



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:02 AM UTC

PDB ID : 3TPF / pdb_00003tpf
Title : Crystal structure of anabolic ornithine carbamoyltransferase from *Campylobacter jejuni* subsp. *jejuni* NCTC 11168
Authors : Shabalin, I.G.; Onopriyenko, O.; Grimshaw, S.; Porebski, P.J.; Grabowski, M.; Savchenko, A.; Chruszcz, M.; Anderson, W.F.; Minor, W.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2011-09-07
Resolution : 2.70 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
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (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

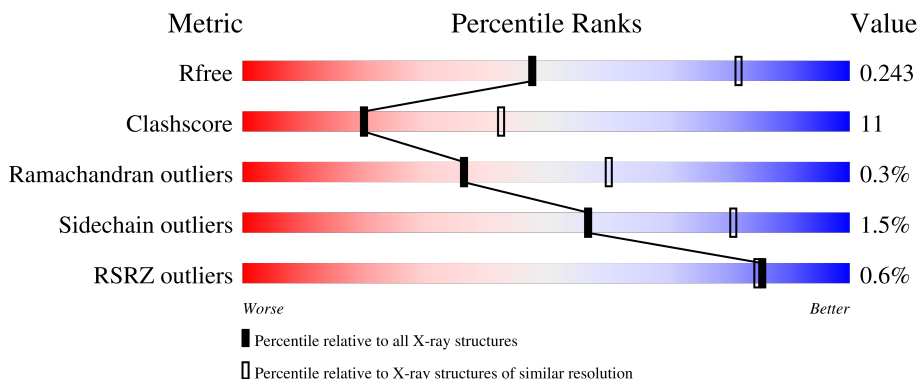
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


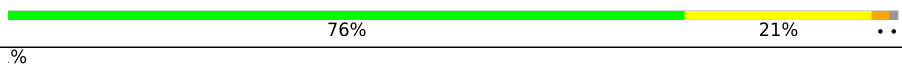
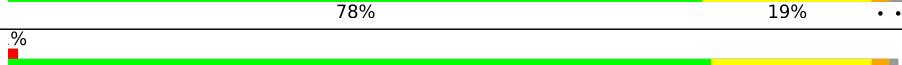

The reported resolution of this entry is 2.70 Å.

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.



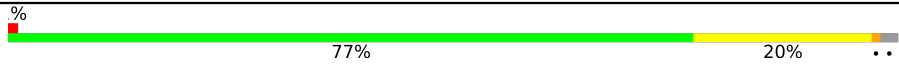

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	307	 79% 18% ..
1	B	307	 76% 21% ..
1	C	307	 78% 19% ..
1	D	307	 79% 18% ..

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Mol	Chain	Length	Quality of chain
1	E	307	 <p>% 77% 20% ..</p>
1	F	307	 <p>% 80% 15% ..</p>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 13981 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 Ornithine carbamoyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	303	2334	1503	385	430	3	13	0	0	0
1	B	305	2358	1523	383	437	3	12	0	0	0
1	C	301	2284	1469	380	420	3	12	0	0	0
1	D	303	2305	1481	376	432	3	13	0	0	0
1	E	300	2281	1470	371	425	3	12	0	0	0
1	F	298	2294	1481	375	423	3	12	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q9PNU6
B	0	GLY	-	expression tag	UNP Q9PNU6
C	0	GLY	-	expression tag	UNP Q9PNU6
D	0	GLY	-	expression tag	UNP Q9PNU6
E	0	GLY	-	expression tag	UNP Q9PNU6
F	0	GLY	-	expression tag	UNP Q9PNU6

- Molecule 2 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 7 4 3	0	0
2	B	1	Total C O 7 4 3	0	0
2	F	1	Total C O 7 4 3	0	0

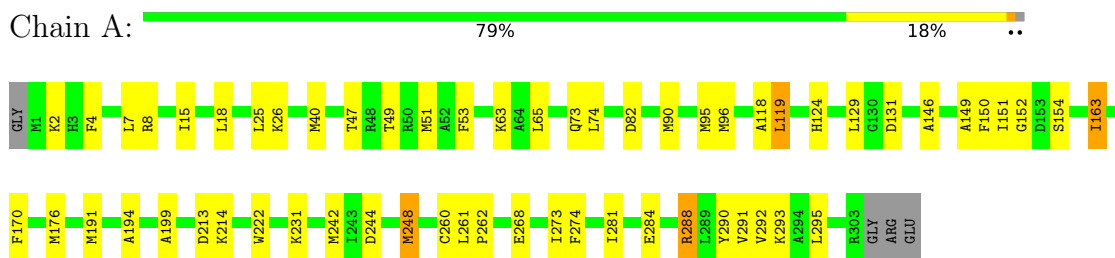
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	15	Total O 15 15	0	0
3	B	32	Total O 32 32	0	0
3	C	16	Total O 16 16	0	0
3	D	13	Total O 13 13	0	0
3	E	14	Total O 14 14	0	0
3	F	14	Total O 14 14	0	0

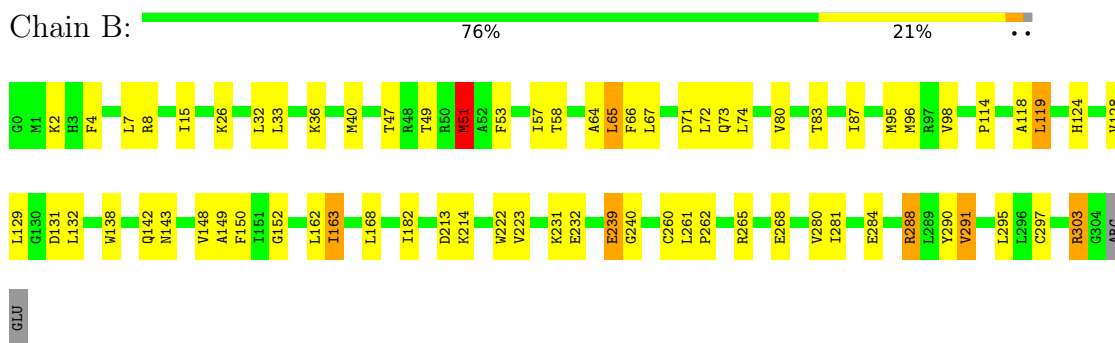
3 Residue-property plots [i](#)

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.

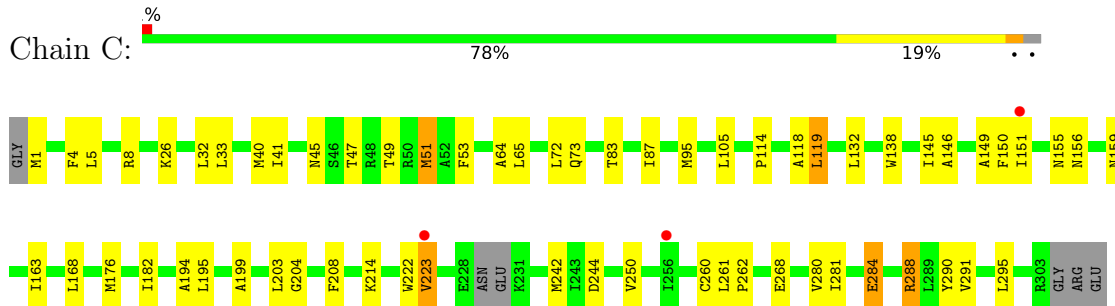
- Molecule 1: Ornithine carbamoyltransferase



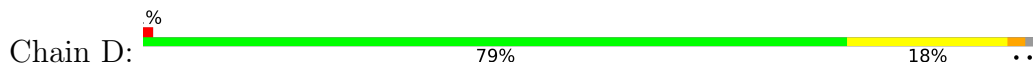
- Molecule 1: Ornithine carbamoyltransferase



- Molecule 1: Ornithine carbamoyltransferase



- Molecule 1: Ornithine carbamoyltransferase



4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	157.16Å 87.40Å 141.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.70 50.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.00-2.70) 99.8 (50.00-2.70)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.99 (at 2.69Å)	Xtrriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.202 , 0.240 0.204 , 0.243	Depositor DCC
R_{free} test set	2758 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	59.8	Xtrriage
Anisotropy	0.771	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 59.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13981	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.40 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.3286e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality i

5.1 Standard geometry i

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

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	1.18	3/2366 (0.1%)	1.15	4/3182 (0.1%)
1	B	1.32	4/2389 (0.2%)	1.23	7/3206 (0.2%)
1	C	1.13	1/2312 (0.0%)	1.13	2/3108 (0.1%)
1	D	1.17	2/2337 (0.1%)	1.13	6/3149 (0.2%)
1	E	1.08	1/2311 (0.0%)	1.12	2/3112 (0.1%)
1	F	1.19	2/2325 (0.1%)	1.14	2/3125 (0.1%)
All	All	1.18	13/14040 (0.1%)	1.15	23/18882 (0.1%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	65	LEU	CA-C	-8.65	1.42	1.52
1	A	248	MSE	SE-CE	6.60	2.15	1.95
1	A	65	LEU	CA-C	-6.25	1.45	1.52
1	F	65	LEU	CA-C	-5.34	1.46	1.52
1	B	128	VAL	CA-CB	-5.30	1.47	1.54
1	B	114	PRO	CA-C	-5.27	1.46	1.52
1	D	37	THR	CA-C	-5.24	1.46	1.52
1	F	66	PHE	C-O	-5.24	1.17	1.24
1	B	36	LYS	N-CA	-5.11	1.39	1.45
1	A	63	LYS	N-CA	5.09	1.52	1.46
1	D	114	PRO	CA-C	-5.05	1.46	1.52
1	E	96	MSE	CA-C	-5.03	1.46	1.52
1	C	114	PRO	C-O	-5.03	1.17	1.23

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	163	ILE	CB-CA-C	-6.47	103.69	111.97
1	B	152	GLY	N-CA-C	6.24	119.00	110.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	148	VAL	CB-CA-C	-5.94	101.78	110.62
1	B	163	ILE	CB-CA-C	-5.86	104.47	111.97
1	D	152	GLY	N-CA-C	5.84	118.47	110.58
1	E	96	MSE	CG-SE-CE	5.84	111.77	98.92
1	A	231	LYS	N-CA-C	5.83	117.31	111.07
1	F	163	ILE	N-CA-C	5.79	115.98	110.42
1	C	51	MSE	CG-SE-CE	-5.77	86.23	98.92
1	D	231	LYS	N-CA-C	5.76	117.23	111.07
1	B	232	GLU	CA-CB-CG	5.67	125.43	114.10
1	B	303	ARG	N-CA-C	5.58	119.40	112.59
1	A	152	GLY	N-CA-C	5.57	118.09	110.58
1	A	163	ILE	CB-CA-C	-5.54	104.88	111.97
1	B	148	VAL	N-CA-C	5.49	116.39	108.48
1	D	157	MSE	CG-SE-CE	-5.49	86.84	98.92
1	D	148	VAL	N-CA-C	5.37	116.21	108.48
1	B	98	VAL	N-CA-C	5.25	116.94	108.85
1	E	152	GLY	N-CA-C	5.20	117.60	110.58
1	A	90	MSE	CA-CB-CG	-5.18	103.73	114.10
1	D	163	ILE	CB-CA-C	-5.08	105.47	111.97
1	B	51	MSE	CG-SE-CE	5.07	110.07	98.92
1	C	284	GLU	N-CA-C	-5.02	104.90	111.02

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	2334	0	2270	50	0
1	B	2358	0	2307	60	0
1	C	2284	0	2219	59	0
1	D	2305	0	2188	53	0
1	E	2281	0	2182	60	0
1	F	2294	0	2223	50	0
2	A	7	0	10	0	0
2	B	7	0	10	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	7	0	10	0	0
3	A	15	0	0	0	0
3	B	32	0	0	2	0
3	C	16	0	0	0	0
3	D	13	0	0	1	0
3	E	14	0	0	1	0
3	F	14	0	0	0	0
All	All	13981	0	13419	312	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (312) 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:248:MSE:SE	1:A:248:MSE:CE	2.15	1.44
1:D:53:PHE:CD2	1:D:95:MSE:CE	2.31	1.13
1:D:53:PHE:CD2	1:D:95:MSE:HE2	1.91	1.03
1:E:53:PHE:CD2	1:E:95:MSE:CE	2.42	1.03
1:F:53:PHE:CD2	1:F:95:MSE:HE2	1.94	1.02
1:B:53:PHE:CD2	1:B:95:MSE:CE	2.50	0.94
1:A:53:PHE:CD2	1:A:95:MSE:CE	2.50	0.93
1:D:53:PHE:CG	1:D:95:MSE:CE	2.52	0.92
1:C:203:LEU:HD12	1:C:204:GLY:N	1.86	0.91
1:C:53:PHE:CD2	1:C:95:MSE:CE	2.52	0.91
1:F:65:LEU:N	1:F:65:LEU:HD12	1.85	0.90
1:D:47:THR:HG21	1:E:72:LEU:HD13	1.54	0.89
1:C:83:THR:HG22	1:C:87:ILE:HD12	1.57	0.86
1:E:53:PHE:CG	1:E:95:MSE:HE3	2.11	0.86
1:B:53:PHE:CD2	1:B:95:MSE:HE2	2.12	0.83
1:D:53:PHE:CE2	1:D:95:MSE:HE2	2.13	0.82
1:D:53:PHE:CG	1:D:95:MSE:HE2	2.14	0.81
1:F:53:PHE:CD2	1:F:95:MSE:CE	2.64	0.81
1:F:53:PHE:CG	1:F:95:MSE:HE2	2.15	0.81
1:C:73:GLN:HG2	1:C:73:GLN:O	1.80	0.81
1:B:223:VAL:HG21	1:B:231:LYS:HG3	1.62	0.80
1:E:53:PHE:CD2	1:E:95:MSE:HE3	2.15	0.80
1:F:4:PHE:CG	1:F:295:LEU:HD11	2.17	0.80
1:A:53:PHE:CG	1:A:95:MSE:CE	2.68	0.77
1:E:53:PHE:CE2	1:E:95:MSE:HE2	2.20	0.77
1:C:53:PHE:CG	1:C:95:MSE:CE	2.68	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:203:LEU:HD12	1:C:204:GLY:H	1.48	0.77
1:B:53:PHE:CG	1:B:95:MSE:CE	2.67	0.76
1:F:64:ALA:C	1:F:65:LEU:HD12	2.11	0.76
1:D:53:PHE:CG	1:D:95:MSE:HE3	2.19	0.76
1:C:65:LEU:HD12	1:C:65:LEU:N	2.00	0.76
1:C:53:PHE:CD2	1:C:95:MSE:HE2	2.22	0.75
1:A:53:PHE:CD2	1:A:95:MSE:HE2	2.21	0.73
1:C:242:MSE:HE3	1:C:244:ASP:HB3	1.69	0.73
1:B:80:VAL:HG23	3:B:331:HOH:O	1.88	0.72
1:D:83:THR:HG22	1:D:87:ILE:HD12	1.72	0.71
1:B:53:PHE:CE2	1:B:95:MSE:HE2	2.26	0.71
1:D:40:MSE:HE2	1:D:64:ALA:HB1	1.73	0.70
1:A:4:PHE:CG	1:A:295:LEU:HD11	2.26	0.70
1:A:4:PHE:CD1	1:A:295:LEU:HD11	2.29	0.68
1:A:51:MSE:HE1	1:C:87:ILE:HG12	1.75	0.68
1:C:132:LEU:HD21	1:C:168:LEU:HD11	1.73	0.68
1:B:239:GLU:HG3	1:B:240:GLY:H	1.57	0.68
1:C:40:MSE:HA	1:C:95:MSE:HG3	1.76	0.68
1:F:4:PHE:CD1	1:F:295:LEU:HD11	2.29	0.68
1:D:53:PHE:CD2	1:D:95:MSE:HE1	2.24	0.67
1:E:53:PHE:CE2	1:E:95:MSE:CE	2.77	0.67
1:B:40:MSE:HA	1:B:95:MSE:HG3	1.76	0.67
1:A:129:LEU:HB3	1:A:291:VAL:HG11	1.75	0.67
1:C:64:ALA:C	1:C:65:LEU:HD12	2.19	0.67
1:D:150:PHE:C	1:D:151:ILE:HD12	2.20	0.67
1:B:53:PHE:CG	1:B:95:MSE:HE3	2.29	0.67
1:E:4:PHE:CG	1:E:295:LEU:HD11	2.29	0.67
1:C:4:PHE:CD1	1:C:295:LEU:HD11	2.31	0.66
1:C:83:THR:CG2	1:C:87:ILE:HD12	2.26	0.66
1:A:53:PHE:CG	1:A:95:MSE:HE3	2.31	0.65
1:F:53:PHE:CG	1:F:95:MSE:CE	2.78	0.65
1:D:4:PHE:CD1	1:D:295:LEU:HD11	2.32	0.65
1:A:40:MSE:HA	1:A:95:MSE:HG3	1.78	0.65
1:B:239:GLU:HG3	1:B:240:GLY:N	2.12	0.65
1:E:53:PHE:CD2	1:E:95:MSE:HE2	2.27	0.65
1:E:53:PHE:CG	1:E:95:MSE:CE	2.76	0.65
1:C:41:ILE:HG23	1:C:72:LEU:HD23	1.79	0.65
1:E:83:THR:HG22	1:E:87:ILE:HD12	1.78	0.64
1:E:279:ASP:HB2	3:E:315:HOH:O	1.96	0.64
1:E:45:ASN:HD22	1:F:71:ASP:HA	1.62	0.64
1:E:4:PHE:CD1	1:E:295:LEU:HD11	2.32	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:THR:HG21	1:C:72:LEU:HD13	1.80	0.64
1:B:66:PHE:C	1:B:67:LEU:HD23	2.23	0.64
1:E:149:ALA:HB2	1:E:214:LYS:HG3	1.80	0.64
1:C:53:PHE:CD2	1:C:95:MSE:HE1	2.34	0.63
1:B:53:PHE:CG	1:B:95:MSE:HE2	2.33	0.63
1:B:132:LEU:HD21	1:B:168:LEU:HD11	1.80	0.63
1:F:65:LEU:N	1:F:65:LEU:CD1	2.59	0.63
1:A:53:PHE:CD2	1:A:95:MSE:HE1	2.33	0.63
1:C:53:PHE:CG	1:C:95:MSE:HE3	2.34	0.63
1:E:203:LEU:HD12	1:E:204:GLY:N	2.15	0.62
1:B:87:ILE:HG12	1:C:51:MSE:HE1	1.82	0.61
1:D:4:PHE:CG	1:D:295:LEU:HD11	2.35	0.61
1:C:40:MSE:HG2	1:C:95:MSE:HG2	1.82	0.60
1:B:72:LEU:HD13	1:C:47:THR:HG21	1.82	0.60
1:C:149:ALA:HB2	1:C:214:LYS:HG3	1.83	0.60
1:B:32:LEU:HD12	1:B:297:CYS:SG	2.42	0.60
1:E:132:LEU:HD21	1:E:168:LEU:HD11	1.84	0.60
1:F:118:ALA:O	1:F:119:LEU:HB2	2.02	0.60
1:C:4:PHE:CG	1:C:295:LEU:HD11	2.37	0.60
1:D:53:PHE:CB	1:D:95:MSE:HE3	2.32	0.60
1:B:40:MSE:HG2	1:B:95:MSE:HG2	1.84	0.59
1:B:83:THR:HG22	1:B:87:ILE:HD12	1.85	0.58
1:B:57:ILE:HG23	1:B:58:THR:N	2.18	0.58
1:C:53:PHE:CE2	1:C:95:MSE:HE2	2.38	0.58
1:B:4:PHE:CD1	1:B:295:LEU:HD11	2.38	0.58
1:D:40:MSE:HA	1:D:95:MSE:HG3	1.85	0.58
1:A:51:MSE:CE	1:C:87:ILE:HG12	2.33	0.58
1:C:53:PHE:CG	1:C:95:MSE:HE2	2.38	0.58
1:D:83:THR:CG2	1:D:87:ILE:HD12	2.33	0.58
1:F:47:THR:HG22	1:F:51:MSE:HE2	1.85	0.58
1:A:40:MSE:HG2	1:A:95:MSE:HG2	1.87	0.57
1:A:53:PHE:CE2	1:A:95:MSE:HE2	2.38	0.57
1:C:118:ALA:O	1:C:119:LEU:HB2	2.05	0.57
1:D:51:MSE:CE	1:E:87:ILE:HG12	2.35	0.57
1:E:65:LEU:HD12	1:E:65:LEU:N	2.20	0.57
1:F:149:ALA:HB2	1:F:214:LYS:HG3	1.86	0.57
1:E:50:ARG:NH1	1:F:66:PHE:O	2.37	0.57
1:A:53:PHE:CG	1:A:95:MSE:HE2	2.39	0.57
1:D:154:SER:HA	1:D:176:MSE:HE1	1.85	0.57
1:E:118:ALA:O	1:E:119:LEU:HB2	2.05	0.57
1:E:40:MSE:HA	1:E:95:MSE:HG3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:ILE:HG12	1:C:51:MSE:CE	2.36	0.56
1:B:118:ALA:O	1:B:119:LEU:HB2	2.05	0.56
1:D:82:ASP:OD2	1:F:265:ARG:NH2	2.38	0.56
1:E:51:MSE:CE	1:F:87:ILE:HG12	2.35	0.56
1:F:40:MSE:HA	1:F:95:MSE:HG3	1.88	0.56
1:F:8:ARG:HG3	1:F:163:ILE:HD13	1.87	0.56
1:A:7:LEU:HD22	1:A:15:ILE:HG12	1.88	0.55
1:A:129:LEU:HB3	1:A:291:VAL:CG1	2.36	0.55
1:C:182:ILE:HD12	1:C:203:LEU:HD22	1.89	0.55
1:F:49:THR:HG22	1:F:95:MSE:HE1	1.88	0.55
1:A:150:PHE:C	1:A:151:ILE:HD12	2.32	0.55
1:F:53:PHE:CE2	1:F:95:MSE:HE2	2.40	0.55
1:B:260:CYS:O	1:B:261:LEU:HB2	2.07	0.54
1:A:118:ALA:O	1:A:119:LEU:HB2	2.08	0.54
1:E:45:ASN:ND2	1:F:71:ASP:OD1	2.40	0.54
1:E:119:LEU:HD12	1:E:124:HIS:CD2	2.42	0.54
1:B:4:PHE:CG	1:B:295:LEU:HD11	2.42	0.54
1:C:32:LEU:O	1:C:33:LEU:HD23	2.08	0.54
1:E:119:LEU:HD11	1:E:156:ASN:HD21	1.73	0.54
1:C:8:ARG:HG3	1:C:163:ILE:HD13	1.90	0.54
1:C:65:LEU:N	1:C:65:LEU:CD1	2.71	0.54
1:C:41:ILE:CG2	1:C:72:LEU:HD23	2.38	0.54
1:E:261:LEU:N	1:E:261:LEU:HD23	2.22	0.53
1:F:222:TRP:CZ2	1:F:268:GLU:HA	2.43	0.53
1:B:65:LEU:N	1:B:65:LEU:HD12	2.23	0.53
1:B:138:TRP:CE2	1:B:280:VAL:HG23	2.43	0.53
1:B:149:ALA:HB2	1:B:214:LYS:HG3	1.90	0.53
1:E:150:PHE:C	1:E:151:ILE:HD12	2.34	0.53
1:C:222:TRP:CZ2	1:C:268:GLU:HA	2.43	0.53
1:A:74:LEU:HD11	1:A:96:MSE:CE	2.38	0.53
1:D:98:VAL:HG21	1:D:104:LEU:HG	1.91	0.53
1:D:149:ALA:HB2	1:D:214:LYS:HG3	1.89	0.53
1:D:40:MSE:HG2	1:D:95:MSE:HG2	1.91	0.53
1:A:154:SER:HA	1:A:176:MSE:HE1	1.89	0.52
1:E:203:LEU:HD12	1:E:204:GLY:H	1.73	0.52
1:F:7:LEU:HD22	1:F:15:ILE:HG12	1.91	0.52
1:B:213:ASP:CG	1:B:213:ASP:O	2.53	0.52
1:E:119:LEU:HD12	1:E:124:HIS:CG	2.45	0.52
1:B:71:ASP:HA	1:C:45:ASN:HD22	1.75	0.51
1:E:222:TRP:CZ2	1:E:268:GLU:HA	2.46	0.51
1:C:260:CYS:O	1:C:261:LEU:HB2	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:ARG:HG3	1:B:163:ILE:HD13	1.92	0.51
1:A:73:GLN:O	1:A:73:GLN:HG3	2.11	0.50
1:A:129:LEU:CB	1:A:291:VAL:HG11	2.40	0.50
1:D:222:TRP:CZ2	1:D:268:GLU:HA	2.45	0.50
1:A:149:ALA:HB2	1:A:214:LYS:HG3	1.94	0.50
1:E:260:CYS:O	1:E:261:LEU:HB2	2.12	0.50
1:D:118:ALA:O	1:D:119:LEU:HB2	2.12	0.50
1:B:83:THR:CG2	1:B:87:ILE:HD12	2.42	0.50
1:F:73:GLN:O	1:F:73:GLN:HG3	2.11	0.50
1:A:18:LEU:HD13	1:A:129:LEU:HD11	1.93	0.50
1:C:208:PHE:HD2	1:C:250:VAL:HG21	1.77	0.50
1:D:223:VAL:HG12	1:D:223:VAL:O	2.12	0.50
1:B:64:ALA:C	1:B:65:LEU:HD12	2.37	0.49
1:C:138:TRP:CE2	1:C:280:VAL:HG23	2.48	0.49
1:B:47:THR:HG22	1:B:51:MSE:CE	2.41	0.49
1:E:73:GLN:HG3	1:E:73:GLN:O	2.12	0.49
1:D:26:LYS:HD2	1:D:290:TYR:CE2	2.47	0.49
1:F:119:LEU:HD12	1:F:124:HIS:CD2	2.47	0.49
1:B:67:LEU:HD23	1:B:67:LEU:N	2.27	0.49
1:E:40:MSE:HG2	1:E:95:MSE:HG2	1.95	0.49
1:F:40:MSE:HG2	1:F:95:MSE:HG2	1.94	0.49
1:A:191:MSE:HE3	1:A:191:MSE:HA	1.94	0.49
1:C:145:ILE:CG2	1:C:146:ALA:N	2.74	0.49
1:F:213:ASP:CG	1:F:213:ASP:O	2.56	0.48
1:A:194:ALA:CB	1:A:199:ALA:HB3	2.43	0.48
1:D:207:LYS:O	1:D:211:LEU:HD12	2.13	0.48
1:F:260:CYS:O	1:F:261:LEU:HB2	2.13	0.48
1:A:260:CYS:O	1:A:261:LEU:HB2	2.14	0.48
1:B:26:LYS:HD2	1:B:290:TYR:CE2	2.49	0.47
1:F:119:LEU:HD12	1:F:124:HIS:CG	2.49	0.47
1:D:260:CYS:O	1:D:261:LEU:HB2	2.14	0.47
1:E:8:ARG:HG3	1:E:163:ILE:HD13	1.95	0.47
1:B:303:ARG:HB2	3:B:327:HOH:O	2.14	0.47
1:C:118:ALA:O	1:C:119:LEU:CB	2.62	0.47
1:B:53:PHE:CD2	1:B:95:MSE:HE1	2.45	0.47
1:D:129:LEU:HB3	1:D:291:VAL:HG11	1.95	0.47
1:C:41:ILE:HG23	1:C:72:LEU:CD2	2.43	0.47
1:C:223:VAL:HG12	1:C:223:VAL:O	2.13	0.47
1:A:119:LEU:HD12	1:A:124:HIS:CD2	2.49	0.47
1:A:222:TRP:CZ2	1:A:268:GLU:HA	2.49	0.47
1:B:222:TRP:CZ2	1:B:268:GLU:HA	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:51:MSE:HE2	1:D:51:MSE:HB2	1.82	0.47
1:E:64:ALA:C	1:E:65:LEU:HD12	2.40	0.47
1:B:261:LEU:HB3	1:B:262:PRO:HA	1.97	0.47
1:B:7:LEU:HD22	1:B:15:ILE:HG12	1.97	0.46
1:E:129:LEU:HB3	1:E:291:VAL:HG11	1.97	0.46
1:A:26:LYS:HD2	1:A:290:TYR:CE2	2.50	0.46
1:B:53:PHE:CB	1:B:95:MSE:HE3	2.45	0.46
1:B:118:ALA:O	1:B:119:LEU:CB	2.63	0.46
1:F:132:LEU:HD21	1:F:168:LEU:HD11	1.97	0.46
1:B:162:LEU:HD22	1:B:182:ILE:HD13	1.98	0.46
1:F:118:ALA:O	1:F:119:LEU:CB	2.64	0.46
1:C:95:MSE:HG3	1:C:95:MSE:O	2.16	0.46
1:D:49:THR:OG1	1:D:288:ARG:NH1	2.49	0.46
1:D:136:LYS:CA	1:D:141:MSE:HE3	2.47	0.45
1:A:273:ILE:CG2	1:A:274:PHE:N	2.78	0.45
1:F:4:PHE:CD1	1:F:295:LEU:CD1	2.98	0.45
1:B:74:LEU:HD11	1:B:96:MSE:CE	2.46	0.45
1:C:53:PHE:CB	1:C:95:MSE:HE3	2.47	0.45
1:D:242:MSE:HE3	1:D:244:ASP:HB3	1.98	0.45
1:D:48:ARG:NH1	1:E:78:GLU:OE1	2.48	0.45
1:E:51:MSE:HE1	1:F:87:ILE:HG12	1.98	0.45
1:F:40:MSE:HE2	1:F:64:ALA:HB1	1.99	0.45
1:F:148:VAL:HG22	1:F:216:VAL:HB	1.99	0.45
1:D:57:ILE:HG23	1:D:58:THR:N	2.31	0.45
1:D:87:ILE:HG12	1:F:51:MSE:CE	2.47	0.45
1:B:129:LEU:HB3	1:B:291:VAL:CG1	2.47	0.45
1:D:48:ARG:HB3	3:D:318:HOH:O	2.16	0.45
1:D:281:ILE:O	1:D:284:GLU:HB3	2.17	0.45
1:E:26:LYS:HD2	1:E:290:TYR:CE2	2.51	0.45
1:A:281:ILE:O	1:A:284:GLU:HB3	2.17	0.45
1:B:57:ILE:CG2	1:B:58:THR:N	2.79	0.45
1:B:142:GLN:O	1:B:143:ASN:CB	2.66	0.44
1:E:49:THR:CG2	1:E:95:MSE:HE1	2.47	0.44
1:D:53:PHE:HB2	1:D:95:MSE:HE3	1.99	0.44
1:B:129:LEU:HB3	1:B:291:VAL:HG11	1.98	0.44
1:E:38:LEU:CD1	1:E:93:PHE:HB2	2.48	0.44
1:C:203:LEU:HD12	1:C:203:LEU:C	2.40	0.44
1:D:129:LEU:HB3	1:D:291:VAL:CG1	2.48	0.44
1:C:26:LYS:HD2	1:C:290:TYR:CE2	2.52	0.44
1:F:131:ASP:OD1	1:F:284:GLU:OE2	2.35	0.44
1:A:25:LEU:HD13	1:A:293:LYS:HB3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:154:SER:CA	1:D:176:MSE:HE1	2.48	0.43
1:D:213:ASP:CG	1:D:213:ASP:O	2.60	0.43
1:E:129:LEU:HB3	1:E:291:VAL:CG1	2.47	0.43
1:A:8:ARG:HG3	1:A:163:ILE:HD13	1.99	0.43
1:C:155:ASN:O	1:C:159:ASN:ND2	2.52	0.43
1:B:131:ASP:OD1	1:B:284:GLU:OE2	2.36	0.43
1:A:273:ILE:HD12	1:A:273:ILE:HA	1.89	0.43
1:E:118:ALA:O	1:E:119:LEU:CB	2.66	0.43
1:A:53:PHE:CB	1:A:95:MSE:HE3	2.49	0.43
1:A:146:ALA:HB3	1:A:170:PHE:CE2	2.54	0.43
1:B:49:THR:OG1	1:B:288:ARG:NH1	2.50	0.43
1:B:66:PHE:O	1:B:67:LEU:HD23	2.18	0.43
1:C:5:LEU:HD13	1:C:105:LEU:HD23	2.01	0.43
1:E:53:PHE:CD2	1:E:95:MSE:HE1	2.45	0.43
1:F:33:LEU:O	1:F:34:GLN:C	2.61	0.43
1:C:156:ASN:HA	1:C:159:ASN:HD22	1.83	0.42
1:C:194:ALA:CB	1:C:199:ALA:HB3	2.49	0.42
1:B:47:THR:HG22	1:B:51:MSE:HE2	2.02	0.42
1:B:32:LEU:O	1:B:33:LEU:HD23	2.19	0.42
1:B:150:PHE:CD2	1:B:150:PHE:C	2.97	0.42
1:B:239:GLU:CG	1:B:240:GLY:N	2.81	0.42
1:D:131:ASP:OD1	1:D:284:GLU:OE2	2.38	0.42
1:C:261:LEU:HB3	1:C:262:PRO:HA	2.01	0.42
1:E:261:LEU:HB3	1:E:262:PRO:HA	2.01	0.42
1:F:26:LYS:HD2	1:F:290:TYR:CE2	2.55	0.42
1:A:261:LEU:HB3	1:A:262:PRO:HA	2.02	0.42
1:E:53:PHE:CZ	1:E:95:MSE:HE2	2.53	0.42
1:C:150:PHE:C	1:C:151:ILE:HD12	2.44	0.42
1:D:261:LEU:HB3	1:D:262:PRO:HA	2.02	0.42
1:F:261:LEU:HB3	1:F:262:PRO:HA	2.02	0.42
1:A:118:ALA:O	1:A:119:LEU:CB	2.67	0.42
1:A:131:ASP:OD1	1:A:284:GLU:OE2	2.38	0.42
1:E:208:PHE:HD2	1:E:250:VAL:HG21	1.84	0.42
1:C:176:MSE:HE2	1:C:176:MSE:HB3	1.93	0.42
1:F:277:HIS:O	1:F:278:ALA:C	2.63	0.42
1:A:242:MSE:HE3	1:A:244:ASP:HB3	2.01	0.42
1:D:148:VAL:CG1	1:D:149:ALA:N	2.80	0.42
1:E:195:LEU:HD23	1:E:195:LEU:HA	1.80	0.42
1:D:136:LYS:CB	1:D:141:MSE:HE3	2.49	0.41
1:E:281:ILE:O	1:E:284:GLU:HB3	2.19	0.41
1:A:49:THR:OG1	1:A:288:ARG:NH1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:176:MSE:HE2	1:E:176:MSE:HB3	1.90	0.41
1:A:4:PHE:CD1	1:A:295:LEU:CD1	3.02	0.41
1:A:82:ASP:OD2	1:B:265:ARG:NH2	2.52	0.41
1:B:119:LEU:HD12	1:B:124:HIS:CD2	2.55	0.41
1:D:151:ILE:HD12	1:D:151:ILE:N	2.35	0.41
1:E:129:LEU:CB	1:E:291:VAL:HG11	2.50	0.41
1:E:151:ILE:HD12	1:E:151:ILE:N	2.35	0.41
1:A:51:MSE:HB2	1:A:51:MSE:HE2	1.81	0.41
1:B:73:GLN:HG3	1:B:73:GLN:O	2.20	0.41
1:C:195:LEU:HD23	1:C:195:LEU:HA	1.81	0.41
1:D:53:PHE:CD1	1:D:95:MSE:HE2	2.52	0.41
1:E:223:VAL:O	1:E:223:VAL:HG12	2.21	0.41
1:B:49:THR:HG22	1:B:95:MSE:HE1	2.02	0.41
1:D:194:ALA:CB	1:D:199:ALA:HB3	2.50	0.41
1:E:175:ALA:HB2	1:E:210:ALA:HB2	2.02	0.41
1:D:118:ALA:O	1:D:119:LEU:CB	2.68	0.41
1:F:245:GLU:HA	1:F:273:ILE:HD13	2.03	0.41
1:C:281:ILE:O	1:C:284:GLU:HB3	2.20	0.41
1:E:265:ARG:NH2	1:F:82:ASP:OD1	2.53	0.41
1:A:53:PHE:CD1	1:A:292:VAL:HG11	2.56	0.41
1:C:49:THR:OG1	1:C:288:ARG:NH1	2.53	0.41
1:D:74:LEU:HD11	1:D:96:MSE:CE	2.51	0.41
1:F:223:VAL:O	1:F:223:VAL:HG12	2.20	0.41
1:F:256:ILE:HD13	1:F:256:ILE:HG21	1.78	0.41
1:C:145:ILE:HG22	1:C:146:ALA:N	2.34	0.41
1:E:277:HIS:O	1:E:278:ALA:C	2.64	0.41
1:F:53:PHE:HB2	1:F:95:MSE:HE3	2.03	0.41
1:C:51:MSE:HE2	1:C:51:MSE:HB2	1.66	0.40
1:F:281:ILE:O	1:F:284:GLU:HB3	2.21	0.40
1:A:146:ALA:HB3	1:A:170:PHE:CD2	2.57	0.40
1:F:49:THR:OG1	1:F:288:ARG:NH1	2.54	0.40
1:F:60:LEU:HD23	1:F:60:LEU:HA	1.95	0.40
1:B:281:ILE:O	1:B:284:GLU:HB3	2.20	0.40
1:D:16:LEU:HD23	1:D:16:LEU:HA	1.93	0.40
1:E:297:CYS:O	1:E:301:ASN:ND2	2.50	0.40
1:A:213:ASP:CG	1:A:213:ASP:O	2.65	0.40
1:D:265:ARG:NH2	1:E:82:ASP:OD2	2.54	0.40
1:E:51:MSE:HE2	1:E:51:MSE:HB2	1.78	0.40
1:E:145:ILE:CG2	1:E:146:ALA:N	2.83	0.40
1:F:101:HIS:O	1:F:102:GLU:C	2.63	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	301/307 (98%)	286 (95%)	14 (5%)	1 (0%)	36	60
1	B	303/307 (99%)	285 (94%)	17 (6%)	1 (0%)	36	60
1	C	297/307 (97%)	284 (96%)	12 (4%)	1 (0%)	36	60
1	D	301/307 (98%)	286 (95%)	14 (5%)	1 (0%)	36	60
1	E	296/307 (96%)	286 (97%)	9 (3%)	1 (0%)	36	60
1	F	292/307 (95%)	279 (96%)	12 (4%)	1 (0%)	36	60
All	All	1790/1842 (97%)	1706 (95%)	78 (4%)	6 (0%)	36	60

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	LEU
1	B	119	LEU
1	C	119	LEU
1	D	119	LEU
1	E	119	LEU
1	F	119	LEU

5.3.2 Protein sidechains [i](#)

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	236/253 (93%)	234 (99%)	2 (1%)	73	88
1	B	238/253 (94%)	233 (98%)	5 (2%)	47	75

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	228/253 (90%)	224 (98%)	4 (2%)	51	78
1	D	228/253 (90%)	224 (98%)	4 (2%)	51	78
1	E	226/253 (89%)	223 (99%)	3 (1%)	61	83
1	F	230/253 (91%)	227 (99%)	3 (1%)	61	83
All	All	1386/1518 (91%)	1365 (98%)	21 (2%)	57	81

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

Mol	Chain	Res	Type
1	A	2	LYS
1	A	288	ARG
1	B	2	LYS
1	B	51	MSE
1	B	239	GLU
1	B	288	ARG
1	B	291	VAL
1	C	1	MSE
1	C	223	VAL
1	C	288	ARG
1	C	291	VAL
1	D	49	THR
1	D	203	LEU
1	D	288	ARG
1	D	291	VAL
1	E	1	MSE
1	E	288	ARG
1	E	291	VAL
1	F	51	MSE
1	F	288	ARG
1	F	291	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	HIS
1	A	45	ASN
1	A	73	GLN
1	A	99	ASN
1	A	301	ASN
1	C	21	HIS

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Mol	Chain	Res	Type
1	C	45	ASN
1	C	127	GLN
1	C	159	ASN
1	C	277	HIS
1	D	21	HIS
1	D	45	ASN
1	D	99	ASN
1	D	179	ASN
1	D	301	ASN
1	E	20	ASN
1	E	21	HIS
1	E	45	ASN
1	E	73	GLN
1	E	127	GLN
1	E	159	ASN
1	F	21	HIS
1	F	73	GLN
1	F	301	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PEG	F	307	-	6,6,6	0.75	0	5,5,5	0.51	0
2	PEG	A	307	-	6,6,6	0.68	0	5,5,5	0.25	0
2	PEG	B	307	-	6,6,6	0.69	0	5,5,5	0.38	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PEG	F	307	-	-	2/4/4/4	-
2	PEG	A	307	-	-	2/4/4/4	-
2	PEG	B	307	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	307	PEG	O1-C1-C2-O2
2	B	307	PEG	O1-C1-C2-O2
2	A	307	PEG	O2-C3-C4-O4
2	B	307	PEG	C4-C3-O2-C2
2	A	307	PEG	C4-C3-O2-C2
2	B	307	PEG	O2-C3-C4-O4
2	F	307	PEG	C4-C3-O2-C2

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

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	290/307 (94%)	-0.37	0 100 100	35, 68, 128, 178	0
1	B	292/307 (95%)	-0.42	0 100 100	30, 56, 110, 157	0
1	C	288/307 (93%)	-0.12	3 (1%) 79 78	41, 83, 171, 212	0
1	D	290/307 (94%)	-0.19	2 (0%) 84 83	34, 84, 149, 170	0
1	E	287/307 (93%)	0.00	2 (0%) 84 83	43, 102, 172, 187	0
1	F	286/307 (93%)	-0.36	3 (1%) 79 78	36, 62, 133, 161	0
All	All	1733/1842 (94%)	-0.24	10 (0%) 85 85	30, 73, 151, 212	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	151	ILE	3.2
1	C	256	ILE	2.8
1	C	223	VAL	2.5
1	F	224	SER	2.2
1	D	170	PHE	2.2
1	F	223	VAL	2.2
1	E	178	LYS	2.1
1	D	223	VAL	2.0
1	E	150	PHE	2.0
1	F	302	GLN	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
2	PEG	F	307	7/7	0.79	0.26	73,79,83,83	0
2	PEG	B	307	7/7	0.80	0.18	69,75,86,95	0
2	PEG	A	307	7/7	0.80	0.09	72,84,96,98	0

6.5 Other polymers [i](#)

There are no such residues in this entry.