



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 10:06 am GMT

PDB ID : 4B7W  
Title : Ligand binding domain human hepatocyte nuclear factor 4alpha: Apo form  
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Deposited on : 2012-08-24  
Resolution : 4.00 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

<http://wwpdb.org/validation/2016/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) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

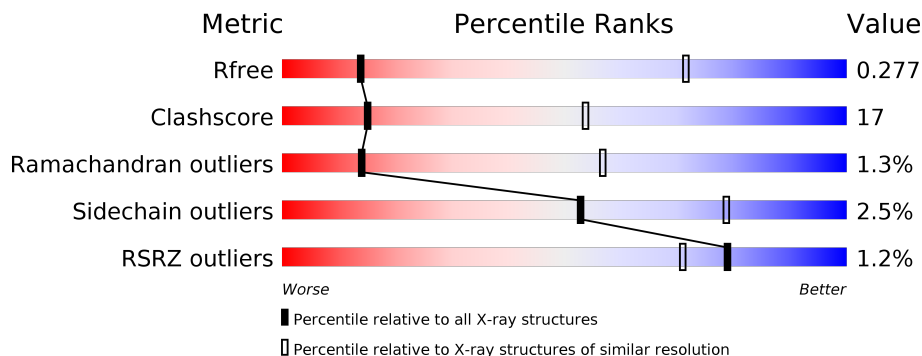
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 4.00 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))
$R_{free}$	100719	1088 (4.40-3.60)
Clashscore	112137	1187 (4.40-3.60)
Ramachandran outliers	110173	1139 (4.40-3.60)
Sidechain outliers	110143	1126 (4.40-3.60)
RSRZ outliers	101464	1099 (4.40-3.60)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	239	
1	B	239	
1	C	239	
1	D	239	

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 6476 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 HEPATOCYTE NUCLEAR FACTOR 4-ALPHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	202	Total	C	N	O	S	0	0	0
			1627	1050	271	297	9			
1	B	200	Total	C	N	O	S	0	0	0
			1611	1040	269	293	9			
1	C	202	Total	C	N	O	S	0	0	0
			1627	1050	271	297	9			
1	D	200	Total	C	N	O	S	0	0	0
			1611	1040	269	293	9			

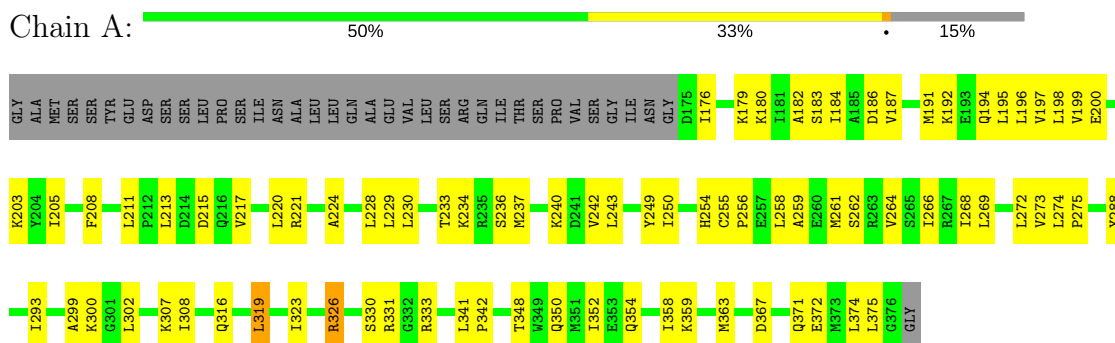
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	139	GLY	-	EXPRESSION TAG	UNP P41235
A	140	ALA	-	EXPRESSION TAG	UNP P41235
A	141	MET	-	EXPRESSION TAG	UNP P41235
B	139	GLY	-	EXPRESSION TAG	UNP P41235
B	140	ALA	-	EXPRESSION TAG	UNP P41235
B	141	MET	-	EXPRESSION TAG	UNP P41235
C	139	GLY	-	EXPRESSION TAG	UNP P41235
C	140	ALA	-	EXPRESSION TAG	UNP P41235
C	141	MET	-	EXPRESSION TAG	UNP P41235
D	139	GLY	-	EXPRESSION TAG	UNP P41235
D	140	ALA	-	EXPRESSION TAG	UNP P41235
D	141	MET	-	EXPRESSION TAG	UNP P41235

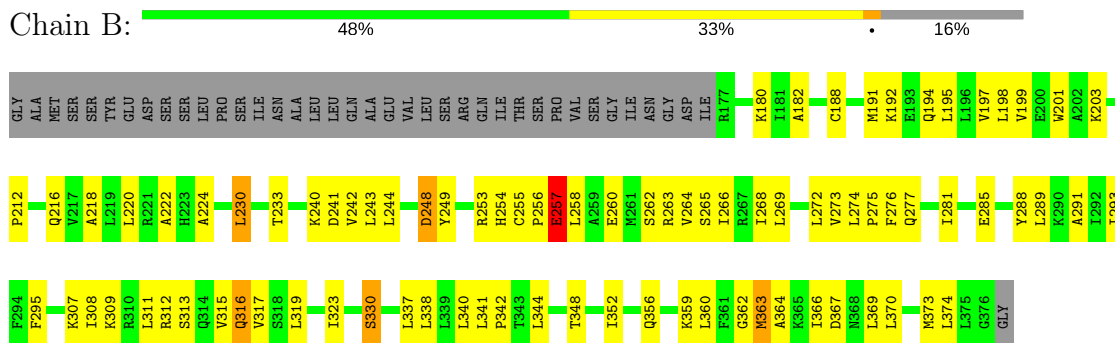
### 3 Residue-property plots

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.

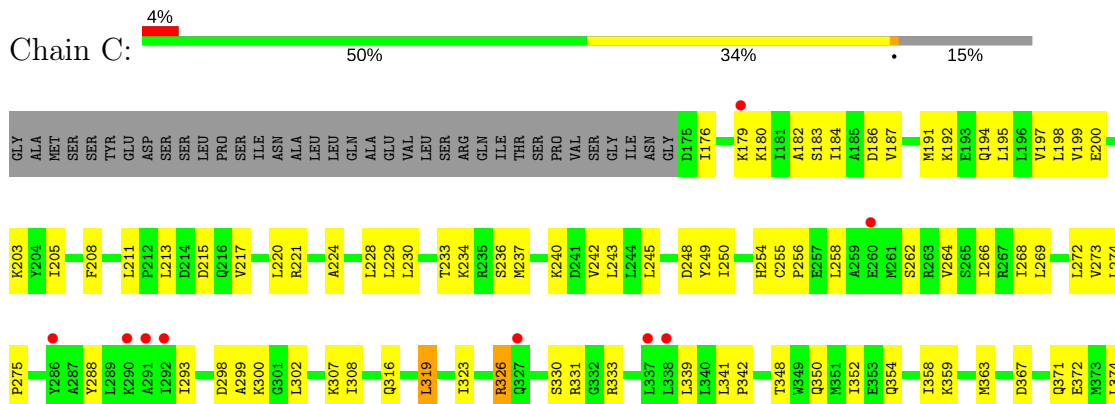
#### • Molecule 1: HEPATOCYTE NUCLEAR FACTOR 4-ALPHA



#### • Molecule 1: HEPATOCYTE NUCLEAR FACTOR 4-ALPHA



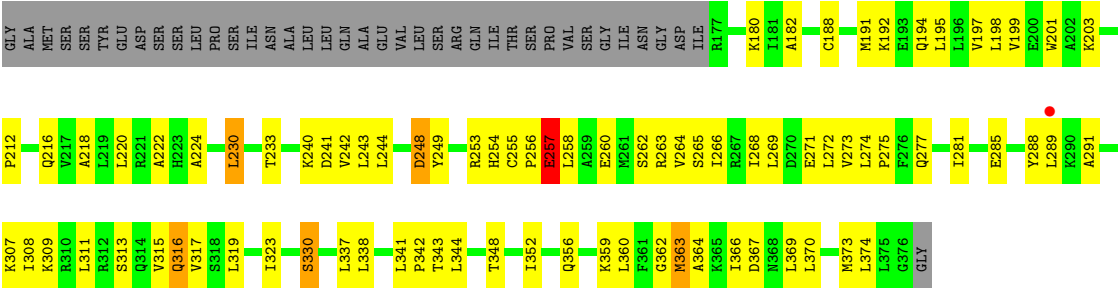
#### • Molecule 1: HEPATOCYTE NUCLEAR FACTOR 4-ALPHA



L375  
G376  
GLY

● Molecule 1: HEPATOCYTE NUCLEAR FACTOR 4-ALPHA

Chain D: 49% 32% 16%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	101.05Å 105.28Å 98.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 4.00 46.69 – 4.02	Depositor EDS
% Data completeness (in resolution range)	85.9 (50.00-4.00) 87.4 (46.69-4.02)	Depositor EDS
$R_{merge}$	0.16	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.05 (at 4.00Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.250 , 0.280 0.258 , 0.277	Depositor DCC
$R_{free}$ test set	374 reflections (4.68%)	DCC
Wilson B-factor (Å <sup>2</sup> )	129.2	Xtriage
Anisotropy	0.554	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 119.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.39$ , $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.079 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	6476	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	180.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.30% 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 [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	0.24	0/1655	0.42	0/2235
1	B	0.24	0/1639	0.42	0/2213
1	C	0.24	0/1655	0.42	0/2235
1	D	0.24	0/1639	0.42	0/2213
All	All	0.24	0/6588	0.42	0/8896

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 [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	1627	0	1675	55	0
1	B	1611	0	1660	61	0
1	C	1627	0	1675	56	0
1	D	1611	0	1660	58	0
All	All	6476	0	6670	221	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 17.

All (221) 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:358:ILE:HG23	1:A:363:MET:HB2	1.63	0.80
1:B:364:ALA:HA	1:B:367:ASP:HB3	1.62	0.80
1:D:364:ALA:HA	1:D:367:ASP:HB3	1.61	0.80
1:A:341:LEU:HB2	1:A:342:PRO:HD3	1.66	0.78
1:C:358:ILE:HG23	1:C:363:MET:HB2	1.64	0.78
1:C:341:LEU:HB2	1:C:342:PRO:HD3	1.66	0.76
1:B:260:GLU:HA	1:B:263:ARG:HH12	1.55	0.72
1:B:241:ASP:HB3	1:B:253:ARG:HB2	1.72	0.72
1:D:260:GLU:HA	1:D:263:ARG:HH12	1.55	0.71
1:D:241:ASP:HB3	1:D:253:ARG:HB2	1.73	0.70
1:D:274:LEU:HB3	1:D:275:PRO:HD3	1.73	0.70
1:B:274:LEU:HB3	1:B:275:PRO:HD3	1.73	0.69
1:A:184:ILE:HD12	1:A:184:ILE:H	1.57	0.69
1:C:184:ILE:HD12	1:C:184:ILE:H	1.59	0.68
1:B:230:LEU:HD23	1:B:344:LEU:HD22	1.75	0.67
1:C:182:ALA:HB2	1:C:249:TYR:HB3	1.75	0.67
1:A:182:ALA:HB2	1:A:249:TYR:HB3	1.77	0.67
1:D:230:LEU:HD23	1:D:344:LEU:HD22	1.75	0.66
1:D:341:LEU:HB2	1:D:342:PRO:HD3	1.78	0.66
1:A:274:LEU:HB3	1:A:275:PRO:HD3	1.78	0.66
1:B:341:LEU:HB2	1:B:342:PRO:HD3	1.77	0.65
1:C:274:LEU:HB3	1:C:275:PRO:HD3	1.78	0.65
1:D:269:LEU:HA	1:D:273:VAL:HB	1.79	0.65
1:B:182:ALA:HB2	1:B:249:TYR:HB3	1.78	0.64
1:D:263:ARG:HA	1:D:266:ILE:HD12	1.79	0.64
1:D:182:ALA:HB2	1:D:249:TYR:HB3	1.79	0.64
1:B:269:LEU:HA	1:B:273:VAL:HB	1.78	0.64
1:A:233:THR:HG21	1:A:272:LEU:HB3	1.79	0.63
1:C:233:THR:HG21	1:C:272:LEU:HB3	1.80	0.63
1:A:230:LEU:HD11	1:A:293:ILE:HD12	1.80	0.62
1:B:263:ARG:HA	1:B:266:ILE:HD12	1.82	0.62
1:B:269:LEU:HD23	1:B:273:VAL:HG21	1.81	0.62
1:D:269:LEU:HD23	1:D:273:VAL:HG21	1.81	0.61
1:D:313:SER:HA	1:D:316:GLN:HG2	1.83	0.61
1:C:230:LEU:HD11	1:C:293:ILE:HD12	1.81	0.60
1:B:244:LEU:HD11	1:B:248:ASP:HA	1.83	0.60
1:A:192:LYS:HE2	1:A:367:ASP:HB2	1.84	0.60
1:C:192:LYS:HE2	1:C:367:ASP:HB2	1.84	0.59
1:D:352:ILE:HG22	1:D:356:GLN:HE21	1.67	0.59
1:B:352:ILE:HG22	1:B:356:GLN:HE21	1.68	0.58
1:B:313:SER:HA	1:B:316:GLN:HG2	1.86	0.58
1:C:299:ALA:HB3	1:C:302:LEU:HG	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:299:ALA:HB3	1:A:302:LEU:HG	1.86	0.56
1:D:244:LEU:HD11	1:D:248:ASP:HA	1.86	0.56
1:C:323:ILE:HG21	1:C:333:ARG:HB3	1.88	0.55
1:A:323:ILE:HG21	1:A:333:ARG:HB3	1.89	0.55
1:A:220:LEU:O	1:A:224:ALA:HB2	2.07	0.54
1:A:217:VAL:HG12	1:A:221:ARG:HD3	1.90	0.54
1:B:289:LEU:HD11	1:B:337:LEU:HD22	1.89	0.54
1:D:289:LEU:HD11	1:D:337:LEU:HD22	1.89	0.54
1:C:220:LEU:O	1:C:224:ALA:HB2	2.07	0.53
1:B:220:LEU:O	1:B:224:ALA:HB2	2.08	0.53
1:D:220:LEU:O	1:D:224:ALA:HB2	2.08	0.53
1:B:233:THR:HG23	1:B:273:VAL:HA	1.91	0.53
1:B:360:LEU:O	1:D:360:LEU:HD13	2.08	0.53
1:D:233:THR:HG23	1:D:273:VAL:HA	1.91	0.53
1:C:268:ILE:HG22	1:C:273:VAL:HG23	1.90	0.53
1:A:194:GLN:HA	1:A:197:VAL:HG22	1.91	0.53
1:C:180:LYS:O	1:C:250:ILE:HG12	2.08	0.53
1:C:217:VAL:HG12	1:C:221:ARG:HD3	1.90	0.53
1:A:180:LYS:O	1:A:250:ILE:HG12	2.08	0.53
1:B:199:VAL:O	1:B:203:LYS:HG2	2.09	0.53
1:B:362:GLY:O	1:B:366:ILE:HG12	2.09	0.53
1:A:268:ILE:HG22	1:A:273:VAL:HG23	1.91	0.52
1:C:194:GLN:HA	1:C:197:VAL:HG22	1.91	0.52
1:B:323:ILE:HG22	1:B:330:SER:HB2	1.91	0.52
1:B:364:ALA:HB3	1:D:360:LEU:HD11	1.91	0.52
1:D:362:GLY:O	1:D:366:ILE:HG12	2.10	0.52
1:D:199:VAL:O	1:D:203:LYS:HG2	2.09	0.52
1:B:240:LYS:O	1:B:242:VAL:HG23	2.10	0.51
1:B:212:PRO:O	1:B:216:GLN:HG3	2.10	0.51
1:D:254:HIS:O	1:D:256:PRO:HD3	2.10	0.51
1:B:360:LEU:HD13	1:D:360:LEU:O	2.11	0.51
1:D:212:PRO:O	1:D:216:GLN:HG3	2.11	0.51
1:D:240:LYS:O	1:D:242:VAL:HG23	2.11	0.51
1:D:323:ILE:HG22	1:D:330:SER:HB2	1.92	0.51
1:A:234:LYS:O	1:A:237:MET:HG2	2.12	0.50
1:B:254:HIS:O	1:B:256:PRO:HD3	2.11	0.50
1:D:264:VAL:O	1:D:268:ILE:HG13	2.12	0.50
1:D:243:LEU:HG	1:D:269:LEU:HD21	1.93	0.50
1:B:272:LEU:C	1:B:275:PRO:HD2	2.32	0.50
1:D:359:LYS:O	1:D:363:MET:HG2	2.12	0.50
1:A:211:LEU:HD23	1:A:307:LYS:HE3	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:LEU:HA	1:A:273:VAL:HB	1.94	0.49
1:C:200:GLU:OE1	1:C:203:LYS:HD2	2.11	0.49
1:A:240:LYS:O	1:A:242:VAL:HG23	2.13	0.49
1:A:183:SER:H	1:A:186:ASP:HB2	1.77	0.49
1:A:350:GLN:O	1:A:354:GLN:HG2	2.13	0.49
1:B:264:VAL:O	1:B:268:ILE:HG13	2.12	0.49
1:B:367:ASP:HA	1:B:370:LEU:HD12	1.95	0.49
1:C:350:GLN:O	1:C:354:GLN:HG2	2.12	0.49
1:C:269:LEU:HA	1:C:273:VAL:HB	1.95	0.49
1:A:200:GLU:OE1	1:A:203:LYS:HD2	2.12	0.49
1:B:288:TYR:CD2	1:B:319:LEU:HD23	2.48	0.49
1:B:359:LYS:O	1:B:363:MET:HG2	2.13	0.49
1:C:359:LYS:NZ	1:C:375:LEU:HD13	2.27	0.49
1:D:367:ASP:HA	1:D:370:LEU:HD12	1.95	0.48
1:C:272:LEU:C	1:C:275:PRO:HD2	2.34	0.48
1:D:272:LEU:C	1:D:275:PRO:HD2	2.33	0.48
1:B:265:SER:HA	1:B:268:ILE:HD12	1.95	0.48
1:C:183:SER:H	1:C:186:ASP:HB2	1.78	0.48
1:C:211:LEU:HD23	1:C:307:LYS:HE3	1.94	0.48
1:C:234:LYS:O	1:C:237:MET:HG2	2.13	0.48
1:C:240:LYS:O	1:C:242:VAL:HG23	2.13	0.48
1:D:308:ILE:HA	1:D:311:LEU:HD12	1.94	0.48
1:B:243:LEU:HG	1:B:269:LEU:HD21	1.93	0.48
1:A:272:LEU:C	1:A:275:PRO:HD2	2.34	0.48
1:B:308:ILE:HA	1:B:311:LEU:HD12	1.95	0.48
1:D:288:TYR:CD2	1:D:319:LEU:HD23	2.48	0.48
1:D:319:LEU:O	1:D:323:ILE:HG12	2.13	0.48
1:D:366:ILE:O	1:D:370:LEU:HG	2.14	0.48
1:A:359:LYS:NZ	1:A:375:LEU:HD13	2.29	0.47
1:D:262:SER:O	1:D:266:ILE:HG13	2.14	0.47
1:B:268:ILE:HG22	1:B:273:VAL:HG23	1.95	0.47
1:A:323:ILE:HG22	1:A:330:SER:HB2	1.97	0.47
1:B:366:ILE:O	1:B:370:LEU:HG	2.14	0.47
1:C:323:ILE:HG22	1:C:330:SER:HB2	1.96	0.47
1:D:268:ILE:HG22	1:D:273:VAL:HG23	1.95	0.47
1:B:262:SER:O	1:B:266:ILE:HG13	2.14	0.47
1:A:236:SER:OG	1:A:243:LEU:HA	2.15	0.47
1:D:291:ALA:HB3	1:D:315:VAL:HG11	1.97	0.47
1:A:187:VAL:O	1:A:191:MET:HG3	2.15	0.47
1:B:194:GLN:HA	1:B:197:VAL:HG22	1.96	0.47
1:D:265:SER:HA	1:D:268:ILE:HD12	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:187:VAL:O	1:C:191:MET:HG3	2.15	0.47
1:C:236:SER:OG	1:C:243:LEU:HA	2.15	0.47
1:C:255:CYS:SG	1:C:258:LEU:HD12	2.55	0.47
1:A:348:THR:O	1:A:352:ILE:HG13	2.16	0.46
1:B:255:CYS:SG	1:B:257:GLU:HG2	2.56	0.46
1:B:256:PRO:C	1:B:258:LEU:H	2.19	0.46
1:B:291:ALA:HB3	1:B:315:VAL:HG11	1.96	0.46
1:B:319:LEU:O	1:B:323:ILE:HG12	2.15	0.46
1:A:308:ILE:H	1:A:308:ILE:HD12	1.79	0.46
1:D:194:GLN:HA	1:D:197:VAL:HG22	1.98	0.46
1:D:255:CYS:SG	1:D:257:GLU:HG2	2.56	0.46
1:B:268:ILE:O	1:B:273:VAL:HG23	2.16	0.46
1:C:195:LEU:HD21	1:C:374:LEU:HD11	1.98	0.46
1:D:256:PRO:C	1:D:258:LEU:H	2.19	0.46
1:C:331:ARG:NH1	1:D:317:VAL:HG11	2.30	0.45
1:C:348:THR:O	1:C:352:ILE:HG13	2.16	0.45
1:A:255:CYS:SG	1:A:258:LEU:HD12	2.56	0.45
1:A:221:ARG:HG2	1:A:372:GLU:O	2.16	0.45
1:C:213:LEU:O	1:C:217:VAL:HG23	2.17	0.45
1:A:299:ALA:HB3	1:A:302:LEU:CG	2.47	0.45
1:A:359:LYS:NZ	1:A:371:GLN:HE22	2.15	0.45
1:C:359:LYS:NZ	1:C:371:GLN:HE22	2.15	0.45
1:C:221:ARG:HG2	1:C:372:GLU:O	2.17	0.45
1:A:195:LEU:HD21	1:A:374:LEU:HD11	1.98	0.44
1:C:308:ILE:H	1:C:308:ILE:HD12	1.81	0.44
1:C:300:LYS:NZ	1:C:300:LYS:HB2	2.32	0.44
1:C:354:GLN:O	1:C:358:ILE:HG13	2.17	0.44
1:C:229:LEU:O	1:C:233:THR:HG23	2.17	0.44
1:A:213:LEU:O	1:A:217:VAL:HG23	2.17	0.44
1:C:195:LEU:HD21	1:C:228:LEU:HD11	1.98	0.44
1:A:354:GLN:O	1:A:358:ILE:HG13	2.17	0.44
1:B:348:THR:O	1:B:352:ILE:HG13	2.17	0.44
1:D:188:CYS:O	1:D:192:LYS:HG3	2.17	0.44
1:B:191:MET:O	1:B:195:LEU:HB2	2.17	0.44
1:D:268:ILE:O	1:D:273:VAL:HG23	2.17	0.44
1:D:370:LEU:O	1:D:374:LEU:HB2	2.18	0.44
1:C:299:ALA:HB3	1:C:302:LEU:CG	2.47	0.44
1:C:319:LEU:HD22	1:C:323:ILE:HD11	2.00	0.44
1:D:198:LEU:O	1:D:201:TRP:HB3	2.18	0.44
1:C:176:ILE:HA	1:C:179:LYS:HD2	2.00	0.44
1:A:195:LEU:HD21	1:A:228:LEU:HD11	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:LYS:NZ	1:A:300:LYS:HB2	2.33	0.44
1:D:307:LYS:O	1:D:311:LEU:HG	2.18	0.44
1:A:176:ILE:HA	1:A:179:LYS:HD2	2.00	0.43
1:A:264:VAL:O	1:A:268:ILE:HG13	2.19	0.43
1:A:268:ILE:O	1:A:273:VAL:HG23	2.18	0.43
1:B:370:LEU:O	1:B:374:LEU:HB2	2.18	0.43
1:A:262:SER:O	1:A:266:ILE:HG13	2.18	0.43
1:A:199:VAL:O	1:A:203:LYS:HG3	2.18	0.43
1:B:188:CYS:O	1:B:192:LYS:HG3	2.19	0.43
1:B:198:LEU:O	1:B:201:TRP:HB3	2.19	0.43
1:B:307:LYS:O	1:B:311:LEU:HG	2.19	0.43
1:C:319:LEU:O	1:C:323:ILE:HG13	2.19	0.43
1:A:229:LEU:O	1:A:233:THR:HG23	2.19	0.43
1:C:199:VAL:O	1:C:203:LYS:HG3	2.18	0.43
1:D:191:MET:O	1:D:195:LEU:HB2	2.19	0.43
1:A:319:LEU:O	1:A:323:ILE:HG13	2.19	0.43
1:D:348:THR:O	1:D:352:ILE:HG13	2.19	0.43
1:A:255:CYS:SG	1:A:258:LEU:HB2	2.59	0.42
1:D:281:ILE:HG23	1:D:285:GLU:HB2	2.01	0.42
1:A:288:TYR:CD2	1:A:319:LEU:HG	2.54	0.42
1:B:364:ALA:HB2	1:D:360:LEU:HD21	2.01	0.42
1:C:262:SER:O	1:C:266:ILE:HG13	2.19	0.42
1:A:342:PRO:HG3	1:B:342:PRO:HG3	2.00	0.42
1:C:198:LEU:HD13	1:C:198:LEU:O	2.19	0.42
1:B:289:LEU:O	1:B:293:ILE:HG13	2.19	0.42
1:A:319:LEU:HD22	1:A:323:ILE:HD11	2.00	0.42
1:D:273:VAL:O	1:D:277:GLN:HG3	2.20	0.42
1:C:268:ILE:O	1:C:273:VAL:HG23	2.19	0.42
1:D:218:ALA:O	1:D:222:ALA:HB2	2.20	0.42
1:A:254:HIS:O	1:A:256:PRO:HD3	2.20	0.41
1:B:218:ALA:O	1:B:222:ALA:HB2	2.20	0.41
1:C:254:HIS:O	1:C:256:PRO:HD3	2.21	0.41
1:A:198:LEU:O	1:A:198:LEU:HD13	2.19	0.41
1:B:276:PHE:CD2	1:B:281:ILE:HD12	2.55	0.41
1:B:281:ILE:HG23	1:B:285:GLU:HB2	2.03	0.41
1:B:369:LEU:O	1:B:373:MET:HG2	2.20	0.41
1:C:205:ILE:HG22	1:C:208:PHE:H	1.84	0.41
1:D:369:LEU:O	1:D:373:MET:HG2	2.20	0.41
1:A:205:ILE:HG22	1:A:208:PHE:H	1.84	0.41
1:C:264:VAL:O	1:C:268:ILE:HG13	2.21	0.41
1:C:288:TYR:CD2	1:C:319:LEU:HG	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:LYS:O	1:A:196:LEU:HG	2.20	0.41
1:B:244:LEU:CD1	1:B:248:ASP:HA	2.51	0.41
1:B:338:LEU:C	1:B:340:LEU:H	2.24	0.41
1:C:255:CYS:SG	1:C:258:LEU:HB2	2.61	0.41
1:C:323:ILE:CG2	1:C:330:SER:HB2	2.51	0.41
1:A:331:ARG:NH1	1:B:317:VAL:HG11	2.35	0.40
1:C:339:LEU:HD21	1:D:338:LEU:HD22	2.02	0.40
1:D:352:ILE:HG22	1:D:356:GLN:NE2	2.35	0.40
1:B:198:LEU:O	1:B:198:LEU:HD23	2.21	0.40
1:C:298:ASP:OD1	1:D:271:GLU:HG2	2.21	0.40
1:A:259:ALA:C	1:A:261:MET:H	2.25	0.40
1:B:273:VAL:O	1:B:277:GLN:HG3	2.21	0.40
1:C:245:LEU:HB2	1:C:248:ASP:O	2.22	0.40
1:D:271:GLU:HB3	1:D:343:THR:HG21	2.04	0.40
1:B:295:PHE:O	1:B:312:ARG:HB2	2.21	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	200/239 (84%)	176 (88%)	23 (12%)	1 (0%)	32	73
1	B	198/239 (83%)	176 (89%)	18 (9%)	4 (2%)	9	49
1	C	200/239 (84%)	175 (88%)	24 (12%)	1 (0%)	32	73
1	D	198/239 (83%)	177 (89%)	17 (9%)	4 (2%)	9	49
All	All	796/956 (83%)	704 (88%)	82 (10%)	10 (1%)	14	57

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	330	SER
1	D	330	SER
1	B	363	MET
1	D	363	MET
1	B	180	LYS
1	D	180	LYS
1	A	326	ARG
1	B	257	GLU
1	C	326	ARG
1	D	257	GLU

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	179/209 (86%)	175 (98%)	4 (2%)	57	81
1	B	177/209 (85%)	172 (97%)	5 (3%)	49	76
1	C	179/209 (86%)	175 (98%)	4 (2%)	57	81
1	D	177/209 (85%)	172 (97%)	5 (3%)	49	76
All	All	712/836 (85%)	694 (98%)	18 (2%)	53	79

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	215	ASP
1	A	316	GLN
1	A	319	LEU
1	A	326	ARG
1	B	230	LEU
1	B	248	ASP
1	B	257	GLU
1	B	309	LYS
1	B	316	GLN
1	C	215	ASP
1	C	316	GLN
1	C	319	LEU

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Mol	Chain	Res	Type
1	C	326	ARG
1	D	230	LEU
1	D	248	ASP
1	D	257	GLU
1	D	309	LYS
1	D	316	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	277	GLN
1	A	284	ASN
1	A	324	ASN
1	A	350	GLN
1	A	356	GLN
1	B	227	HIS
1	B	277	GLN
1	B	354	GLN
1	B	356	GLN
1	C	277	GLN
1	C	284	ASN
1	C	324	ASN
1	C	350	GLN
1	C	356	GLN
1	C	371	GLN
1	D	227	HIS
1	D	277	GLN
1	D	345	GLN
1	D	354	GLN
1	D	356	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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 [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, 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	202/239 (84%)	-0.32	0	100 100	81, 158, 228, 285	0
1	B	200/239 (83%)	-0.45	0	100 100	53, 134, 205, 267	0
1	C	202/239 (84%)	0.12	9 (4%)	34 27	135, 204, 262, 306	0
1	D	200/239 (83%)	-0.13	1 (0%)	90 86	134, 210, 254, 286	0
All	All	804/956 (84%)	-0.19	10 (1%)	79 71	53, 184, 252, 306	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	179	LYS	3.9
1	C	290	LYS	3.7
1	C	291	ALA	3.0
1	C	327	GLN	2.6
1	C	286	TYR	2.5
1	C	338	LEU	2.4
1	C	292	ILE	2.3
1	C	260	GLU	2.3
1	C	337	LEU	2.1
1	D	289	LEU	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 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.