



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 11:26 am BST

PDB ID : 3N50  
Title : Human Early B-cell factor 3 (EBF3) IPT/TIG and HLHLH domains  
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Deposited on : 2010-05-24  
Resolution : 3.10 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

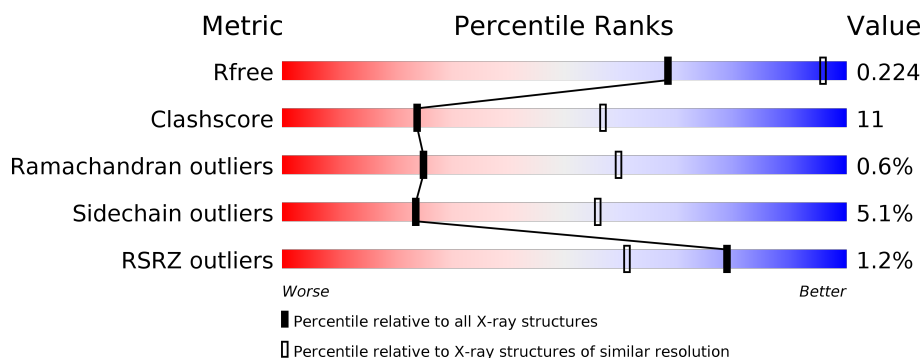
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.10 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))
$R_{free}$	130704	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	159	<div> <div>%</div> <div> <div></div> <div>55%</div> <div>30%</div> <div>16%</div> </div> </div>
1	B	159	<div> <div>%</div> <div> <div></div> <div>79%</div> <div>16%</div> <div>• •</div> </div> </div>
1	C	159	<div> <div>4%</div> <div> <div></div> <div>62%</div> <div>28%</div> <div>• 8%</div> </div> </div>
1	D	159	<div> <div></div> <div> <div></div> <div>72%</div> <div>24%</div> <div>• •</div> </div> </div>
1	E	159	<div> <div>%</div> <div> <div></div> <div>60%</div> <div>31%</div> <div>• 8%</div> </div> </div>
1	F	159	<div> <div>%</div> <div> <div></div> <div>69%</div> <div>25%</div> <div>• •</div> </div> </div>

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 6892 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 Transcription factor COE3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	134	Total	C	N	O	S	0	0	0
			1038	673	177	184	4			
1	B	154	Total	C	N	O	S	0	0	0
			1199	774	206	215	4			
1	C	146	Total	C	N	O	S	0	0	0
			1127	729	192	202	4			
1	D	154	Total	C	N	O	S	0	0	0
			1199	774	206	215	4			
1	E	146	Total	C	N	O	S	0	0	0
			1130	733	193	200	4			
1	F	154	Total	C	N	O	S	0	0	0
			1199	774	206	215	4			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	248	SER	-	EXPRESSION TAG	UNP Q9H4W6
A	249	MET	-	EXPRESSION TAG	UNP Q9H4W6
A	250	SER	-	EXPRESSION TAG	UNP Q9H4W6
A	251	GLU	-	EXPRESSION TAG	UNP Q9H4W6
B	248	SER	-	EXPRESSION TAG	UNP Q9H4W6
B	249	MET	-	EXPRESSION TAG	UNP Q9H4W6
B	250	SER	-	EXPRESSION TAG	UNP Q9H4W6
B	251	GLU	-	EXPRESSION TAG	UNP Q9H4W6
C	248	SER	-	EXPRESSION TAG	UNP Q9H4W6
C	249	MET	-	EXPRESSION TAG	UNP Q9H4W6
C	250	SER	-	EXPRESSION TAG	UNP Q9H4W6
C	251	GLU	-	EXPRESSION TAG	UNP Q9H4W6
D	248	SER	-	EXPRESSION TAG	UNP Q9H4W6
D	249	MET	-	EXPRESSION TAG	UNP Q9H4W6
D	250	SER	-	EXPRESSION TAG	UNP Q9H4W6
D	251	GLU	-	EXPRESSION TAG	UNP Q9H4W6
E	248	SER	-	EXPRESSION TAG	UNP Q9H4W6

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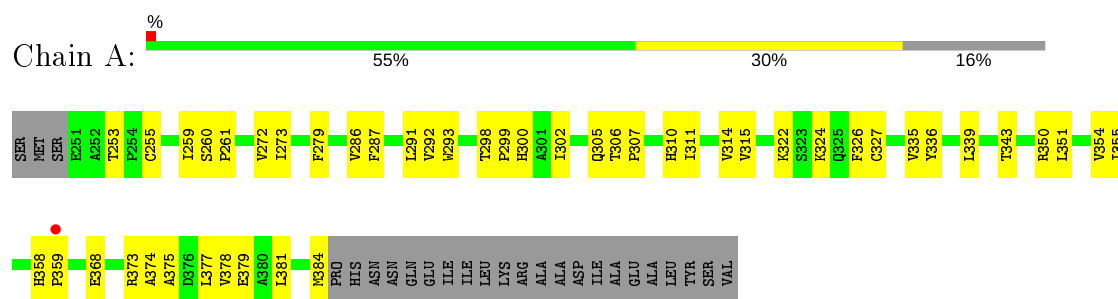
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Chain	Residue	Modelled	Actual	Comment	Reference
E	249	MET	-	EXPRESSION TAG	UNP Q9H4W6
E	250	SER	-	EXPRESSION TAG	UNP Q9H4W6
E	251	GLU	-	EXPRESSION TAG	UNP Q9H4W6
F	248	SER	-	EXPRESSION TAG	UNP Q9H4W6
F	249	MET	-	EXPRESSION TAG	UNP Q9H4W6
F	250	SER	-	EXPRESSION TAG	UNP Q9H4W6
F	251	GLU	-	EXPRESSION TAG	UNP Q9H4W6

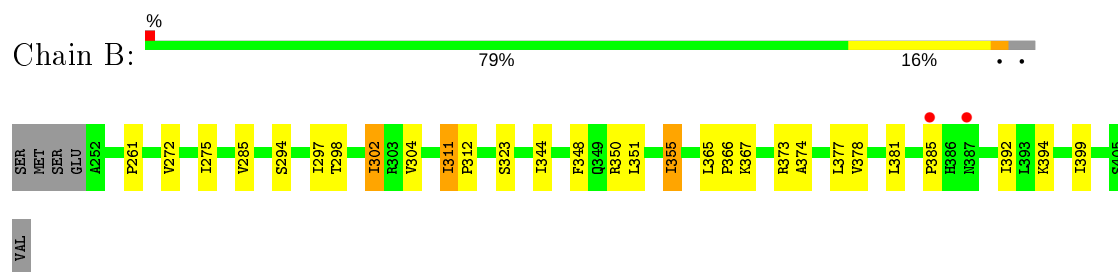
### 3 Residue-property plots [i](#)

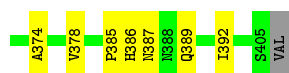
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Transcription factor COE3

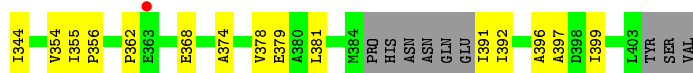
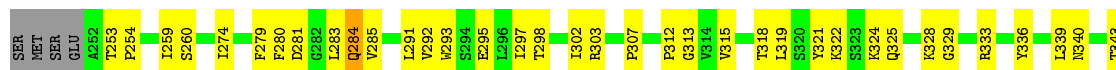


#### • Molecule 1: Transcription factor COE3





• Molecule 1: Transcription factor COE3



• Molecule 1: Transcription factor COE3



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	229.64Å 229.64Å 229.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.84 – 3.10 19.84 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.5 (19.84-3.10) 99.5 (19.84-3.10)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.73 (at 3.09Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, $R_{free}$	0.188 , 0.227 0.184 , 0.224	Depositor DCC
$R_{free}$ test set	2001 reflections (5.53%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	82.3	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 43.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.018 for -l,-k,-h	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6892	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.82% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.30	0/1066	0.49	0/1453
1	B	0.30	0/1230	0.48	0/1676
1	C	0.32	0/1154	0.50	0/1571
1	D	0.31	0/1230	0.49	0/1676
1	E	0.30	0/1157	0.48	0/1574
1	F	0.30	0/1230	0.47	0/1676
All	All	0.30	0/7067	0.49	0/9626

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1038	0	1055	35	0
1	B	1199	0	1218	22	0
1	C	1127	0	1147	39	0
1	D	1199	0	1218	23	0
1	E	1130	0	1163	36	0
1	F	1199	0	1218	31	0
All	All	6892	0	7019	154	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 (154) 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:E:292:VAL:HG12	1:E:307:PRO:HG3	1.62	0.81
1:C:307:PRO:HD2	1:C:336:TYR:OH	1.88	0.72
1:E:297:ILE:HG22	1:E:298:THR:HG23	1.72	0.72
1:E:399:ILE:HD12	1:F:356:PRO:HG3	1.71	0.71
1:D:386:HIS:HA	1:D:389:GLN:HB2	1.72	0.71
1:F:297:ILE:HG22	1:F:298:THR:HG23	1.76	0.68
1:C:392:ILE:HG23	1:C:393:LEU:H	1.59	0.66
1:A:307:PRO:HD2	1:A:336:TYR:OH	1.97	0.64
1:D:351:LEU:HD22	1:D:355:ILE:HD13	1.78	0.64
1:C:399:ILE:HD12	1:D:356:PRO:HG3	1.79	0.64
1:E:354:VAL:HG12	1:F:378:VAL:HG12	1.79	0.64
1:E:280:PHE:CE1	1:E:283:LEU:HB2	2.33	0.64
1:C:349:GLN:OE1	1:E:333:ARG:HD2	1.99	0.63
1:A:255:CYS:HA	1:A:327:CYS:SG	2.40	0.62
1:E:368:GLU:HG3	1:F:344:ILE:HG13	1.80	0.62
1:D:261:PRO:HD2	1:D:272:VAL:HG22	1.82	0.62
1:B:294:SER:HB2	1:B:302:ILE:HG23	1.83	0.61
1:A:314:VAL:HG22	1:A:335:VAL:HG22	1.83	0.61
1:C:297:ILE:HG22	1:C:298:THR:HG23	1.83	0.60
1:E:379:GLU:OE2	1:F:350:ARG:NH2	2.35	0.60
1:C:381:LEU:HD23	1:D:392:ILE:HD13	1.84	0.60
1:A:261:PRO:HD2	1:A:272:VAL:HG22	1.83	0.59
1:A:377:LEU:O	1:A:381:LEU:HG	2.03	0.59
1:C:354:VAL:HG12	1:D:378:VAL:HG12	1.84	0.59
1:C:255:CYS:HA	1:C:327:CYS:SG	2.42	0.59
1:F:351:LEU:HD22	1:F:355:ILE:HD13	1.85	0.58
1:E:279:PHE:CD2	1:E:302:ILE:HG13	2.38	0.58
1:E:307:PRO:HD2	1:E:336:TYR:OH	2.04	0.58
1:F:386:HIS:HA	1:F:389:GLN:HB2	1.84	0.58
1:A:259:ILE:HG22	1:A:261:PRO:O	2.04	0.57
1:C:358:HIS:HB3	1:C:359:PRO:HD2	1.86	0.57
1:A:354:VAL:HG12	1:B:378:VAL:HG12	1.86	0.56
1:A:351:LEU:O	1:A:355:ILE:HG12	2.04	0.56
1:B:351:LEU:HD22	1:B:355:ILE:HD13	1.88	0.56
1:A:381:LEU:HD21	1:B:381:LEU:HD12	1.86	0.56
1:A:374:ALA:O	1:A:378:VAL:HG23	2.05	0.55
1:A:379:GLU:OE2	1:B:399:ILE:HD13	2.06	0.55
1:A:381:LEU:CD2	1:B:392:ILE:HD13	2.36	0.55
1:C:396:ALA:O	1:C:397:ALA:HB3	2.08	0.53
1:F:261:PRO:HD2	1:F:272:VAL:HG22	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:279:PHE:CD2	1:C:302:ILE:HG13	2.43	0.53
1:C:312:PRO:HG3	1:C:338:ALA:HB2	1.90	0.52
1:E:392:ILE:HG23	1:E:392:ILE:O	2.08	0.52
1:C:261:PRO:HD2	1:C:272:VAL:HG22	1.91	0.52
1:A:339:LEU:HD11	1:B:344:ILE:HD13	1.90	0.52
1:C:368:GLU:HG3	1:D:344:ILE:HG13	1.92	0.52
1:C:379:GLU:OE2	1:D:350:ARG:NH2	2.42	0.52
1:C:292:VAL:HG12	1:C:307:PRO:HG3	1.92	0.52
1:E:344:ILE:HG13	1:F:368:GLU:HG2	1.91	0.52
1:F:314:VAL:HG22	1:F:335:VAL:HG22	1.92	0.52
1:E:391:ILE:HG22	1:E:391:ILE:O	2.09	0.52
1:A:273:ILE:HG21	1:B:275:ILE:HD11	1.91	0.52
1:D:348:PHE:CE1	1:D:367:LYS:HB3	2.45	0.52
1:E:280:PHE:HD1	1:E:281:ASP:O	1.93	0.51
1:F:366:PRO:HD2	1:F:369:VAL:HG21	1.92	0.51
1:C:286:VAL:HG22	1:C:291:LEU:HD23	1.92	0.51
1:E:381:LEU:HD23	1:F:392:ILE:HD13	1.92	0.51
1:F:329:GLY:C	1:F:331:PRO:HD3	2.31	0.51
1:C:381:LEU:CD2	1:D:392:ILE:HD13	2.40	0.50
1:C:392:ILE:HG23	1:C:393:LEU:N	2.26	0.50
1:A:350:ARG:O	1:A:354:VAL:HG23	2.11	0.50
1:A:292:VAL:HG12	1:A:307:PRO:HG3	1.93	0.50
1:F:294:SER:HB2	1:F:302:ILE:HG23	1.93	0.50
1:E:396:ALA:O	1:E:397:ALA:HB3	2.12	0.50
1:E:399:ILE:HD12	1:F:356:PRO:CG	2.42	0.49
1:C:358:HIS:HB3	1:C:359:PRO:CD	2.42	0.49
1:A:368:GLU:HG3	1:B:344:ILE:HG13	1.94	0.49
1:A:381:LEU:HD23	1:B:392:ILE:HD13	1.95	0.49
1:C:259:ILE:HG22	1:C:261:PRO:O	2.12	0.49
1:C:326:PHE:O	1:C:327:CYS:HB2	2.12	0.48
1:B:373:ARG:O	1:B:377:LEU:HG	2.13	0.48
1:E:328:LYS:HA	1:E:328:LYS:HD3	1.68	0.48
1:F:259:ILE:HG22	1:F:261:PRO:O	2.13	0.48
1:E:374:ALA:O	1:E:378:VAL:HG23	2.14	0.48
1:E:293:TRP:CZ2	1:E:295:GLU:HG3	2.50	0.47
1:C:251:GLU:HG2	1:C:252:ALA:H	1.79	0.47
1:B:297:ILE:HG22	1:B:298:THR:HG23	1.95	0.47
1:E:283:LEU:HD11	1:E:319:LEU:HB3	1.96	0.47
1:E:259:ILE:HD12	1:E:274:ILE:HG12	1.97	0.47
1:D:374:ALA:O	1:D:378:VAL:HG23	2.15	0.47
1:D:366:PRO:HD2	1:D:369:VAL:HG21	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:309:ARG:NH1	1:C:311:ILE:O	2.49	0.46
1:A:384:MET:HG2	1:B:392:ILE:HD11	1.97	0.46
1:E:285:VAL:O	1:E:292:VAL:HG22	2.15	0.46
1:A:286:VAL:HG22	1:A:291:LEU:HD23	1.97	0.46
1:C:251:GLU:CG	1:C:252:ALA:H	2.28	0.46
1:A:293:TRP:HE1	1:C:284:GLN:HG2	1.81	0.46
1:A:311:ILE:HD12	1:A:311:ILE:N	2.31	0.46
1:A:326:PHE:O	1:A:327:CYS:HB2	2.15	0.46
1:B:261:PRO:HD2	1:B:272:VAL:HG22	1.97	0.46
1:A:298:THR:C	1:A:300:HIS:H	2.19	0.46
1:E:391:ILE:O	1:E:392:ILE:HG22	2.16	0.46
1:F:375:ALA:O	1:F:379:GLU:HG3	2.16	0.46
1:F:385:PRO:C	1:F:387:ASN:H	2.19	0.45
1:C:314:VAL:HG22	1:C:335:VAL:HG22	1.97	0.45
1:F:385:PRO:O	1:F:386:HIS:HB3	2.16	0.45
1:D:329:GLY:C	1:D:331:PRO:HD3	2.36	0.45
1:D:297:ILE:HG22	1:D:298:THR:HG23	1.99	0.45
1:D:365:LEU:HA	1:D:366:PRO:HD3	1.82	0.45
1:F:283:LEU:HD23	1:F:302:ILE:CD1	2.47	0.45
1:C:340:ASN:O	1:C:341:GLU:C	2.55	0.44
1:E:292:VAL:CG1	1:E:307:PRO:HG3	2.39	0.44
1:A:305:GLN:NE2	1:C:291:LEU:H	2.15	0.44
1:A:305:GLN:HE22	1:C:291:LEU:H	1.65	0.44
1:F:266:THR:HB	1:F:309:ARG:O	2.17	0.44
1:D:311:ILE:HA	1:D:312:PRO:HD3	1.80	0.44
1:B:285:VAL:HG21	1:B:304:VAL:HG11	2.00	0.44
1:F:297:ILE:HG22	1:F:298:THR:CG2	2.47	0.44
1:F:352:GLN:NE2	1:F:370:LEU:HD11	2.33	0.44
1:A:358:HIS:HB3	1:A:359:PRO:HD2	2.00	0.44
1:C:311:ILE:HG13	1:E:312:PRO:HG2	1.99	0.43
1:D:385:PRO:C	1:D:387:ASN:H	2.21	0.43
1:C:340:ASN:HB3	1:E:313:GLY:HA3	2.00	0.43
1:B:351:LEU:HD22	1:B:355:ILE:CD1	2.49	0.43
1:A:291:LEU:HD13	1:C:293:TRP:CE2	2.54	0.43
1:F:378:VAL:O	1:F:381:LEU:HB3	2.19	0.43
1:A:306:THR:HA	1:A:307:PRO:HD3	1.75	0.43
1:C:307:PRO:HD2	1:C:336:TYR:HH	1.81	0.42
1:E:253:THR:HA	1:E:254:PRO:HD3	1.86	0.42
1:E:333:ARG:HG3	1:E:333:ARG:HH11	1.83	0.42
1:A:358:HIS:HD2	1:A:373:ARG:NH2	2.17	0.42
1:F:382:TYR:C	1:F:384:MET:H	2.21	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:355:ILE:HA	1:E:356:PRO:HD3	1.94	0.42
1:A:279:PHE:CD2	1:A:302:ILE:HG13	2.54	0.42
1:B:348:PHE:CE1	1:B:367:LYS:HB3	2.55	0.42
1:D:266:THR:HG23	1:D:337:THR:O	2.20	0.42
1:D:368:GLU:H	1:D:368:GLU:HG3	1.59	0.42
1:E:321:TYR:CZ	1:E:322:LYS:HE3	2.54	0.42
1:D:254:PRO:HA	1:D:278:ASN:HB2	2.01	0.42
1:F:366:PRO:HB2	1:F:369:VAL:HG23	2.01	0.41
1:C:363:GLU:HG2	1:E:329:GLY:HA2	2.02	0.41
1:A:287:PHE:CD1	1:A:292:VAL:HG11	2.55	0.41
1:F:298:THR:C	1:F:300:HIS:H	2.23	0.41
1:E:284:GLN:CG	1:E:291:LEU:HD22	2.50	0.41
1:E:322:LYS:O	1:E:324:LYS:N	2.45	0.41
1:E:339:LEU:HD11	1:F:344:ILE:HD13	2.02	0.41
1:B:374:ALA:O	1:B:378:VAL:HG23	2.20	0.41
1:A:322:LYS:C	1:A:324:LYS:H	2.23	0.41
1:A:375:ALA:CB	1:B:350:ARG:HG2	2.51	0.41
1:B:365:LEU:HA	1:B:366:PRO:HD3	1.92	0.41
1:C:368:GLU:H	1:C:368:GLU:CD	2.24	0.41
1:D:280:PHE:CZ	1:D:283:LEU:HD13	2.56	0.41
1:E:297:ILE:HD11	1:E:303:ARG:CZ	2.51	0.41
1:F:365:LEU:HA	1:F:366:PRO:HD3	1.93	0.41
1:F:395:ARG:O	1:F:399:ILE:HG13	2.21	0.41
1:C:298:THR:C	1:C:300:HIS:H	2.25	0.41
1:D:355:ILE:HA	1:D:356:PRO:HD3	1.82	0.41
1:A:310:HIS:HB2	1:A:311:ILE:HD12	2.02	0.40
1:B:367:LYS:H	1:B:367:LYS:HG2	1.69	0.40
1:F:355:ILE:HA	1:F:356:PRO:HD3	1.91	0.40
1:B:311:ILE:HA	1:B:312:PRO:HD3	1.81	0.40
1:C:374:ALA:O	1:C:378:VAL:HG23	2.21	0.40
1:C:346:TYR:O	1:C:349:GLN:HB2	2.21	0.40
1:D:255:CYS:O	1:D:276:GLY:HA3	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	132/159 (83%)	123 (93%)	8 (6%)	1 (1%)	19	54
1	B	152/159 (96%)	142 (93%)	9 (6%)	1 (1%)	22	57
1	C	142/159 (89%)	128 (90%)	14 (10%)	0	100	100
1	D	152/159 (96%)	144 (95%)	7 (5%)	1 (1%)	22	57
1	E	142/159 (89%)	126 (89%)	15 (11%)	1 (1%)	22	57
1	F	152/159 (96%)	144 (95%)	7 (5%)	1 (1%)	22	57
All	All	872/954 (91%)	807 (92%)	60 (7%)	5 (1%)	25	59

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	385	PRO
1	B	385	PRO
1	D	362	PRO
1	E	362	PRO
1	A	299	PRO

### 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	112/134 (84%)	108 (96%)	4 (4%)	35	67
1	B	129/134 (96%)	124 (96%)	5 (4%)	32	65
1	C	120/134 (90%)	113 (94%)	7 (6%)	20	51
1	D	129/134 (96%)	122 (95%)	7 (5%)	22	53
1	E	121/134 (90%)	114 (94%)	7 (6%)	20	51
1	F	129/134 (96%)	121 (94%)	8 (6%)	18	49
All	All	740/804 (92%)	702 (95%)	38 (5%)	24	56

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

Mol	Chain	Res	Type
1	A	253	THR
1	A	260	SER
1	A	315	VAL
1	A	343	THR
1	B	302	ILE
1	B	311	ILE
1	B	323	SER
1	B	355	ILE
1	B	394	LYS
1	C	251	GLU
1	C	260	SER
1	C	281	ASP
1	C	304	VAL
1	C	309	ARG
1	C	337	THR
1	C	393	LEU
1	D	260	SER
1	D	289	THR
1	D	311	ILE
1	D	323	SER
1	D	339	LEU
1	D	340	ASN
1	D	355	ILE
1	E	260	SER
1	E	284	GLN
1	E	315	VAL
1	E	318	THR
1	E	325	GLN
1	E	340	ASN
1	E	343	THR
1	F	260	SER
1	F	281	ASP
1	F	289	THR
1	F	298	THR
1	F	302	ILE
1	F	349	GLN
1	F	355	ILE
1	F	394	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	305	GLN
1	A	340	ASN
1	A	358	HIS
1	B	389	GLN
1	E	310	HIS
1	F	305	GLN
1	F	349	GLN

### 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 carbohydrates 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	134/159 (84%)	-0.21	1 (0%) 87 75	52, 81, 132, 144	0
1	B	154/159 (96%)	-0.36	2 (1%) 77 59	53, 73, 140, 162	0
1	C	146/159 (91%)	-0.23	6 (4%) 37 18	52, 72, 113, 143	0
1	D	154/159 (96%)	-0.45	0 100 100	52, 69, 119, 139	0
1	E	146/159 (91%)	-0.14	1 (0%) 87 75	52, 88, 119, 133	0
1	F	154/159 (96%)	-0.38	1 (0%) 89 78	52, 78, 132, 153	0
All	All	888/954 (93%)	-0.30	11 (1%) 79 61	52, 76, 126, 162	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	359	PRO	3.4
1	C	359	PRO	2.9
1	B	387	ASN	2.5
1	C	252	ALA	2.4
1	B	385	PRO	2.3
1	E	363	GLU	2.3
1	C	401	GLU	2.2
1	F	387	ASN	2.2
1	C	402	ALA	2.1
1	C	396	ALA	2.1
1	C	251	GLU	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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



### 6.3 Carbohydrates [i](#)

There are no carbohydrates 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.