



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 19, 2017 – 11:41 PM EDT

PDB ID : 3EXZ  
Title : Crystal structure of the MaoC-like dehydratase from Rhodospirillum rubrum.  
Northeast Structural Genomics Consortium target RrR103a.  
Authors : Vorobiev, S.M.; Su, M.; Seetharaman, J.; Wang, H.; Foote, E.L.; Ciccocanti,  
C.; Janjua, H.; Xiao, R.; Acton, T.B.; Montelione, G.T.; Tong, L.; Hunt, J.F.;  
Northeast Structural Genomics Consortium (NESG)  
Deposited on : unknown  
Resolution : 2.30 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	:	rb-20030345
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)	:	rb-20030345

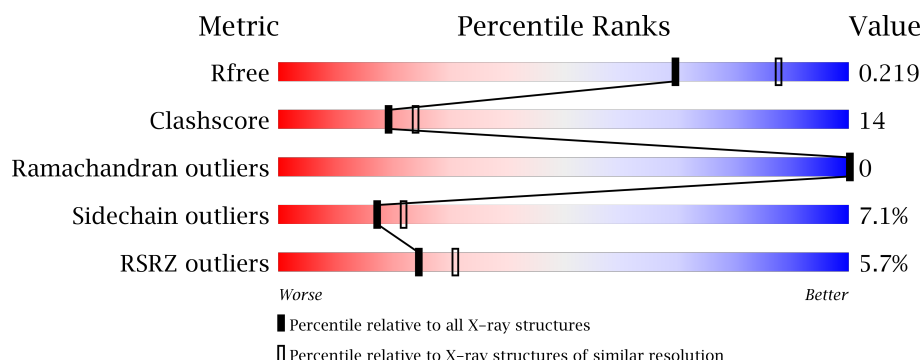
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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	154	<div> <div>2%</div> <div>78% 13% 5%</div> </div>
1	B	154	<div> <div>%</div> <div>80% 16% 5%</div> </div>
1	C	154	<div> <div>22%</div> <div>61% 29% 7%</div> </div>
1	D	154	<div> <div>75% 17% 5%</div> </div>
1	E	154	<div> <div>2%</div> <div>73% 19% 5%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5843 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 MaoC-like dehydratase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	146	Total	C	N	O	S	Se	0	0	0
			1096	690	199	205	1	1			
1	B	147	Total	C	N	O	S	Se	0	0	0
			1116	701	205	208	1	1			
1	C	143	Total	C	N	O	S	Se	0	0	0
			1088	683	201	202	1	1			
1	D	146	Total	C	N	O	S	Se	0	0	0
			1106	695	202	207	1	1			
1	E	146	Total	C	N	O	S	Se	0	0	0
			1088	685	199	202	1	1			

There are 45 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	14	MSE	-	expression tag	UNP Q2RSA1
A	160	LEU	-	expression tag	UNP Q2RSA1
A	161	GLU	-	expression tag	UNP Q2RSA1
A	162	HIS	-	expression tag	UNP Q2RSA1
A	163	HIS	-	expression tag	UNP Q2RSA1
A	164	HIS	-	expression tag	UNP Q2RSA1
A	165	HIS	-	expression tag	UNP Q2RSA1
A	166	HIS	-	expression tag	UNP Q2RSA1
A	167	HIS	-	expression tag	UNP Q2RSA1
B	14	MSE	-	expression tag	UNP Q2RSA1
B	160	LEU	-	expression tag	UNP Q2RSA1
B	161	GLU	-	expression tag	UNP Q2RSA1
B	162	HIS	-	expression tag	UNP Q2RSA1
B	163	HIS	-	expression tag	UNP Q2RSA1
B	164	HIS	-	expression tag	UNP Q2RSA1
B	165	HIS	-	expression tag	UNP Q2RSA1
B	166	HIS	-	expression tag	UNP Q2RSA1
B	167	HIS	-	expression tag	UNP Q2RSA1
C	14	MSE	-	expression tag	UNP Q2RSA1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	160	LEU	-	expression tag	UNP Q2RSA1
C	161	GLU	-	expression tag	UNP Q2RSA1
C	162	HIS	-	expression tag	UNP Q2RSA1
C	163	HIS	-	expression tag	UNP Q2RSA1
C	164	HIS	-	expression tag	UNP Q2RSA1
C	165	HIS	-	expression tag	UNP Q2RSA1
C	166	HIS	-	expression tag	UNP Q2RSA1
C	167	HIS	-	expression tag	UNP Q2RSA1
D	14	MSE	-	expression tag	UNP Q2RSA1
D	160	LEU	-	expression tag	UNP Q2RSA1
D	161	GLU	-	expression tag	UNP Q2RSA1
D	162	HIS	-	expression tag	UNP Q2RSA1
D	163	HIS	-	expression tag	UNP Q2RSA1
D	164	HIS	-	expression tag	UNP Q2RSA1
D	165	HIS	-	expression tag	UNP Q2RSA1
D	166	HIS	-	expression tag	UNP Q2RSA1
D	167	HIS	-	expression tag	UNP Q2RSA1
E	14	MSE	-	expression tag	UNP Q2RSA1
E	160	LEU	-	expression tag	UNP Q2RSA1
E	161	GLU	-	expression tag	UNP Q2RSA1
E	162	HIS	-	expression tag	UNP Q2RSA1
E	163	HIS	-	expression tag	UNP Q2RSA1
E	164	HIS	-	expression tag	UNP Q2RSA1
E	165	HIS	-	expression tag	UNP Q2RSA1
E	166	HIS	-	expression tag	UNP Q2RSA1
E	167	HIS	-	expression tag	UNP Q2RSA1

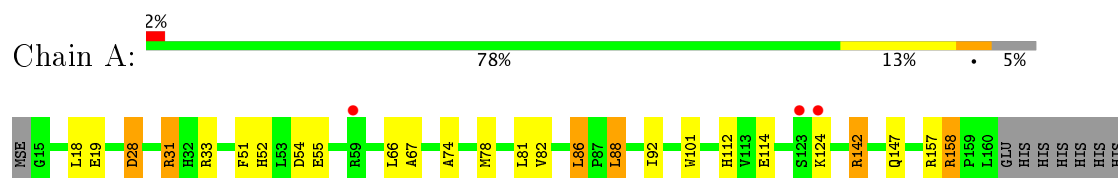
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	47	Total O 47 47	0	0
2	B	91	Total O 91 91	0	0
2	C	27	Total O 27 27	0	0
2	D	106	Total O 106 106	0	0
2	E	78	Total O 78 78	0	0

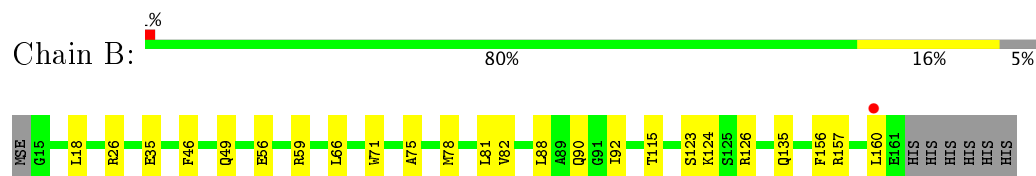
### 3 Residue-property plots [i](#)

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

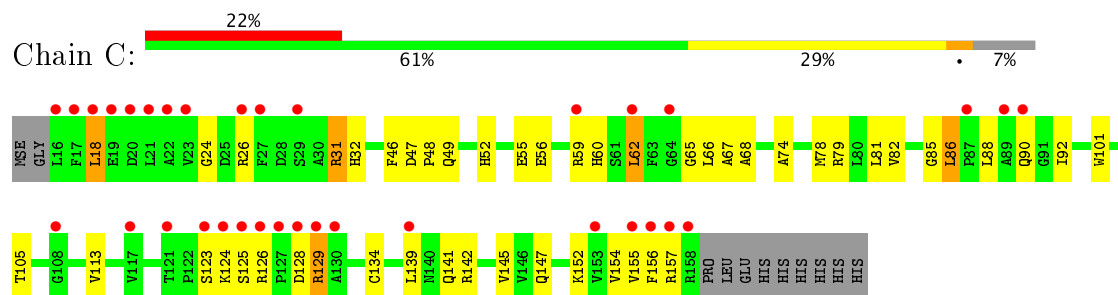
- Molecule 1: MaoC-like dehydratase



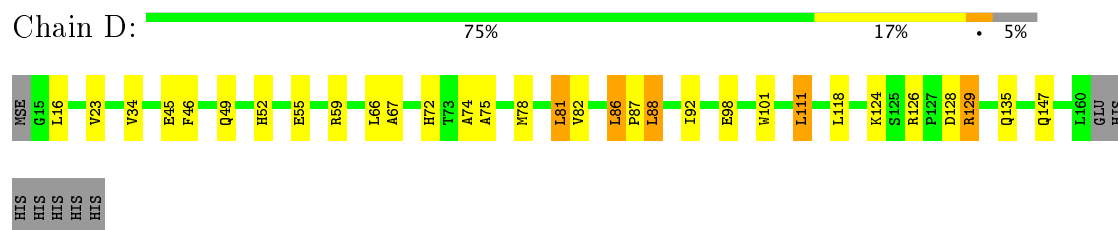
- Molecule 1: MaoC-like dehydratase



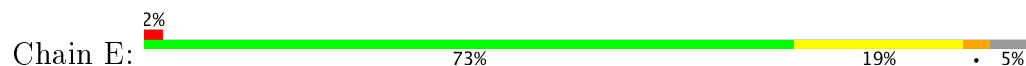
- Molecule 1: MaoC-like dehydratase

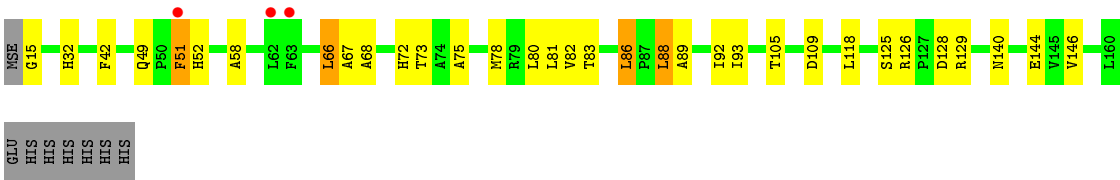


- Molecule 1: MaoC-like dehydratase



- Molecule 1: MaoC-like dehydratase





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	142.95Å 70.51Å 87.86Å 90.00° 93.75° 90.00°	Depositor
Resolution (Å)	38.49 – 2.30 38.49 – 2.29	Depositor EDS
% Data completeness (in resolution range)	91.9 (38.49-2.30) 97.2 (38.49-2.29)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.74 (at 2.29Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.193 , 0.212 0.202 , 0.219	Depositor DCC
$R_{free}$ test set	1874 reflections (4.93%)	DCC
Wilson B-factor (Å <sup>2</sup> )	32.3	Xtriage
Anisotropy	0.257	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 42.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5843	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 5.26% 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	0.42	0/1118	0.67	0/1521
1	B	0.43	0/1139	0.68	0/1548
1	C	0.39	0/1110	0.63	0/1508
1	D	0.43	0/1128	0.70	0/1533
1	E	0.39	0/1110	0.67	0/1511
All	All	0.41	0/5605	0.67	0/7621

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	1096	0	1085	25	0
1	B	1116	0	1107	22	0
1	C	1088	0	1080	57	0
1	D	1106	0	1100	47	0
1	E	1088	0	1069	28	0
2	A	47	0	0	1	0
2	B	91	0	0	0	0
2	C	27	0	0	3	0
2	D	106	0	0	7	0
2	E	78	0	0	1	0
All	All	5843	0	5441	158	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 14.

All (158) 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:78:MSE:HE2	1:B:82:VAL:HG23	1.34	1.09
1:E:78:MSE:HE2	1:E:82:VAL:HG23	1.13	1.06
1:A:78:MSE:CE	1:A:82:VAL:HG23	1.85	1.06
1:A:78:MSE:HE2	1:A:82:VAL:CG2	1.86	1.05
1:A:78:MSE:HE2	1:A:82:VAL:HG23	1.03	1.01
1:B:49:GLN:HE22	1:D:92:ILE:H	1.03	1.00
1:B:92:ILE:H	1:D:49:GLN:HE22	1.01	0.98
1:C:49:GLN:HE22	1:E:92:ILE:H	1.06	0.96
1:D:78:MSE:HE2	1:D:82:VAL:HG23	1.48	0.93
1:D:59:ARG:HD2	2:D:583:HOH:O	1.73	0.89
1:B:78:MSE:HE3	1:B:81:LEU:HB2	1.57	0.85
1:E:78:MSE:HE2	1:E:82:VAL:CG2	2.05	0.82
1:C:78:MSE:HE3	1:C:82:VAL:HG23	1.64	0.80
1:D:124:LYS:HB3	1:D:124:LYS:NZ	1.97	0.79
1:C:78:MSE:HE1	1:C:92:ILE:HD12	1.64	0.79
1:B:78:MSE:HE2	1:B:82:VAL:CG2	2.12	0.78
1:A:55:GLU:OE2	1:C:124:LYS:HB2	1.84	0.78
1:A:101:TRP:HE1	1:A:147:GLN:HE21	1.32	0.78
1:E:78:MSE:HE3	1:E:81:LEU:HB2	1.68	0.75
1:C:49:GLN:NE2	1:E:92:ILE:H	1.82	0.74
1:C:101:TRP:HE1	1:C:147:GLN:HE21	1.33	0.74
1:C:68:ALA:HB3	1:C:105:THR:HB	1.71	0.72
1:C:59:ARG:NH1	1:C:60:HIS:CE1	2.58	0.72
1:E:78:MSE:CE	1:E:82:VAL:HG23	2.06	0.72
1:C:62:LEU:HD12	1:E:93:ILE:HD13	1.71	0.71
1:C:92:ILE:H	1:E:49:GLN:HE22	1.37	0.71
1:C:24:GLY:O	1:C:26:ARG:NH1	2.24	0.70
1:E:78:MSE:O	1:E:78:MSE:HE3	1.91	0.70
1:A:19:GLU:OE2	1:A:157:ARG:HB2	1.91	0.70
1:D:86:LEU:HD22	1:D:88:LEU:HD13	1.73	0.70
1:D:101:TRP:HE1	1:D:147:GLN:HE21	1.39	0.68
1:D:78:MSE:HE2	1:D:82:VAL:CG2	2.24	0.67
1:B:49:GLN:HE22	1:D:92:ILE:N	1.85	0.66
1:C:18:LEU:HB3	1:C:157:ARG:HD2	1.78	0.64
1:D:55:GLU:HG3	2:D:583:HOH:O	1.98	0.64
1:C:59:ARG:NH1	1:C:60:HIS:HE1	1.97	0.62
1:A:19:GLU:OE1	1:A:158:ARG:HD2	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:78:MSE:CE	1:C:82:VAL:CG2	2.77	0.62
1:C:78:MSE:CE	1:C:82:VAL:HG23	2.29	0.62
1:D:124:LYS:HZ3	1:D:124:LYS:HB3	1.65	0.62
1:C:81:LEU:CD1	1:C:113:VAL:HG11	2.30	0.62
1:A:78:MSE:O	1:A:78:MSE:HE3	1.99	0.62
1:C:86:LEU:HD22	1:C:88:LEU:HD13	1.81	0.61
1:B:78:MSE:CE	1:B:82:VAL:HG23	2.22	0.60
1:C:49:GLN:HE22	1:E:92:ILE:N	1.87	0.60
1:C:81:LEU:HD11	1:C:113:VAL:HG11	1.82	0.60
1:D:52:HIS:HD2	1:D:67:ALA:O	1.84	0.60
1:C:79:ARG:HD3	2:C:534:HOH:O	2.02	0.60
1:C:78:MSE:CE	1:C:92:ILE:HD12	2.30	0.59
1:B:92:ILE:N	1:D:49:GLN:HE22	1.85	0.59
1:C:79:ARG:CD	2:C:534:HOH:O	2.51	0.58
1:E:86:LEU:HD22	1:E:88:LEU:HD13	1.85	0.58
1:D:78:MSE:HE1	1:D:92:ILE:HD12	1.85	0.57
1:D:124:LYS:HB3	1:D:124:LYS:HZ2	1.67	0.57
1:D:124:LYS:CB	1:D:124:LYS:NZ	2.67	0.57
1:D:98:GLU:HG3	2:D:239:HOH:O	2.04	0.57
1:D:126:ARG:C	1:D:128:ASP:H	2.07	0.57
1:D:78:MSE:O	1:D:78:MSE:HE3	2.05	0.57
1:E:68:ALA:HB3	1:E:105:THR:HB	1.87	0.56
1:B:123:SER:HB3	1:B:126:ARG:O	2.06	0.56
1:C:78:MSE:HE3	1:C:82:VAL:CG2	2.34	0.55
1:C:123:SER:HB2	1:C:126:ARG:O	2.05	0.55
1:A:19:GLU:CD	1:A:158:ARG:HD2	2.28	0.54
1:C:78:MSE:HE2	1:C:82:VAL:CG2	2.38	0.54
1:B:78:MSE:O	1:B:78:MSE:HE3	2.07	0.54
1:C:126:ARG:C	1:C:128:ASP:H	2.11	0.54
1:A:74:ALA:CA	1:A:147:GLN:HE22	2.20	0.54
1:D:16:LEU:HD12	1:D:16:LEU:N	2.23	0.54
1:C:78:MSE:HE1	1:C:92:ILE:CD1	2.36	0.54
1:E:42:PHE:CD1	1:E:72:HIS:HB2	2.43	0.53
1:C:31:ARG:HD2	1:C:31:ARG:N	2.24	0.53
1:A:78:MSE:HE1	1:A:92:ILE:HD12	1.90	0.53
1:D:16:LEU:HD13	1:D:87:PRO:O	2.09	0.53
1:A:78:MSE:HE1	1:A:92:ILE:CD1	2.40	0.52
1:D:45:GLU:HG2	1:D:46:PHE:CE2	2.44	0.52
1:A:28:ASP:OD2	1:A:114:GLU:HG2	2.10	0.51
1:A:52:HIS:HD2	1:A:67:ALA:O	1.93	0.51
1:D:45:GLU:HG2	1:D:46:PHE:CD2	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:ASP:OD1	1:C:124:LYS:CE	2.59	0.51
1:C:101:TRP:HE1	1:C:147:GLN:NE2	2.04	0.51
1:D:78:MSE:HE3	1:D:81:LEU:CB	2.41	0.51
1:B:78:MSE:CE	1:B:81:LEU:HB2	2.37	0.51
1:E:73:THR:HG21	1:E:146:VAL:CG1	2.41	0.50
1:A:55:GLU:HG3	1:C:125:SER:HB3	1.94	0.50
1:D:126:ARG:C	1:D:128:ASP:N	2.65	0.50
1:A:54:ASP:OD1	1:C:124:LYS:NZ	2.45	0.50
1:D:78:MSE:CE	1:D:81:LEU:HB3	2.42	0.50
1:B:78:MSE:HE1	1:B:92:ILE:CD1	2.42	0.49
1:B:78:MSE:HE1	1:B:92:ILE:HD12	1.93	0.49
1:E:78:MSE:CE	1:E:81:LEU:HB2	2.40	0.49
1:C:18:LEU:HB2	1:C:156:PHE:O	2.13	0.49
1:A:55:GLU:OE2	1:C:124:LYS:CB	2.60	0.48
1:C:139:LEU:CD2	1:C:145:VAL:HG22	2.43	0.48
1:D:129:ARG:NH1	2:D:556:HOH:O	2.46	0.48
1:C:59:ARG:HA	1:C:65:GLY:N	2.29	0.48
1:D:78:MSE:HE1	1:D:92:ILE:CD1	2.44	0.47
1:E:78:MSE:HE1	1:E:92:ILE:CD1	2.44	0.47
1:B:56:GLU:HA	1:B:59:ARG:NH1	2.29	0.47
1:C:126:ARG:C	1:C:128:ASP:N	2.67	0.47
1:C:74:ALA:HA	1:C:147:GLN:HE22	1.80	0.47
1:C:31:ARG:HG2	1:C:31:ARG:HH11	1.79	0.47
1:A:54:ASP:OD1	1:C:124:LYS:HE3	2.14	0.47
1:A:101:TRP:HE1	1:A:147:GLN:NE2	2.07	0.47
1:A:86:LEU:HD22	1:A:88:LEU:HD13	1.97	0.47
1:E:78:MSE:O	1:E:78:MSE:CE	2.63	0.47
1:A:31:ARG:HG3	1:A:112:HIS:HB3	1.97	0.46
1:D:74:ALA:CA	1:D:147:GLN:HE22	2.29	0.46
1:C:124:LYS:HB3	1:C:124:LYS:HZ3	1.80	0.46
1:C:141:GLN:OE1	1:C:141:GLN:N	2.45	0.46
1:C:46:PHE:CG	1:E:75:ALA:HB1	2.51	0.46
1:B:115:THR:HA	1:B:135:GLN:O	2.15	0.45
1:C:129:ARG:HB2	1:C:155:VAL:O	2.17	0.45
1:D:126:ARG:O	1:D:128:ASP:N	2.49	0.45
1:D:98:GLU:O	1:D:98:GLU:HG3	2.16	0.45
1:C:56:GLU:O	1:C:59:ARG:HG2	2.17	0.44
1:D:78:MSE:HE3	1:D:81:LEU:HB2	1.98	0.44
1:A:33:ARG:NH1	2:A:540:HOH:O	2.50	0.44
1:C:78:MSE:CE	1:C:82:VAL:HG22	2.46	0.44
1:C:78:MSE:HE2	1:C:82:VAL:HG22	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:78:MSE:CE	1:D:92:ILE:HD12	2.47	0.44
1:A:74:ALA:HA	1:A:147:GLN:HE22	1.81	0.44
1:B:92:ILE:H	1:D:49:GLN:NE2	1.86	0.44
1:C:52:HIS:HD2	1:C:67:ALA:O	2.01	0.44
1:E:73:THR:HG21	1:E:146:VAL:HG12	1.99	0.44
1:D:23:VAL:HG13	1:E:83:THR:O	2.18	0.43
1:C:31:ARG:CG	1:C:31:ARG:HH11	2.32	0.43
1:E:52:HIS:HD2	1:E:67:ALA:O	2.01	0.43
1:C:74:ALA:CA	1:C:147:GLN:HE22	2.31	0.43
1:C:32:HIS:CE1	2:C:534:HOH:O	2.71	0.43
1:D:129:ARG:CD	2:D:344:HOH:O	2.66	0.43
1:C:47:ASP:N	1:C:48:PRO:HD3	2.31	0.43
1:E:66:LEU:HA	1:E:66:LEU:HD23	1.89	0.43
1:C:78:MSE:O	1:C:82:VAL:HG23	2.19	0.43
1:C:18:LEU:HD23	1:C:18:LEU:O	2.19	0.42
1:B:26:ARG:NH1	1:B:26:ARG:HG2	2.34	0.42
1:C:154:VAL:HG23	1:C:154:VAL:O	2.20	0.42
1:B:90:GLN:NE2	1:B:156:PHE:HB2	2.35	0.42
1:D:16:LEU:N	1:D:16:LEU:CD1	2.82	0.42
1:D:74:ALA:HA	1:D:147:GLN:HE22	1.84	0.42
1:D:78:MSE:HE3	1:D:81:LEU:HB3	2.01	0.42
1:D:118:LEU:HD11	1:D:135:GLN:HB2	2.02	0.42
1:B:75:ALA:HB1	1:D:46:PHE:CG	2.55	0.42
1:C:90:GLN:OE1	1:C:156:PHE:HB2	2.19	0.42
1:B:35:GLU:HA	1:B:35:GLU:OE1	2.20	0.42
1:E:32:HIS:HB2	1:E:80:LEU:HD21	2.02	0.42
1:D:124:LYS:CB	1:D:124:LYS:HZ2	2.30	0.41
1:B:71:TRP:CZ3	1:D:74:ALA:HB1	2.55	0.41
1:D:34:VAL:CG2	1:D:111:LEU:HD22	2.51	0.41
1:B:46:PHE:CG	1:D:75:ALA:HB1	2.56	0.41
1:E:140:ASN:HD21	1:E:144:GLU:HB2	1.85	0.41
1:C:81:LEU:HA	1:C:85:GLY:HA3	2.03	0.41
1:E:15:GLY:HA3	1:E:89:ALA:HB2	2.02	0.40
1:A:142:ARG:HG2	1:A:142:ARG:HH11	1.87	0.40
1:E:126:ARG:C	1:E:128:ASP:H	2.25	0.40
1:E:51:PHE:HA	1:E:58:ALA:HB2	2.03	0.40
1:D:72:HIS:HE1	2:D:270:HOH:O	2.04	0.40
1:D:129:ARG:NE	2:D:344:HOH:O	2.54	0.40
1:E:109:ASP:OD2	2:E:357:HOH:O	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	144/154 (94%)	142 (99%)	2 (1%)	0	100	100
1	B	145/154 (94%)	143 (99%)	2 (1%)	0	100	100
1	C	141/154 (92%)	132 (94%)	9 (6%)	0	100	100
1	D	144/154 (94%)	142 (99%)	2 (1%)	0	100	100
1	E	144/154 (94%)	139 (96%)	5 (4%)	0	100	100
All	All	718/770 (93%)	698 (97%)	20 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	113/122 (93%)	102 (90%)	11 (10%)	9	11
1	B	116/122 (95%)	110 (95%)	6 (5%)	27	36
1	C	113/122 (93%)	103 (91%)	10 (9%)	12	14
1	D	115/122 (94%)	109 (95%)	6 (5%)	27	36
1	E	110/122 (90%)	103 (94%)	7 (6%)	20	27
All	All	567/610 (93%)	527 (93%)	40 (7%)	17	22

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	LEU
1	A	28	ASP
1	A	31	ARG
1	A	51	PHE
1	A	66	LEU
1	A	81	LEU
1	A	86	LEU
1	A	88	LEU
1	A	124	LYS
1	A	142	ARG
1	A	158	ARG
1	B	18	LEU
1	B	66	LEU
1	B	88	LEU
1	B	124	LYS
1	B	157	ARG
1	B	160	LEU
1	C	18	LEU
1	C	31	ARG
1	C	55	GLU
1	C	62	LEU
1	C	66	LEU
1	C	86	LEU
1	C	129	ARG
1	C	134	CYS
1	C	142	ARG
1	C	152	LYS
1	D	66	LEU
1	D	81	LEU
1	D	86	LEU
1	D	88	LEU
1	D	111	LEU
1	D	129	ARG
1	E	51	PHE
1	E	66	LEU
1	E	86	LEU
1	E	88	LEU
1	E	118	LEU
1	E	125	SER
1	E	129	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	HIS
1	A	147	GLN
1	B	49	GLN
1	B	90	GLN
1	C	49	GLN
1	C	52	HIS
1	C	60	HIS
1	C	72	HIS
1	C	103	ASN
1	C	147	GLN
1	D	49	GLN
1	D	52	HIS
1	D	72	HIS
1	D	147	GLN
1	E	49	GLN
1	E	52	HIS
1	E	72	HIS

### 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 ⓘ

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	145/154 (94%)	-0.02	3 (2%) 64 70	22, 38, 56, 68	0
1	B	146/154 (94%)	-0.14	1 (0%) 87 90	17, 29, 51, 59	0
1	C	142/154 (92%)	1.33	34 (23%) 1 1	28, 55, 88, 99	0
1	D	145/154 (94%)	-0.14	0 100 100	18, 27, 48, 58	0
1	E	145/154 (94%)	0.05	3 (2%) 64 70	23, 34, 65, 75	0
All	All	723/770 (93%)	0.21	41 (5%) 24 31	17, 35, 73, 99	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	156	PHE	7.9
1	C	125	SER	7.5
1	C	126	ARG	7.4
1	C	157	ARG	7.1
1	C	124	LYS	7.0
1	C	128	ASP	6.9
1	C	129	ARG	6.6
1	C	21	LEU	5.3
1	C	123	SER	4.8
1	C	130	ALA	4.7
1	C	158	ARG	4.6
1	C	155	VAL	4.3
1	C	27	PHE	4.0
1	C	22	ALA	3.9
1	C	121	THR	3.7
1	C	17	PHE	3.6
1	C	153	VAL	3.5
1	C	20	ASP	3.4
1	C	127	PRO	3.4
1	C	16	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
1	C	89	ALA	3.2
1	C	18	LEU	3.1
1	A	124	LYS	3.0
1	C	90	GLN	3.0
1	B	160	LEU	2.9
1	C	117	VAL	2.7
1	C	26	ARG	2.7
1	C	19	GLU	2.5
1	C	64	GLY	2.5
1	E	51	PHE	2.4
1	E	62	LEU	2.3
1	C	59	ARG	2.3
1	C	108	GLY	2.3
1	C	62	LEU	2.3
1	C	87	PRO	2.2
1	C	139	LEU	2.2
1	C	23	VAL	2.2
1	A	59	ARG	2.1
1	A	123	SER	2.0
1	C	29	SER	2.0
1	E	63	PHE	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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.