



# Full wwPDB X-ray Structure Validation Report ⓘ

May 8, 2016 – 11:33 PM EDT

PDB ID : 5E9F  
Title : Structural insights of isocitrate lyases from *Magnaporthe oryzae*  
Authors : Park, Y.; Cho, Y.; Lee, Y.-H.; Lee, Y.-W.; Rhee, S.  
Deposited on : 2015-10-15  
Resolution : 2.80 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  
Mogul : unknown  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20027457  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : rb-20027457

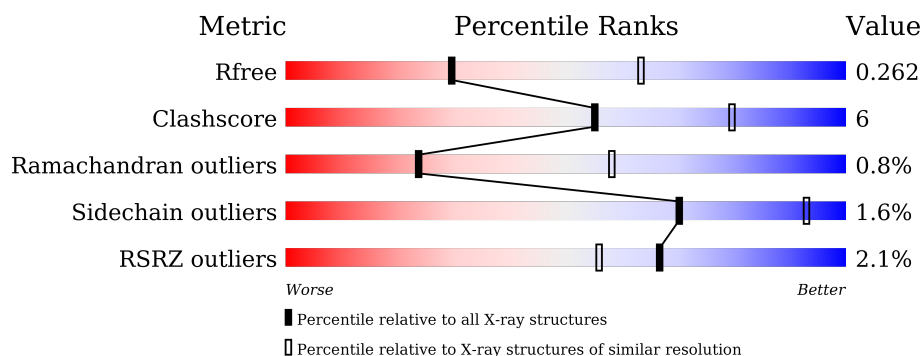
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
$R_{free}$	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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	560	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">%</span> </div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>78%</span> <span>12%</span> <span>• 9%</span> </div> </div>
1	B	560	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">%</span> </div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>77%</span> <span>11%</span> <span>• 11%</span> </div> </div>
1	C	560	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">2%</span> </div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>73%</span> <span>16%</span> <span>• 11%</span> </div> </div>
1	D	560	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">3%</span> </div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>65%</span> <span>15%</span> <span>• 19%</span> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MG	A	601	-	-	-	X
2	MG	B	601	-	-	-	X
2	MG	C	601	-	-	-	X
2	MG	D	601	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14983 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 Isocitrate lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	507	Total	C	N	O	S	0	0	0
			3851	2440	662	731	18			
1	B	501	Total	C	N	O	S	0	0	0
			3853	2444	662	729	18			
1	C	496	Total	C	N	O	S	0	0	0
			3736	2368	642	709	17			
1	D	453	Total	C	N	O	S	0	0	0
			3474	2200	597	662	15			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	548	ALA	-	expression tag	UNP P0CT06
A	549	ALA	-	expression tag	UNP P0CT06
A	550	ALA	-	expression tag	UNP P0CT06
A	551	LEU	-	expression tag	UNP P0CT06
A	552	GLU	-	expression tag	UNP P0CT06
A	553	HIS	-	expression tag	UNP P0CT06
A	554	HIS	-	expression tag	UNP P0CT06
A	555	HIS	-	expression tag	UNP P0CT06
A	556	HIS	-	expression tag	UNP P0CT06
A	557	HIS	-	expression tag	UNP P0CT06
A	558	HIS	-	expression tag	UNP P0CT06
A	559	HIS	-	expression tag	UNP P0CT06
A	560	HIS	-	expression tag	UNP P0CT06
B	548	ALA	-	expression tag	UNP P0CT06
B	549	ALA	-	expression tag	UNP P0CT06
B	550	ALA	-	expression tag	UNP P0CT06
B	551	LEU	-	expression tag	UNP P0CT06
B	552	GLU	-	expression tag	UNP P0CT06
B	553	HIS	-	expression tag	UNP P0CT06
B	554	HIS	-	expression tag	UNP P0CT06
B	555	HIS	-	expression tag	UNP P0CT06

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Chain	Residue	Modelled	Actual	Comment	Reference
B	556	HIS	-	expression tag	UNP P0CT06
B	557	HIS	-	expression tag	UNP P0CT06
B	558	HIS	-	expression tag	UNP P0CT06
B	559	HIS	-	expression tag	UNP P0CT06
B	560	HIS	-	expression tag	UNP P0CT06
C	548	ALA	-	expression tag	UNP P0CT06
C	549	ALA	-	expression tag	UNP P0CT06
C	550	ALA	-	expression tag	UNP P0CT06
C	551	LEU	-	expression tag	UNP P0CT06
C	552	GLU	-	expression tag	UNP P0CT06
C	553	HIS	-	expression tag	UNP P0CT06
C	554	HIS	-	expression tag	UNP P0CT06
C	555	HIS	-	expression tag	UNP P0CT06
C	556	HIS	-	expression tag	UNP P0CT06
C	557	HIS	-	expression tag	UNP P0CT06
C	558	HIS	-	expression tag	UNP P0CT06
C	559	HIS	-	expression tag	UNP P0CT06
C	560	HIS	-	expression tag	UNP P0CT06
D	548	ALA	-	expression tag	UNP P0CT06
D	549	ALA	-	expression tag	UNP P0CT06
D	550	ALA	-	expression tag	UNP P0CT06
D	551	LEU	-	expression tag	UNP P0CT06
D	552	GLU	-	expression tag	UNP P0CT06
D	553	HIS	-	expression tag	UNP P0CT06
D	554	HIS	-	expression tag	UNP P0CT06
D	555	HIS	-	expression tag	UNP P0CT06
D	556	HIS	-	expression tag	UNP P0CT06
D	557	HIS	-	expression tag	UNP P0CT06
D	558	HIS	-	expression tag	UNP P0CT06
D	559	HIS	-	expression tag	UNP P0CT06
D	560	HIS	-	expression tag	UNP P0CT06

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Mg 1 1	0	0
2	A	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0

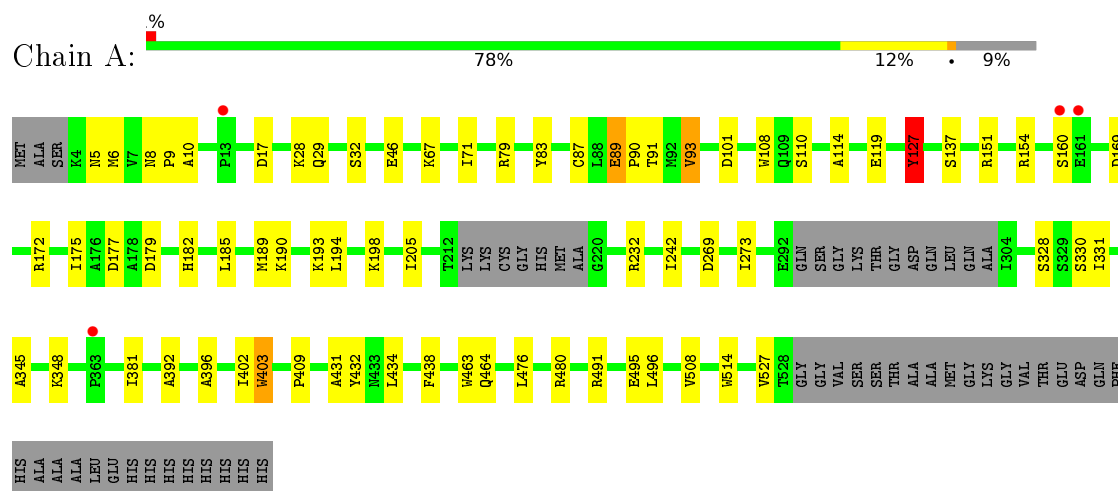
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	20	Total 20	O 20	0	0
3	B	16	Total 16	O 16	0	0
3	C	17	Total 17	O 17	0	0
3	D	12	Total 12	O 12	0	0

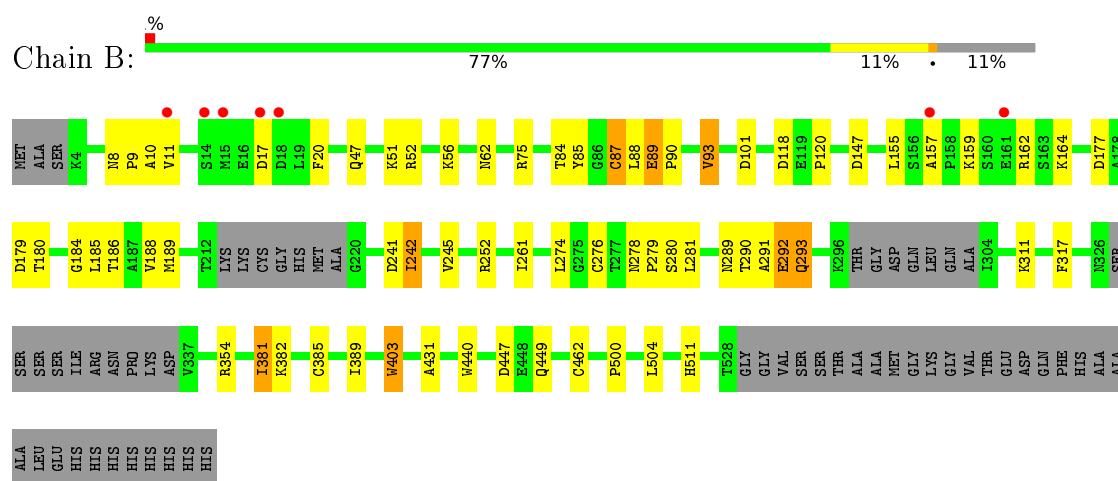
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

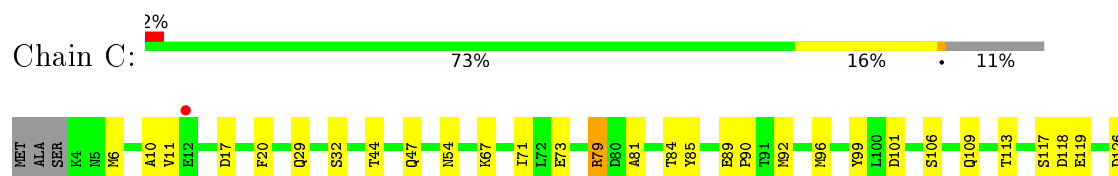
#### • Molecule 1: Isocitrate lyase

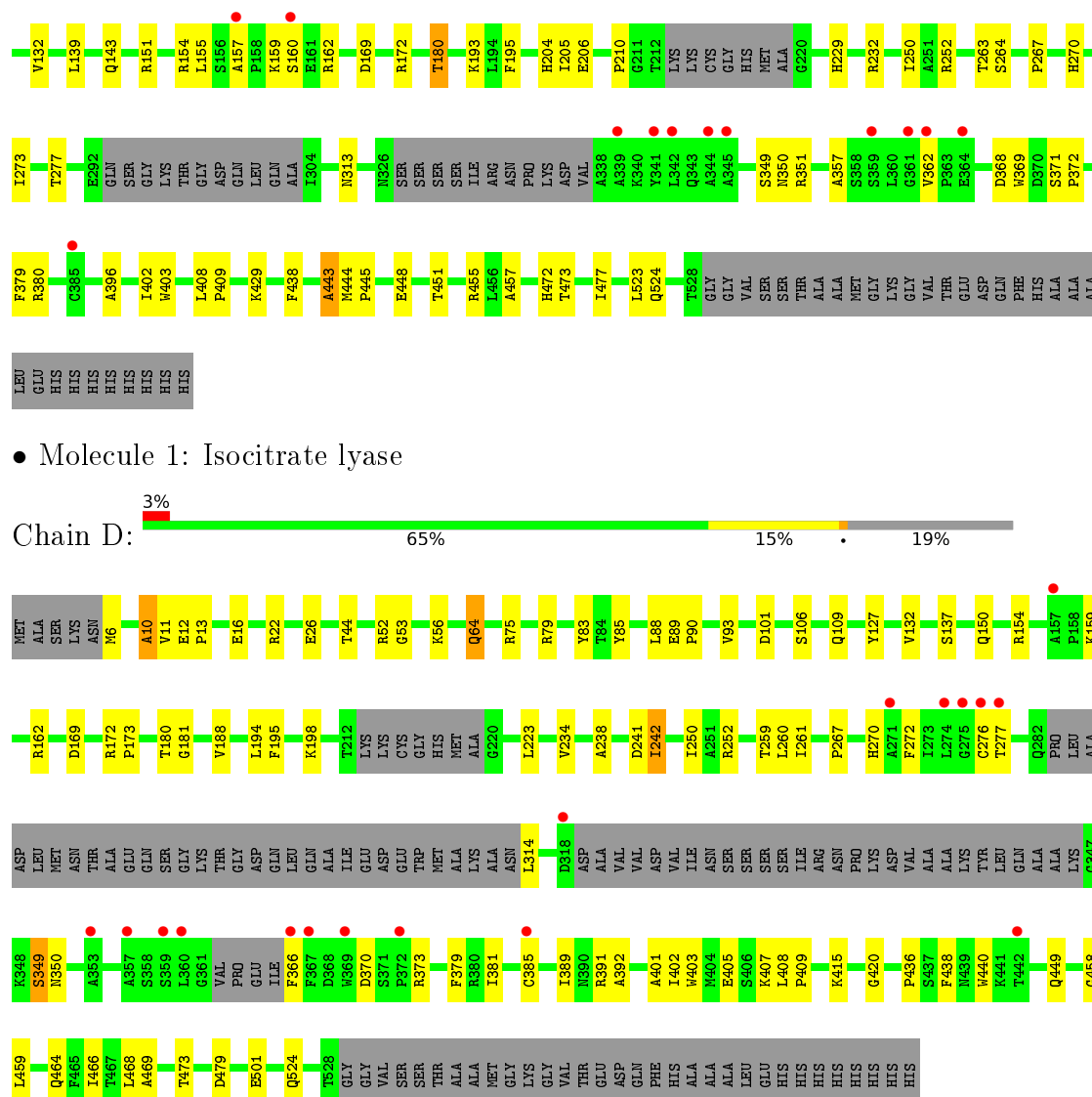


#### • Molecule 1: Isocitrate lyase



#### • Molecule 1: Isocitrate lyase







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	121.58Å 135.31Å 158.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.40 – 2.80 43.39 – 2.81	Depositor EDS
% Data completeness (in resolution range)	99.9 (32.40-2.80) 95.3 (43.39-2.81)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.72 (at 2.81Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.9_1692)	Depositor
R, $R_{free}$	0.186 , 0.226 0.205 , 0.262	Depositor DCC
$R_{free}$ test set	1905 reflections (3.10%)	DCC
Wilson B-factor (Å <sup>2</sup> )	47.6	Xtriage
Anisotropy	0.569	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 44.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	14983	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.62% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: MG

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	0.47	1/3942 (0.0%)	0.59	0/5371
1	B	0.43	1/3942 (0.0%)	0.59	2/5361 (0.0%)
1	C	0.43	0/3822	0.57	0/5213
1	D	0.44	0/3556	0.57	0/4840
All	All	0.44	2/15262 (0.0%)	0.58	2/20785 (0.0%)

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	B	0	1
1	C	0	1
1	D	0	1
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	127	TYR	CD1-CE1	-6.59	1.29	1.39
1	B	87	CYS	CB-SG	-5.38	1.73	1.81

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	292	GLU	N-CA-C	5.50	125.85	111.00
1	B	403	TRP	CA-CB-CG	5.21	123.60	113.70

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	293	GLN	Peptide
1	C	443	ALA	Peptide
1	D	88	LEU	Peptide

## 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3851	0	3626	53	0
1	B	3853	0	3687	47	0
1	C	3736	0	3502	58	0
1	D	3474	0	3268	61	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	20	0	0	0	0
3	B	16	0	0	0	0
3	C	17	0	0	1	0
3	D	12	0	0	0	0
All	All	14983	0	14083	184	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 6.

All (184) 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:D:159:LYS:HA	1:D:162:ARG:HG3	1.58	0.84
1:A:127:TYR:OH	1:A:179:ASP:OD2	1.96	0.82
1:C:444:MET:HG2	1:C:448:GLU:HB2	1.63	0.81
1:D:64:GLN:HE21	1:D:64:GLN:N	1.79	0.80
1:D:389:ILE:HD13	1:D:420:GLY:HA3	1.68	0.76
1:C:157:ALA:O	1:C:162:ARG:NH1	2.21	0.74
1:B:440:TRP:O	1:B:449:GLN:NE2	2.20	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:370:ASP:HB3	1:D:373:ARG:HH21	1.54	0.72
1:C:349:SER:O	1:C:351:ARG:N	2.24	0.69
1:A:193:LYS:NZ	1:C:119:GLU:OE1	2.26	0.69
1:A:101:ASP:H	1:D:6:MET:HE2	1.62	0.65
1:A:83:TYR:O	1:D:6:MET:HE1	1.96	0.65
1:C:277:THR:OG1	1:C:313:ASN:O	2.16	0.63
1:B:10:ALA:HB2	1:C:169:ASP:HA	1.81	0.63
1:B:274:LEU:HD21	1:B:382:LYS:HB2	1.78	0.63
1:D:64:GLN:CA	1:D:64:GLN:HE21	2.12	0.62
1:B:84:THR:OG1	1:B:85:TYR:N	2.28	0.61
1:B:89:GLU:HG2	1:B:90:PRO:HD2	1.82	0.61
1:B:87:CYS:SG	1:B:93:VAL:HG23	2.44	0.58
1:A:403:TRP:HB3	1:A:431:ALA:O	2.03	0.58
1:B:185:LEU:O	1:B:189:MET:HG3	2.03	0.58
1:A:108:TRP:CH2	1:B:511:HIS:HB2	2.39	0.57
1:C:206:GLU:O	1:C:232:ARG:NH1	2.37	0.57
1:B:261:ILE:HB	1:B:381:ILE:HD12	1.87	0.57
1:C:472:HIS:ND1	1:D:501:GLU:OE2	2.38	0.57
1:A:110:SER:HA	1:A:114:ALA:HB3	1.86	0.56
1:B:101:ASP:H	1:C:6:MET:HE2	1.70	0.56
1:C:159:LYS:HA	1:C:162:ARG:HG3	1.88	0.56
1:A:119:GLU:OE1	1:C:193:LYS:NZ	2.38	0.56
1:D:440:TRP:O	1:D:449:GLN:NE2	2.38	0.56
1:D:276:CYS:HA	1:D:314:LEU:HD23	1.87	0.55
1:B:75:ARG:HB3	1:B:462:CYS:HB2	1.89	0.55
1:B:289:ASN:O	1:B:291:ALA:N	2.39	0.55
1:D:349:SER:OG	1:D:350:ASN:N	2.40	0.55
1:C:229:HIS:HA	1:C:232:ARG:HD2	1.90	0.54
1:A:151:ARG:HD3	1:D:52:ARG:O	2.07	0.54
1:B:289:ASN:C	1:B:291:ALA:H	2.11	0.54
1:D:64:GLN:CA	1:D:64:GLN:NE2	2.70	0.54
1:A:137:SER:HB2	1:A:198:LYS:HB3	1.89	0.53
1:A:89:GLU:OE1	1:B:88:LEU:HD13	2.08	0.53
1:A:83:TYR:HA	1:A:464:GLN:O	2.09	0.53
1:A:177:ASP:OD1	1:A:179:ASP:N	2.38	0.52
1:B:157:ALA:O	1:B:162:ARG:NH1	2.43	0.52
1:B:317:PHE:CZ	1:B:354:ARG:HG2	2.45	0.52
1:B:62:ASN:HB2	1:B:245:VAL:HG12	1.92	0.52
1:A:527:VAL:HG13	1:D:234:VAL:HG12	1.92	0.52
1:C:252:ARG:HD2	1:C:403:TRP:CZ2	2.45	0.52
1:A:89:GLU:HG3	1:A:90:PRO:N	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17:ASP:OD1	1:C:162:ARG:NH2	2.44	0.51
1:C:263:THR:HA	1:C:379:PHE:CE1	2.45	0.51
1:C:89:GLU:HG2	1:C:90:PRO:HD2	1.91	0.51
1:A:17:ASP:OD1	1:D:162:ARG:NH2	2.44	0.51
1:D:154:ARG:O	1:D:162:ARG:HD3	2.10	0.51
1:D:22:ARG:O	1:D:26:GLU:HG3	2.11	0.51
1:D:267:PRO:HA	1:D:270:HIS:CE1	2.46	0.50
1:D:436:PRO:HG3	1:D:466:ILE:HG22	1.93	0.50
1:D:64:GLN:HA	1:D:64:GLN:NE2	2.26	0.50
1:A:190:LYS:NZ	1:C:126:ASP:OD1	2.44	0.50
1:B:241:ASP:O	1:B:242:ILE:HB	2.12	0.50
1:C:44:THR:OG1	1:C:47:GLN:HG3	2.11	0.50
1:C:205:ILE:HG23	1:C:229:HIS:CE1	2.47	0.50
1:B:186:THR:HG23	1:C:524:GLN:HE21	1.77	0.50
1:A:508:VAL:HG22	1:A:514:TRP:CE2	2.47	0.50
1:B:189:MET:HE1	1:C:523:LEU:HB3	1.93	0.49
1:C:368:ASP:OD2	1:C:371:SER:OG	2.31	0.49
1:D:106:SER:HB3	1:D:109:GLN:HB2	1.94	0.49
1:D:83:TYR:HA	1:D:464:GLN:O	2.13	0.49
1:B:162:ARG:NH2	1:C:17:ASP:OD1	2.46	0.48
1:C:101:ASP:HA	1:C:172:ARG:HD2	1.95	0.48
1:C:477:ILE:HD13	1:D:473:THR:O	2.13	0.48
1:A:495:GLU:HB3	1:A:496:LEU:HD12	1.95	0.48
1:B:164:LYS:HA	1:B:164:LYS:HD2	1.61	0.48
1:B:155:LEU:HD13	1:C:20:PHE:CD1	2.48	0.48
1:A:409:PRO:HD3	1:A:438:PHE:CG	2.49	0.48
1:A:205:ILE:HD13	1:A:232:ARG:HB3	1.95	0.48
1:D:267:PRO:HA	1:D:270:HIS:ND1	2.28	0.48
1:A:28:LYS:NZ	1:A:46:GLU:OE2	2.47	0.48
1:C:180:THR:HG21	1:C:210:PRO:HB3	1.94	0.47
1:A:8:ASN:ND2	1:D:169:ASP:OD2	2.47	0.47
1:B:385:CYS:O	1:B:389:ILE:HG13	2.15	0.47
1:D:385:CYS:O	1:D:389:ILE:HG13	2.15	0.47
1:B:500:PRO:O	1:B:504:LEU:HG	2.14	0.47
1:A:101:ASP:HA	1:A:172:ARG:HD2	1.97	0.47
1:B:147:ASP:OD1	1:C:54:ASN:ND2	2.47	0.47
1:C:372:PRO:HB2	1:C:380:ARG:HG3	1.96	0.47
1:A:5:ASN:O	1:D:75:ARG:NH2	2.44	0.47
1:A:87:CYS:SG	1:A:93:VAL:HG23	2.55	0.46
1:C:451:THR:O	1:C:455:ARG:HG3	2.14	0.46
1:B:47:GLN:O	1:B:51:LYS:HG2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:259:THR:O	1:D:260:LEU:HD23	2.16	0.46
1:C:73:GLU:OE2	1:C:429:LYS:NZ	2.39	0.46
1:D:89:GLU:HG2	1:D:90:PRO:HD2	1.97	0.46
1:A:396:ALA:HA	1:A:402:ILE:HD11	1.97	0.46
1:A:269:ASP:O	1:A:273:ILE:HG13	2.15	0.46
1:A:29:GLN:O	1:A:32:SER:OG	2.24	0.46
1:C:273:ILE:HG22	1:C:369:TRP:CD1	2.51	0.46
1:B:159:LYS:HA	1:B:162:ARG:HG3	1.97	0.45
1:D:261:ILE:HG22	1:D:379:PHE:HB2	1.98	0.45
1:A:169:ASP:HA	1:D:10:ALA:HB2	1.97	0.45
1:C:132:VAL:HG12	1:C:195:PHE:HE1	1.82	0.45
1:C:84:THR:OG1	1:C:85:TYR:N	2.48	0.45
1:B:101:ASP:H	1:C:6:MET:CE	2.29	0.45
1:B:118:ASP:O	1:B:120:PRO:HD3	2.17	0.45
1:B:52:ARG:O	1:C:151:ARG:HD3	2.17	0.45
1:A:476:LEU:O	1:A:480:ARG:HG2	2.17	0.45
1:D:194:LEU:O	1:D:198:LYS:HG2	2.17	0.45
1:B:184:GLY:O	1:B:188:VAL:HG23	2.17	0.45
1:B:8:ASN:OD1	1:C:172:ARG:NH1	2.47	0.44
1:C:139:LEU:O	1:C:143:GLN:HG3	2.17	0.44
1:A:409:PRO:HD3	1:A:438:PHE:CD2	2.53	0.44
1:B:403:TRP:HB3	1:B:431:ALA:O	2.16	0.44
1:C:396:ALA:HA	1:C:402:ILE:HD11	2.00	0.44
1:D:79:ARG:HB3	1:D:458:GLY:HA2	1.99	0.44
1:A:137:SER:HB2	1:A:198:LYS:HD2	1.99	0.44
1:A:67:LYS:O	1:A:71:ILE:HG13	2.16	0.44
1:A:345:ALA:HA	1:A:348:LYS:HG3	2.00	0.44
1:A:79:ARG:HD3	1:A:79:ARG:HA	1.78	0.44
1:C:445:PRO:HD2	1:C:448:GLU:HG3	2.00	0.44
1:D:241:ASP:O	1:D:242:ILE:HB	2.17	0.43
1:D:392:ALA:HB1	1:D:402:ILE:HG21	2.01	0.43
1:A:432:TYR:HB3	1:A:464:GLN:HG2	1.99	0.43
1:C:357:ALA:HB1	1:C:362:VAL:HB	2.00	0.43
1:A:194:LEU:O	1:A:198:LYS:HG2	2.19	0.43
1:B:278:ASN:ND2	1:B:311:LYS:O	2.48	0.43
1:D:85:TYR:CD2	1:D:469:ALA:HB3	2.54	0.43
1:D:137:SER:HB2	1:D:198:LYS:HD2	2.00	0.43
1:A:127:TYR:HE1	1:A:182:HIS:CE1	2.36	0.43
1:A:6:MET:HE2	1:D:101:ASP:OD1	2.19	0.43
1:C:263:THR:OG1	1:C:264:SER:N	2.51	0.43
1:A:185:LEU:O	1:A:189:MET:HG3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:29:GLN:O	1:C:32:SER:OG	2.34	0.43
1:D:272:PHE:O	1:D:381:ILE:HG13	2.19	0.43
1:C:79:ARG:NH1	3:C:701:HOH:O	2.52	0.42
1:C:109:GLN:O	1:C:113:THR:OG1	2.29	0.42
1:C:117:SER:O	1:C:118:ASP:HB2	2.19	0.42
1:D:172:ARG:HA	1:D:173:PRO:HD3	1.91	0.42
1:B:252:ARG:HD2	1:B:403:TRP:CZ2	2.54	0.42
1:B:56:LYS:HB3	1:B:56:LYS:HE3	1.79	0.42
1:C:67:LYS:O	1:C:71:ILE:HG13	2.19	0.42
1:C:92:MET:O	1:C:96:MET:HG3	2.19	0.42
1:D:276:CYS:SG	1:D:277:THR:N	2.92	0.42
1:B:278:ASN:OD1	1:B:280:SER:OG	2.31	0.42
1:D:52:ARG:HE	1:D:242:ILE:HA	1.85	0.42
1:D:252:ARG:HD2	1:D:403:TRP:CZ2	2.55	0.42
1:D:407:LYS:O	1:D:408:LEU:HD12	2.19	0.42
1:B:177:ASP:OD1	1:B:179:ASP:N	2.41	0.42
1:D:250:ILE:HG12	1:D:401:ALA:HB3	2.01	0.42
1:C:408:LEU:HD23	1:C:409:PRO:HD2	2.02	0.42
1:C:473:THR:O	1:C:477:ILE:HG12	2.20	0.42
1:A:175:ILE:HD12	1:A:463:TRP:CZ2	2.55	0.42
1:A:328:SER:O	1:A:330:SER:HA	2.20	0.42
1:D:238:ALA:O	1:D:242:ILE:HG13	2.19	0.42
1:A:273:ILE:HA	1:A:381:ILE:HG13	2.02	0.42
1:A:89:GLU:CD	1:B:88:LEU:HD13	2.41	0.42
1:D:12:GLU:HA	1:D:13:PRO:HD3	1.94	0.41
1:D:409:PRO:HD3	1:D:438:PHE:CG	2.54	0.41
1:A:154:ARG:NE	1:D:16:GLU:OE2	2.38	0.41
1:A:392:ALA:HB1	1:A:402:ILE:HG21	2.02	0.41
1:C:154:ARG:O	1:C:162:ARG:HD3	2.19	0.41
1:D:436:PRO:HG2	1:D:468:LEU:HB2	2.02	0.41
1:B:278:ASN:HA	1:B:279:PRO:HD3	1.86	0.41
1:D:132:VAL:HG12	1:D:195:PHE:HE1	1.84	0.41
1:D:181:GLY:HA3	1:D:188:VAL:HG22	2.02	0.41
1:C:438:PHE:CE1	1:C:443:ALA:HB2	2.56	0.41
1:C:477:ILE:HD12	1:D:473:THR:HB	2.03	0.41
1:C:81:ALA:HB2	1:C:457:ALA:HB2	2.03	0.41
1:A:6:MET:HE2	1:D:101:ASP:H	1.85	0.41
1:B:20:PHE:CD1	1:C:155:LEU:HD13	2.55	0.41
1:D:223:LEU:HD13	1:D:391:ARG:NH1	2.36	0.41
1:C:267:PRO:HA	1:C:270:HIS:CE1	2.56	0.41
1:C:99:TYR:OH	1:D:479:ASP:OD1	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:277:THR:HA	1:D:366:PHE:HE2	1.86	0.41
1:A:151:ARG:NH1	1:D:53:GLY:O	2.50	0.41
1:B:9:PRO:O	1:B:10:ALA:HB3	2.20	0.40
1:B:274:LEU:HD21	1:B:382:LYS:CB	2.47	0.40
1:D:56:LYS:HA	1:D:56:LYS:HD3	1.96	0.40
1:A:6:MET:HE3	1:D:83:TYR:H	1.86	0.40
1:B:281:LEU:HD23	1:B:281:LEU:HA	1.91	0.40
1:A:127:TYR:HD1	1:A:127:TYR:C	2.25	0.40
1:A:491:ARG:O	1:A:495:GLU:HB2	2.21	0.40
1:B:52:ARG:HE	1:B:242:ILE:HA	1.86	0.40
1:C:204:HIS:HB3	1:C:250:ILE:HB	2.04	0.40
1:A:89:GLU:OE2	1:A:91:THR:OG1	2.19	0.40
1:D:415:LYS:HG3	1:D:459:LEU:HD21	2.04	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	501/560 (90%)	478 (95%)	18 (4%)	5 (1%)	19	52
1	B	493/560 (88%)	472 (96%)	17 (3%)	4 (1%)	24	58
1	C	488/560 (87%)	463 (95%)	21 (4%)	4 (1%)	24	58
1	D	443/560 (79%)	415 (94%)	25 (6%)	3 (1%)	26	62
All	All	1925/2240 (86%)	1828 (95%)	81 (4%)	16 (1%)	24	58

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	9	PRO
1	A	242	ILE

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Mol	Chain	Res	Type
1	B	292	GLU
1	C	350	ASN
1	A	10	ALA
1	B	242	ILE
1	B	290	THR
1	A	434	LEU
1	D	10	ALA
1	B	11	VAL
1	C	10	ALA
1	C	160	SER
1	C	11	VAL
1	D	11	VAL
1	D	242	ILE
1	A	331	ILE

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	381/462 (82%)	376 (99%)	5 (1%)	76	94
1	B	389/462 (84%)	382 (98%)	7 (2%)	66	91
1	C	364/462 (79%)	361 (99%)	3 (1%)	86	97
1	D	347/462 (75%)	338 (97%)	9 (3%)	54	86
All	All	1481/1848 (80%)	1457 (98%)	24 (2%)	70	93

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	89	GLU
1	A	93	VAL
1	A	127	TYR
1	A	160	SER
1	A	403	TRP
1	B	89	GLU
1	B	93	VAL

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Mol	Chain	Res	Type
1	B	180	THR
1	B	276	CYS
1	B	293	GLN
1	B	381	ILE
1	B	447	ASP
1	C	79	ARG
1	C	106	SER
1	C	180	THR
1	D	44	THR
1	D	64	GLN
1	D	93	VAL
1	D	127	TYR
1	D	150	GLN
1	D	180	THR
1	D	349	SER
1	D	405	GLU
1	D	524	GLN

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

Mol	Chain	Res	Type
1	A	64	GLN
1	B	524	GLN
1	C	313	ASN
1	C	413	GLN
1	C	524	GLN
1	D	64	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

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers

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 ⓘ

### 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	507/560 (90%)	-0.17	4 (0%) 87 81	20, 34, 53, 68	0
1	B	501/560 (89%)	-0.08	7 (1%) 78 69	22, 39, 57, 82	0
1	C	496/560 (88%)	-0.08	13 (2%) 59 47	22, 37, 67, 85	0
1	D	453/560 (80%)	0.00	17 (3%) 44 32	23, 38, 73, 93	0
All	All	1957/2240 (87%)	-0.08	41 (2%) 67 56	20, 37, 62, 93	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	277	THR	4.7
1	D	276	CYS	4.6
1	D	367	PHE	4.3
1	D	275	GLY	4.2
1	D	360	LEU	3.9
1	D	385	CYS	3.5
1	D	369	TRP	3.5
1	C	361	GLY	3.4
1	D	353	ALA	3.3
1	C	341	TYR	3.0
1	B	18	ASP	2.9
1	C	345	ALA	2.8
1	C	342	LEU	2.8
1	C	157	ALA	2.8
1	B	14	SER	2.7
1	D	157	ALA	2.7
1	B	15	MET	2.7
1	D	366	PHE	2.6
1	C	359	SER	2.6
1	D	359	SER	2.6
1	A	363	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	12	GLU	2.5
1	C	385	CYS	2.5
1	D	271	ALA	2.5
1	D	357	ALA	2.4
1	D	442	THR	2.4
1	D	318	ASP	2.4
1	A	13	PRO	2.4
1	B	11	VAL	2.3
1	C	339	ALA	2.2
1	C	160	SER	2.2
1	C	344	ALA	2.2
1	D	372	PRO	2.2
1	B	161	GLU	2.2
1	C	362	VAL	2.2
1	B	17	ASP	2.1
1	C	364	GLU	2.1
1	A	161	GLU	2.0
1	A	160	SER	2.0
1	B	157	ALA	2.0
1	D	274	LEU	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	MG	D	601	1/1	0.87	0.53	11.24	36,36,36,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	MG	A	601	1/1	0.82	0.38	7.28	41,41,41,41	0
2	MG	B	601	1/1	0.91	0.26	5.56	30,30,30,30	0
2	MG	C	601	1/1	0.92	0.26	3.35	28,28,28,28	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.