



# Full wwPDB X-ray Structure Validation Report ⓘ

Jul 6, 2019 – 11:10 PM EDT

PDB ID : 3MFB  
Title : Crystal Structure of the S-type Pyocin domain of ECA1669 protein from *Erwinia carotovora*, Northeast Structural Genomics Consortium Target EwR82C  
Authors : Forouhar, F.; Neely, H.; Seetharaman, J.; Sahdev, S.; Xiao, R.; Ciccocanti, C.; Lee, D.; Everett, J.K.; Nair, R.; Acton, T.B.; Rost, B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)  
Deposited on : 2010-04-01  
Resolution : 2.20 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  
Mogul : 1.8.0 (224370), CSD as540be (2019)  
Xtriage (Phenix) : 1.13  
EDS : 2.3.2  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Refmac : 5.8.0158  
CCP4 : 7.0 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.3.2

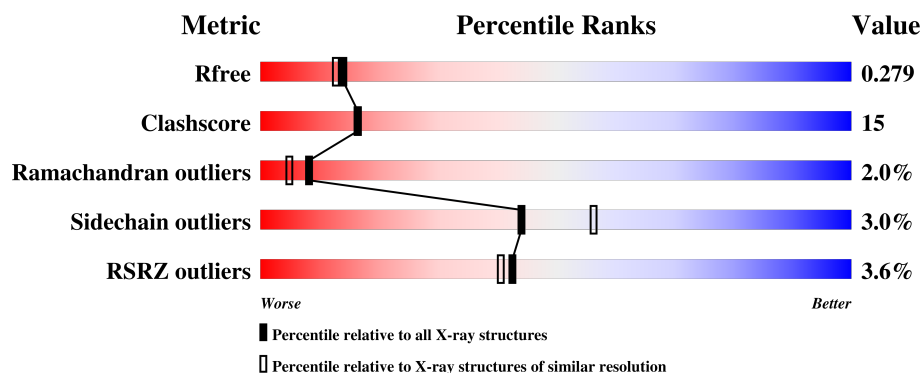
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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))
$R_{free}$	111664	4343 (2.20-2.20)
Clashscore	122126	5027 (2.20-2.20)
Ramachandran outliers	120053	4952 (2.20-2.20)
Sidechain outliers	120020	4953 (2.20-2.20)
RSRZ outliers	108989	4245 (2.20-2.20)

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	157	<div> <div>6%</div> <div>73%19% . .</div> </div>
1	B	157	<div> <div>2%</div> <div>74%18% . .</div> </div>
1	C	157	<div> <div>3%</div> <div>75%17% . .</div> </div>
1	D	157	<div> <div>2%</div> <div>75%17% . .</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5153 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 Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	150	Total	C	N	O	Se	0	0	0
			1205	771	198	231	5			
1	B	150	Total	C	N	O	Se	0	0	0
			1205	771	198	231	5			
1	C	150	Total	C	N	O	Se	0	0	0
			1205	771	198	231	5			
1	D	150	Total	C	N	O	Se	0	0	0
			1205	771	198	231	5			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	331	MSE	-	INITIATING METHIONINE	UNP Q6D6L3
A	480	LEU	-	EXPRESSION TAG	UNP Q6D6L3
A	481	GLU	-	EXPRESSION TAG	UNP Q6D6L3
A	482	HIS	-	EXPRESSION TAG	UNP Q6D6L3
A	483	HIS	-	EXPRESSION TAG	UNP Q6D6L3
A	484	HIS	-	EXPRESSION TAG	UNP Q6D6L3
A	485	HIS	-	EXPRESSION TAG	UNP Q6D6L3
A	486	HIS	-	EXPRESSION TAG	UNP Q6D6L3
A	487	HIS	-	EXPRESSION TAG	UNP Q6D6L3
B	331	MSE	-	INITIATING METHIONINE	UNP Q6D6L3
B	480	LEU	-	EXPRESSION TAG	UNP Q6D6L3
B	481	GLU	-	EXPRESSION TAG	UNP Q6D6L3
B	482	HIS	-	EXPRESSION TAG	UNP Q6D6L3
B	483	HIS	-	EXPRESSION TAG	UNP Q6D6L3
B	484	HIS	-	EXPRESSION TAG	UNP Q6D6L3
B	485	HIS	-	EXPRESSION TAG	UNP Q6D6L3
B	486	HIS	-	EXPRESSION TAG	UNP Q6D6L3
B	487	HIS	-	EXPRESSION TAG	UNP Q6D6L3
C	331	MSE	-	INITIATING METHIONINE	UNP Q6D6L3
C	480	LEU	-	EXPRESSION TAG	UNP Q6D6L3
C	481	GLU	-	EXPRESSION TAG	UNP Q6D6L3

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Chain	Residue	Modelled	Actual	Comment	Reference
C	482	HIS	-	EXPRESSION TAG	UNP Q6D6L3
C	483	HIS	-	EXPRESSION TAG	UNP Q6D6L3
C	484	HIS	-	EXPRESSION TAG	UNP Q6D6L3
C	485	HIS	-	EXPRESSION TAG	UNP Q6D6L3
C	486	HIS	-	EXPRESSION TAG	UNP Q6D6L3
C	487	HIS	-	EXPRESSION TAG	UNP Q6D6L3
D	331	MSE	-	INITIATING METHIONINE	UNP Q6D6L3
D	480	LEU	-	EXPRESSION TAG	UNP Q6D6L3
D	481	GLU	-	EXPRESSION TAG	UNP Q6D6L3
D	482	HIS	-	EXPRESSION TAG	UNP Q6D6L3
D	483	HIS	-	EXPRESSION TAG	UNP Q6D6L3
D	484	HIS	-	EXPRESSION TAG	UNP Q6D6L3
D	485	HIS	-	EXPRESSION TAG	UNP Q6D6L3
D	486	HIS	-	EXPRESSION TAG	UNP Q6D6L3
D	487	HIS	-	EXPRESSION TAG	UNP Q6D6L3

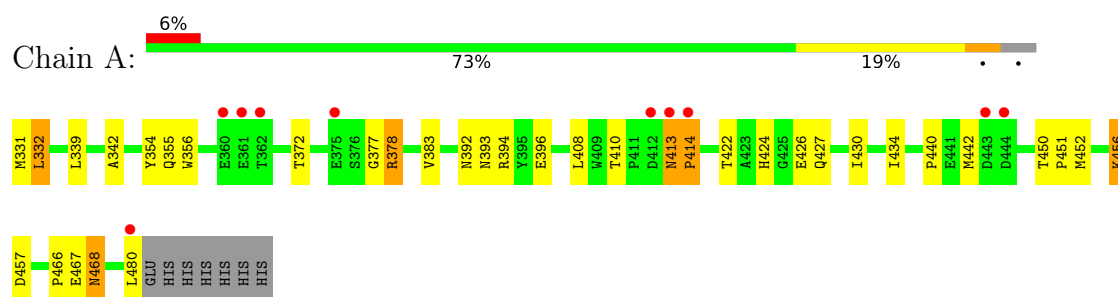
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	78	Total O 78 78	0	0
2	B	90	Total O 90 90	0	0
2	C	73	Total O 73 73	0	0
2	D	92	Total O 92 92	0	0

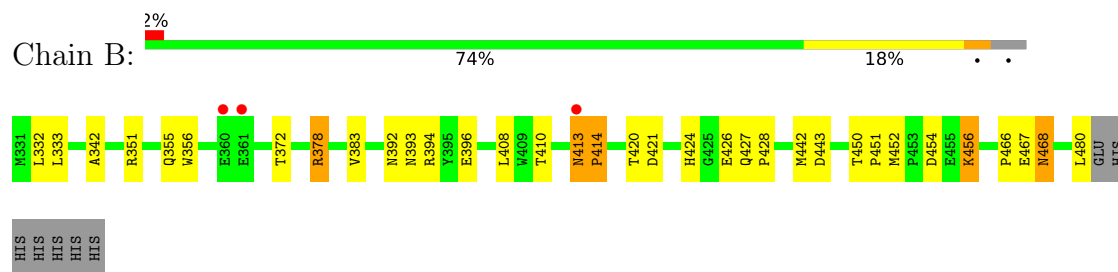
### 3 Residue-property plots

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.

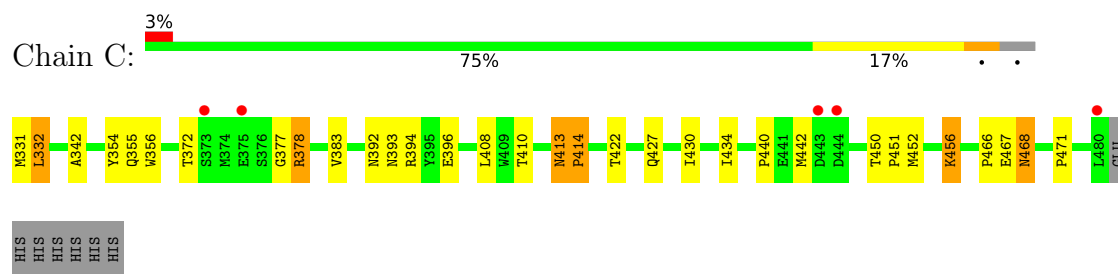
- Molecule 1: Uncharacterized protein



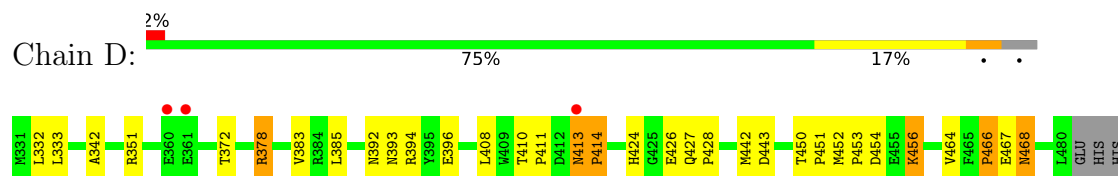
- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	50.08Å 52.31Å 62.16Å 69.50° 90.04° 90.04°	Depositor
Resolution (Å)	19.41 – 2.20 27.16 – 2.19	Depositor EDS
% Data completeness (in resolution range)	87.2 (19.41-2.20) 93.3 (27.16-2.19)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.36 (at 2.20Å)	Xtriage
Refinement program	CNS 1.2 & XtalView, REFMAC	Depositor
R, $R_{free}$	0.232 , 0.265 0.245 , 0.279	Depositor DCC
$R_{free}$ test set	2734 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	14.9	Xtriage
Anisotropy	0.770	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 19.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.477 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	5153	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.73% 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	0.37	0/1236	0.61	0/1681
1	B	0.37	0/1236	0.60	0/1681
1	C	0.37	0/1236	0.60	0/1681
1	D	0.37	0/1236	0.60	0/1681
All	All	0.37	0/4944	0.60	0/6724

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	1205	0	1170	39	0
1	B	1205	0	1170	38	0
1	C	1205	0	1170	37	0
1	D	1205	0	1170	42	0
2	A	78	0	0	6	0
2	B	90	0	0	4	0
2	C	73	0	0	6	0
2	D	92	0	0	8	0
All	All	5153	0	4680	145	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 15.



All (145) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:378:ARG:HH21	1:B:378:ARG:HB2	1.20	1.06
1:A:378:ARG:HB2	1:A:378:ARG:HH21	1.20	1.05
1:D:378:ARG:HH21	1:D:378:ARG:HB2	1.21	1.05
1:C:378:ARG:HB2	1:C:378:ARG:HH21	1.21	1.04
1:A:456:LYS:H	1:A:456:LYS:HE2	1.32	0.94
1:C:456:LYS:H	1:C:456:LYS:HE2	1.32	0.94
1:B:456:LYS:HE2	1:B:456:LYS:H	1.32	0.93
1:D:456:LYS:HE2	1:D:456:LYS:H	1.33	0.92
1:B:378:ARG:CB	1:B:378:ARG:HH21	1.97	0.78
1:D:378:ARG:HH21	1:D:378:ARG:CB	1.97	0.78
1:C:378:ARG:CB	1:C:378:ARG:HH21	1.96	0.77
1:A:378:ARG:HH21	1:A:378:ARG:CB	1.97	0.76
1:D:450:THR:OG1	1:D:451:PRO:HD2	1.86	0.75
1:B:450:THR:OG1	1:B:451:PRO:HD2	1.87	0.73
1:C:450:THR:OG1	1:C:451:PRO:HD2	1.88	0.73
1:A:450:THR:OG1	1:A:451:PRO:HD2	1.89	0.72
1:A:413:ASN:HB2	1:A:414:PRO:HA	1.75	0.68
1:C:413:ASN:HB2	1:C:414:PRO:HA	1.76	0.68
1:A:378:ARG:HB2	1:A:378:ARG:NH2	2.04	0.67
1:C:378:ARG:HB2	1:C:378:ARG:NH2	2.04	0.67
1:D:413:ASN:HB2	1:D:414:PRO:HA	1.76	0.67
1:B:413:ASN:HB2	1:B:414:PRO:HA	1.76	0.66
1:B:392:ASN:O	1:B:394:ARG:HG3	1.96	0.66
1:B:443:ASP:OD1	1:C:331:MSE:N	2.29	0.66
1:D:392:ASN:O	1:D:394:ARG:HG3	1.96	0.65
1:A:339:LEU:HD13	2:A:267:HOH:O	1.95	0.65
1:A:331:MSE:N	1:D:443:ASP:OD1	2.31	0.64
1:A:392:ASN:O	1:A:394:ARG:HG3	1.97	0.64
1:C:450:THR:HG21	2:C:102:HOH:O	1.99	0.63
1:B:378:ARG:NH2	1:B:378:ARG:HB2	2.04	0.63
1:A:422:THR:HG21	2:A:212:HOH:O	1.98	0.63
1:B:428:PRO:HG2	1:C:430:ILE:HG21	1.80	0.63
1:A:450:THR:HG21	2:A:49:HOH:O	1.99	0.63
1:D:378:ARG:NH2	1:D:378:ARG:HB2	2.04	0.62
1:A:430:ILE:HG21	1:D:428:PRO:HG2	1.80	0.62
1:C:392:ASN:O	1:C:394:ARG:HG3	1.99	0.62
1:D:378:ARG:HD2	1:D:452:MSE:CG	2.30	0.62
1:B:378:ARG:HD2	1:B:452:MSE:CG	2.30	0.62
1:B:421:ASP:HB2	1:D:467:GLU:CD	2.20	0.61
1:C:378:ARG:HD2	1:C:452:MSE:CG	2.29	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:452:MSE:HE3	2:D:154:HOH:O	1.99	0.61
1:A:378:ARG:HD2	1:A:452:MSE:CG	2.30	0.61
1:C:378:ARG:HD2	1:C:452:MSE:HG3	1.83	0.59
1:D:378:ARG:HD2	1:D:452:MSE:HG3	1.84	0.59
1:A:378:ARG:HD2	1:A:452:MSE:HG3	1.83	0.59
1:B:378:ARG:HD2	1:B:452:MSE:HG3	1.84	0.59
1:D:411:PRO:HG2	2:D:23:HOH:O	2.03	0.58
1:B:351:ARG:HG3	2:B:143:HOH:O	2.05	0.57
1:A:396:GLU:HB3	1:A:408:LEU:HD23	1.87	0.57
1:A:413:ASN:HB2	1:A:414:PRO:CA	2.35	0.56
1:C:413:ASN:HB2	1:C:414:PRO:CA	2.35	0.56
1:A:427:GLN:HB3	1:A:442:MSE:HE1	1.88	0.56
1:C:396:GLU:HB3	1:C:408:LEU:HD23	1.88	0.56
1:C:422:THR:HG21	2:C:222:HOH:O	2.05	0.56
1:D:413:ASN:HB2	1:D:414:PRO:CA	2.35	0.56
1:B:413:ASN:HB2	1:B:414:PRO:CA	2.36	0.55
1:C:427:GLN:HB3	1:C:442:MSE:HE1	1.89	0.55
1:D:396:GLU:HB3	1:D:408:LEU:HD23	1.89	0.55
1:B:396:GLU:HB3	1:B:408:LEU:HD23	1.89	0.55
1:A:393:ASN:O	1:A:410:THR:HG23	2.06	0.54
2:B:326:HOH:O	1:C:332:LEU:HD22	2.07	0.54
1:A:427:GLN:HB2	1:A:442:MSE:SE	2.58	0.54
1:D:427:GLN:HB2	1:D:442:MSE:SE	2.59	0.53
1:B:421:ASP:HB2	1:D:467:GLU:OE1	2.09	0.53
1:C:393:ASN:O	1:C:410:THR:HG23	2.08	0.53
1:B:421:ASP:HB2	1:D:467:GLU:HG2	1.91	0.52
1:B:427:GLN:HB2	1:B:442:MSE:SE	2.60	0.52
1:D:351:ARG:HG3	2:D:121:HOH:O	2.10	0.52
1:C:427:GLN:HB2	1:C:442:MSE:SE	2.60	0.52
1:B:427:GLN:HB3	1:B:442:MSE:HE1	1.92	0.51
1:D:427:GLN:HB3	1:D:442:MSE:HE1	1.92	0.51
1:A:372:THR:HA	1:A:452:MSE:CE	2.41	0.51
1:B:372:THR:HA	1:B:452:MSE:CE	2.41	0.51
1:D:393:ASN:O	1:D:410:THR:HG23	2.10	0.51
1:C:372:THR:HA	1:C:452:MSE:CE	2.41	0.51
1:B:393:ASN:O	1:B:410:THR:HG23	2.10	0.50
1:D:453:PRO:HG3	2:D:100:HOH:O	2.11	0.50
1:B:421:ASP:HB2	1:D:467:GLU:CG	2.42	0.50
1:D:372:THR:HA	1:D:452:MSE:CE	2.41	0.50
1:D:466:PRO:HG2	2:D:229:HOH:O	2.12	0.49
1:A:468:ASN:ND2	2:A:189:HOH:O	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:332:LEU:HD22	2:D:220:HOH:O	2.12	0.49
1:B:427:GLN:NE2	1:C:434:ILE:HG12	2.28	0.49
1:A:434:ILE:HG12	1:D:427:GLN:NE2	2.28	0.48
1:A:480:LEU:HD23	2:A:292:HOH:O	2.13	0.48
1:C:377:GLY:HA3	2:C:54:HOH:O	2.12	0.47
1:D:467:GLU:N	2:D:229:HOH:O	2.46	0.47
1:B:420:THR:HA	1:D:467:GLU:OE2	2.15	0.47
1:A:456:LYS:CE	1:A:456:LYS:H	2.16	0.47
1:C:471:PRO:HB2	2:C:328:HOH:O	2.15	0.46
1:C:456:LYS:CE	1:C:456:LYS:H	2.16	0.46
1:B:450:THR:HG21	2:B:59:HOH:O	2.15	0.46
1:C:378:ARG:HD2	1:C:452:MSE:HG2	1.98	0.46
1:B:378:ARG:HD2	1:B:452:MSE:HG2	1.98	0.45
1:D:378:ARG:HD2	1:D:452:MSE:HG2	1.98	0.45
1:D:342:ALA:HB2	1:D:383:VAL:HG12	1.99	0.45
1:B:342:ALA:HB2	1:B:383:VAL:HG12	1.99	0.45
1:D:351:ARG:NH1	2:D:111:HOH:O	2.50	0.45
1:D:450:THR:OG1	1:D:451:PRO:CD	2.62	0.45
1:C:372:THR:HA	1:C:452:MSE:HE1	1.99	0.44
1:A:372:THR:HA	1:A:452:MSE:HE1	1.99	0.44
1:B:372:THR:HA	1:B:452:MSE:HE1	1.99	0.44
1:B:450:THR:OG1	1:B:451:PRO:CD	2.62	0.44
1:C:450:THR:OG1	1:C:451:PRO:CD	2.63	0.44
1:C:440:PRO:O	1:C:442:MSE:HE2	2.17	0.44
1:C:393:ASN:ND2	2:C:223:HOH:O	2.51	0.44
1:C:378:ARG:CG	1:C:378:ARG:HH21	2.30	0.43
1:D:378:ARG:HH21	1:D:378:ARG:CG	2.31	0.43
1:D:413:ASN:CB	1:D:414:PRO:CA	2.97	0.43
1:B:378:ARG:HH21	1:B:378:ARG:CG	2.31	0.43
1:C:342:ALA:HB2	1:C:383:VAL:HG12	2.01	0.43
1:A:377:GLY:HA3	2:A:243:HOH:O	2.16	0.43
1:A:450:THR:OG1	1:A:451:PRO:CD	2.64	0.43
1:C:354:TYR:CE1	1:C:442:MSE:HG2	2.54	0.43
1:B:413:ASN:CB	1:B:414:PRO:CA	2.97	0.43
1:D:372:THR:HA	1:D:452:MSE:HE1	2.00	0.43
1:A:378:ARG:HH21	1:A:378:ARG:CG	2.31	0.43
1:B:480:LEU:HD11	2:B:128:HOH:O	2.19	0.42
1:A:354:TYR:CE1	1:A:442:MSE:HG2	2.55	0.42
1:A:440:PRO:O	1:A:442:MSE:HE2	2.19	0.42
1:D:424:HIS:HD2	1:D:426:GLU:OE1	2.03	0.42
1:A:413:ASN:CB	1:A:414:PRO:CA	2.97	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:413:ASN:CB	1:C:414:PRO:CA	2.97	0.42
1:D:333:LEU:HD12	1:D:333:LEU:HA	1.84	0.42
1:B:467:GLU:O	1:B:468:ASN:HB2	2.20	0.42
1:A:378:ARG:HD2	1:A:452:MSE:HG2	1.98	0.41
1:B:333:LEU:HD12	1:B:333:LEU:HA	1.84	0.41
1:C:452:MSE:HE2	2:C:225:HOH:O	2.20	0.41
1:C:467:GLU:O	1:C:468:ASN:HB2	2.20	0.41
1:A:342:ALA:HB2	1:A:383:VAL:HG12	2.02	0.41
1:B:424:HIS:HD2	1:B:426:GLU:OE1	2.03	0.41
1:D:467:GLU:O	1:D:468:ASN:HB2	2.20	0.41
1:D:454:ASP:HB3	1:D:456:LYS:CE	2.51	0.41
1:A:424:HIS:HD2	1:A:426:GLU:OE1	2.02	0.41
1:A:355:GLN:NE2	1:A:457:ASP:HB3	2.35	0.41
1:B:454:ASP:HB3	1:B:456:LYS:CE	2.51	0.41
1:C:355:GLN:HG2	1:C:356:TRP:N	2.36	0.41
1:A:355:GLN:HG2	1:A:356:TRP:N	2.36	0.40
1:A:467:GLU:O	1:A:468:ASN:HB2	2.21	0.40
1:B:372:THR:HA	1:B:452:MSE:HE3	2.03	0.40
1:B:355:GLN:HG2	1:B:356:TRP:N	2.35	0.40
1:D:385:LEU:HD12	1:D:464:VAL:O	2.22	0.40
1:A:456:LYS:N	1:A:456:LYS:HE2	2.15	0.40
1:C:456:LYS:N	1:C:456:LYS:HE2	2.16	0.40
1:D:372:THR:HA	1:D:452:MSE:HE3	2.03	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	148/157 (94%)	139 (94%)	6 (4%)	3 (2%)	8	5
1	B	148/157 (94%)	139 (94%)	6 (4%)	3 (2%)	8	5

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	148/157 (94%)	139 (94%)	6 (4%)	3 (2%)	8	5
1	D	148/157 (94%)	139 (94%)	6 (4%)	3 (2%)	8	5
All	All	592/628 (94%)	556 (94%)	24 (4%)	12 (2%)	8	5

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	414	PRO
1	B	414	PRO
1	C	414	PRO
1	D	414	PRO
1	A	466	PRO
1	B	466	PRO
1	C	466	PRO
1	D	466	PRO
1	A	413	ASN
1	B	413	ASN
1	D	413	ASN
1	C	413	ASN

### 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	133/135 (98%)	129 (97%)	4 (3%)	44	56
1	B	133/135 (98%)	129 (97%)	4 (3%)	44	56
1	C	133/135 (98%)	129 (97%)	4 (3%)	44	56
1	D	133/135 (98%)	129 (97%)	4 (3%)	44	56
All	All	532/540 (98%)	516 (97%)	16 (3%)	44	56

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	332	LEU
1	A	378	ARG
1	A	456	LYS
1	A	468	ASN
1	B	332	LEU
1	B	378	ARG
1	B	456	LYS
1	B	468	ASN
1	C	332	LEU
1	C	378	ARG
1	C	456	LYS
1	C	468	ASN
1	D	332	LEU
1	D	378	ARG
1	D	456	LYS
1	D	468	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	424	HIS
1	B	345	GLN
1	B	424	HIS
1	B	427	GLN
1	C	345	GLN
1	C	393	ASN
1	C	424	HIS
1	D	345	GLN
1	D	388	HIS
1	D	424	HIS
1	D	427	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

### 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	145/157 (92%)	0.07	10 (6%) 17 15	7, 21, 46, 58	0
1	B	145/157 (92%)	-0.09	3 (2%) 63 61	6, 18, 43, 60	0
1	C	145/157 (92%)	0.04	5 (3%) 45 43	6, 21, 45, 57	0
1	D	145/157 (92%)	-0.10	3 (2%) 63 61	6, 18, 43, 60	0
All	All	580/628 (92%)	-0.02	21 (3%) 42 41	6, 20, 45, 60	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	443	ASP	4.5
1	B	360	GLU	4.2
1	A	375	GLU	4.1
1	C	443	ASP	3.8
1	B	413	ASN	3.6
1	D	361	GLU	3.5
1	A	480	LEU	3.4
1	B	361	GLU	3.1
1	C	375	GLU	3.1
1	A	362	THR	2.9
1	A	412	ASP	2.8
1	C	480	LEU	2.7
1	A	444	ASP	2.5
1	A	413	ASN	2.5
1	A	361	GLU	2.4
1	A	414	PRO	2.3
1	A	360	GLU	2.3
1	C	444	ASP	2.3
1	D	360	GLU	2.3
1	D	413	ASN	2.2
1	C	373	SER	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.