.67	110.16
1	C	217	ASP	CA-CB-CG	7.64	120.24	112.60
1	A	269	LYS	CB-CG-CD	7.44	128.41	111.30
1	A	244	THR	CA-CB-OG1	-7.38	98.53	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	174	THR	CA-CB-OG1	-7.30	98.65	109.60
1	C	313	TYR	N-CA-CB	-7.28	99.41	111.20
1	B	222	THR	CA-CB-OG1	-7.24	98.75	109.60
1	A	132	ASP	CA-CB-CG	7.21	119.81	112.60
1	A	313	TYR	N-CA-CB	-7.11	99.68	111.20
1	A	282	ASP	CA-CB-CG	7.07	119.67	112.60
1	B	236	GLU	CB-CG-CD	7.04	124.57	112.60
1	B	313	TYR	CB-CA-C	7.03	123.30	109.66
1	B	244	THR	CA-CB-OG1	-6.96	99.17	109.60
1	C	234	LYS	N-CA-CB	-6.95	99.88	110.16
1	B	117	GLN	N-CA-CB	6.93	120.12	110.07
1	B	132	ASP	CA-CB-CG	6.88	119.48	112.60
1	C	224	ASN	CB-CA-C	6.84	121.70	109.38
1	A	94	LYS	CB-CA-C	-6.83	100.16	110.88
1	A	259	ASN	CB-CA-C	-6.79	102.25	111.89
1	A	114	ASP	CA-CB-CG	6.78	119.38	112.60
1	B	94	LYS	CB-CA-C	-6.76	100.27	110.88
1	C	101	ASP	CA-CB-CG	6.69	119.29	112.60
1	A	161	LYS	CB-CA-C	6.69	121.95	109.71
1	C	313	TYR	CB-CA-C	6.66	122.58	109.66
1	B	236	GLU	N-CA-CB	6.65	119.90	110.12
1	B	139	VAL	N-CA-CB	-6.57	103.53	111.21
1	C	282	ASP	CA-CB-CG	6.54	119.14	112.60
1	C	222	THR	CA-CB-OG1	-6.49	99.86	109.60
1	B	140	ARG	CB-CG-CD	6.46	126.16	111.30
1	B	247	THR	OG1-CB-CG2	6.41	122.13	109.30
1	C	161	LYS	CB-CA-C	6.41	121.44	109.71
1	C	166	GLU	CB-CA-C	-6.39	98.74	109.48
1	A	94	LYS	N-CA-CB	6.34	119.21	110.01
1	B	234	LYS	N-CA-CB	-6.31	100.82	110.16
1	A	140	ARG	CB-CG-CD	6.26	125.70	111.30
1	B	216	ASP	CA-CB-CG	6.25	118.86	112.60
1	A	313	TYR	CB-CA-C	6.20	121.69	109.66
1	C	108	THR	CA-CB-OG1	6.17	118.86	109.60
1	B	193	LYS	CA-CB-CG	-6.16	101.77	114.10
1	B	94	LYS	N-CA-CB	6.14	118.91	110.01
1	B	152	THR	CA-CB-OG1	-6.13	100.40	109.60
1	A	236	GLU	N-CA-CB	6.12	119.12	110.12
1	B	220	GLN	CB-CA-C	6.02	119.68	109.75
1	C	212	TYR	CB-CA-C	6.01	119.52	111.82
1	A	269	LYS	CG-CD-CE	5.94	124.97	111.30
1	B	225	ASP	CB-CA-C	5.91	120.18	109.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	215	GLU	CB-CA-C	-5.77	98.26	110.31
1	C	191	GLU	CB-CA-C	-5.71	103.77	112.11
1	C	114	ASP	CA-CB-CG	5.68	118.28	112.60
1	A	140	ARG	CB-CA-C	-5.65	104.42	111.43
1	A	247	THR	OG1-CB-CG2	5.64	120.59	109.30
1	A	240	LYS	CG-CD-CE	5.64	124.26	111.30
1	A	193	LYS	CA-CB-CG	-5.63	102.83	114.10
1	B	174	THR	CA-CB-OG1	-5.61	101.19	109.60
1	A	235	VAL	N-CA-C	-5.61	105.04	110.42
1	B	222	THR	OG1-CB-CG2	5.60	120.50	109.30
1	A	233	LYS	CB-CG-CD	5.60	124.17	111.30
1	B	174	THR	N-CA-CB	-5.56	100.99	110.16
1	C	217	ASP	CB-CA-C	-5.54	101.20	110.56
1	A	152	THR	CA-CB-OG1	-5.53	101.31	109.60
1	A	234	LYS	CB-CA-C	5.48	120.16	110.85
1	C	152	THR	CA-CB-OG1	-5.46	101.41	109.60
1	A	166	GLU	CB-CG-CD	5.45	121.87	112.60
1	B	108	THR	CA-CB-OG1	5.44	117.76	109.60
1	C	223	ALA	O-C-N	5.43	128.72	123.29
1	B	259	ASN	CB-CA-C	-5.42	104.19	111.89
1	A	151	VAL	N-CA-C	-5.42	107.15	111.81
1	A	225	ASP	CB-CA-C	5.32	119.14	109.83
1	B	101	ASP	CA-CB-CG	5.27	117.87	112.60
1	B	219	LEU	N-CA-CB	5.26	119.35	110.46
1	A	224	ASN	CA-CB-CG	-5.16	107.44	112.60
1	B	191	GLU	CB-CA-C	-5.15	104.58	112.11
1	A	325	ASN	OD1-CG-ND2	-5.13	117.47	122.60
1	B	309	LYS	CB-CG-CD	5.07	122.97	111.30
1	C	236	GLU	N-CA-CB	5.05	117.54	110.12
1	C	140	ARG	CG-CD-NE	5.03	123.07	112.00
1	B	243	PHE	N-CA-CB	5.03	117.51	110.12
1	C	165	LYS	N-CA-CB	5.02	119.57	110.83
1	A	329	ASP	CA-CB-CG	5.02	117.62	112.60

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	140	ARG	Sidechain
1	A	224	ASN	Sidechain
1	A	280	ARG	Sidechain
1	B	136	GLN	Peptide

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Mol	Chain	Res	Type	Group
1	B	140	ARG	Sidechain
1	B	280	ARG	Sidechain
1	C	140	ARG	Sidechain
1	C	222	THR	Peptide
1	C	280	ARG	Sidechain
1	C	89	ARG	Sidechain

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	1983	1840	1898	23	0
1	B	1958	1819	1877	25	1
1	C	1971	1827	1884	33	0
All	All	5912	5486	5659	76	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (76) 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:217:ASP:HA	1:B:324:LEU:HD11	1.61	0.82
1:C:280:ARG:HH11	1:C:280:ARG:HG3	1.45	0.82
1:A:112:GLU:HB3	1:A:114:ASP:OD1	1.93	0.68
1:A:301:ASN:HB3	1:A:322:ILE:HG13	1.77	0.66
1:B:301:ASN:HB3	1:B:322:ILE:HG13	1.79	0.64
1:C:280:ARG:HH11	1:C:280:ARG:CG	2.11	0.62
1:C:301:ASN:HB3	1:C:322:ILE:HG13	1.79	0.62
1:A:109:LEU:HD11	1:A:116:GLN:HG3	1.83	0.61
1:A:144:THR:OG1	1:A:167:MET:CE	2.50	0.60
1:C:189:GLU:HA	1:C:193:LYS:O	2.03	0.58
1:C:150:LEU:HD12	1:C:151:VAL:N	2.19	0.57
1:B:150:LEU:HD12	1:B:151:VAL:N	2.19	0.57
1:A:278:LYS:HE3	1:A:280:ARG:O	2.05	0.57
1:A:117:GLN:HA	1:A:117:GLN:OE1	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:GLU:OE2	1:A:192:GLY:HA2	2.06	0.56
1:B:189:GLU:HA	1:B:193:LYS:O	2.07	0.55
1:C:140:ARG:HG3	1:C:140:ARG:HH21	1.70	0.55
1:C:150:LEU:HD12	1:C:150:LEU:C	2.31	0.55
1:A:151:VAL:HG23	1:A:152:THR:HG23	1.89	0.55
1:B:117:GLN:OE1	1:B:117:GLN:HA	2.06	0.54
1:C:280:ARG:CG	1:C:280:ARG:NH1	2.70	0.53
1:A:189:GLU:HA	1:A:193:LYS:O	2.08	0.53
1:B:217:ASP:CA	1:B:324:LEU:HD11	2.37	0.52
1:A:144:THR:OG1	1:A:167:MET:HE3	2.09	0.52
1:C:240:LYS:O	1:C:244:THR:OG1	2.23	0.51
1:B:189:GLU:OE2	1:B:192:GLY:HA2	2.12	0.49
1:C:167:MET:HE2	1:C:176:GLU:OE2	2.12	0.49
1:C:171:ASP:OD1	1:C:171:ASP:N	2.29	0.49
1:C:128:ASN:OD1	1:C:129:PHE:N	2.45	0.48
1:C:253:LEU:HD23	1:C:260:VAL:O	2.13	0.48
1:A:128:ASN:HA	1:C:213:PHE:CE2	2.48	0.48
1:B:128:ASN:OD1	1:B:129:PHE:N	2.46	0.48
1:B:267:THR:HG22	1:B:268:PHE:O	2.14	0.47
1:B:109:LEU:HD11	1:B:116:GLN:HG3	1.95	0.47
1:C:284:ILE:HG21	1:C:286:TYR:CZ	2.48	0.47
1:B:211:GLY:O	1:C:190:LYS:NZ	2.48	0.47
1:C:242:PHE:HA	1:C:256:ILE:HD11	1.97	0.47
1:B:242:PHE:HA	1:B:256:ILE:HD11	1.96	0.47
1:B:284:ILE:HG21	1:B:286:TYR:CZ	2.51	0.46
1:C:162:VAL:O	1:C:180:VAL:HA	2.16	0.46
1:C:267:THR:HG22	1:C:268:PHE:O	2.15	0.46
1:B:150:LEU:HD12	1:B:150:LEU:C	2.41	0.46
1:B:107:PHE:O	1:B:143:SER:OG	2.30	0.45
1:A:267:THR:HG22	1:A:268:PHE:O	2.17	0.45
1:C:243:PHE:O	1:C:247:THR:HG23	2.17	0.45
1:C:284:ILE:N	1:C:284:ILE:HD13	2.32	0.44
1:B:324:LEU:HD23	1:B:324:LEU:HA	1.83	0.44
1:B:104:TYR:HA	1:B:145:LEU:HD22	2.00	0.44
1:C:111:GLN:O	1:C:112:GLU:C	2.60	0.43
1:A:188:ASP:HB3	1:A:195:TYR:CZ	2.54	0.43
1:A:278:LYS:CE	1:A:280:ARG:O	2.66	0.43
1:C:109:LEU:HD11	1:C:116:GLN:HG3	2.01	0.43
1:A:136:GLN:HA	1:A:136:GLN:OE1	2.19	0.43
1:A:209:GLN:OE1	1:B:126:PHE:HA	2.19	0.43
1:B:162:VAL:O	1:B:180:VAL:HA	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:258:LYS:O	1:C:259:ASN:C	2.61	0.42
1:B:253:LEU:HD23	1:B:260:VAL:O	2.18	0.42
1:C:286:TYR:CE1	1:C:303:THR:HG23	2.54	0.42
1:C:104:TYR:HA	1:C:145:LEU:HD22	2.01	0.42
1:C:225:ASP:CG	1:C:232:HIS:HE2	2.26	0.42
1:B:209:GLN:OE1	1:C:126:PHE:HA	2.20	0.42
1:C:188:ASP:HB3	1:C:195:TYR:CZ	2.55	0.41
1:A:128:ASN:OD1	1:A:129:PHE:N	2.52	0.41
1:B:151:VAL:HG23	1:B:152:THR:HG23	2.01	0.41
1:B:127:TYR:HA	1:B:196:VAL:O	2.20	0.41
1:A:90:LYS:HD3	1:C:95:LEU:HD23	2.02	0.41
1:B:188:ASP:HB3	1:B:195:TYR:CZ	2.56	0.41
1:A:109:LEU:HD22	1:A:140:ARG:HD2	2.02	0.41
1:A:269:LYS:C	1:A:270:THR:HG23	2.46	0.41
1:A:162:VAL:O	1:A:180:VAL:HA	2.21	0.41
1:C:111:GLN:H	1:C:111:GLN:HG2	1.62	0.41
1:A:168:ILE:HG13	1:A:175:GLU:HB2	2.02	0.41
1:B:258:LYS:O	1:B:259:ASN:C	2.64	0.40
1:C:156:LYS:HB3	1:C:187:PHE:O	2.22	0.40
1:A:286:TYR:CE1	1:A:303:THR:HG23	2.57	0.40
1:C:150:LEU:C	1:C:150:LEU:CD1	2.94	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:175:GLU:OE2	1:B:175:GLU:OE2[2_555]	2.19	0.01

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	237/268 (88%)	219 (92%)	17 (7%)	1 (0%)	30	51
1	B	234/268 (87%)	219 (94%)	15 (6%)	0	100	100
1	C	238/268 (89%)	224 (94%)	14 (6%)	0	100	100
All	All	709/804 (88%)	662 (93%)	46 (6%)	1 (0%)	48	70

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	228	SER

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	220/242 (91%)	207 (94%)	13 (6%)	18	38
1	B	217/242 (90%)	205 (94%)	12 (6%)	19	41
1	C	218/242 (90%)	205 (94%)	13 (6%)	17	37
All	All	655/726 (90%)	617 (94%)	38 (6%)	18	39

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	111	GLN
1	A	114	ASP
1	A	139	VAL
1	A	140	ARG
1	A	151	VAL
1	A	168	ILE
1	A	228	SER
1	A	229	ASP
1	A	269	LYS
1	A	280	ARG
1	A	307	SER
1	A	311	LYS

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Mol	Chain	Res	Type
1	A	312	SER
1	B	89	ARG
1	B	136	GLN
1	B	140	ARG
1	B	143	SER
1	B	150	LEU
1	B	151	VAL
1	B	228	SER
1	B	237	LYS
1	B	309	LYS
1	B	312	SER
1	B	324	LEU
1	B	328	ASN
1	C	111	GLN
1	C	114	ASP
1	C	140	ARG
1	C	150	LEU
1	C	151	VAL
1	C	161	LYS
1	C	165	LYS
1	C	171	ASP
1	C	177	LYS
1	C	228	SER
1	C	280	ARG
1	C	284	ILE
1	C	307	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	96	GLN
1	A	116	GLN
1	B	136	GLN
1	B	328	ASN
1	C	96	GLN
1	C	116	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](#)

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	241/268 (89%)	0.40	8 (3%)	49	43	61, 81, 119, 153	0
1	B	238/268 (88%)	0.57	17 (7%)	22	17	61, 85, 131, 152	0
1	C	240/268 (89%)	0.10	9 (3%)	44	38	63, 82, 125, 158	0
All	All	719/804 (89%)	0.35	34 (4%)	36	30	61, 82, 125, 158	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	170	HIS	4.9
1	B	170	HIS	4.7
1	B	169	GLN	4.3
1	C	139	VAL	4.2
1	B	113	GLY	3.4
1	B	168	ILE	3.4
1	C	328	ASN	3.4
1	A	255	LEU	3.3
1	B	328	ASN	3.2
1	A	173	ASP	3.2
1	A	311	LYS	3.1
1	A	226	HIS	3.1
1	B	139	VAL	3.0
1	C	222	THR	2.9
1	B	226	HIS	2.9
1	B	311	LYS	2.8
1	A	139	VAL	2.8
1	A	167	MET	2.7
1	B	212	TYR	2.7
1	C	226	HIS	2.6
1	B	115	LYS	2.5
1	B	173	ASP	2.4
1	A	168	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	138	GLN	2.3
1	C	114	ASP	2.3
1	B	135	ALA	2.2
1	C	135	ALA	2.1
1	B	175	GLU	2.1
1	B	309	LYS	2.1
1	B	295	GLY	2.1
1	B	114	ASP	2.1
1	C	223	ALA	2.1
1	C	167	MET	2.1
1	B	265	ASN	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.