



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 15, 2017 – 06:28 AM EDT

PDB ID : 2FQL  
Title : Crystal structure of trimeric frataxin from the yeast *Saccharomyces cerevisiae*  
Authors : Al-Karadaghi, S.; Karlberg, T.  
Deposited on : unknown  
Resolution : 3.01 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

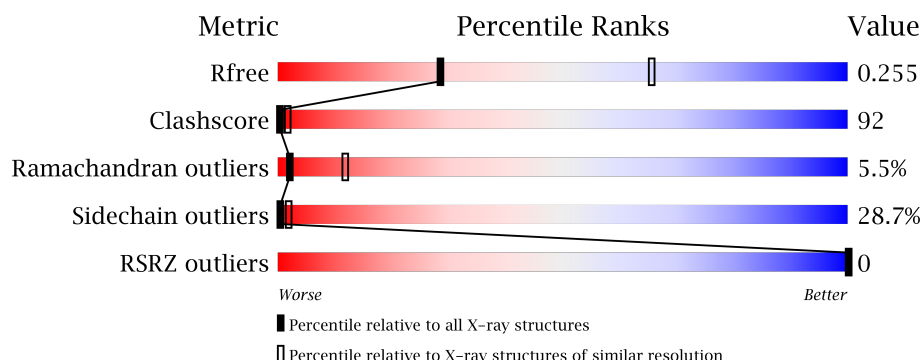
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.01 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	1924 (3.04-3.00)
Clashscore	112137	2279 (3.04-3.00)
Ramachandran outliers	110173	2207 (3.04-3.00)
Sidechain outliers	110143	2210 (3.04-3.00)
RSRZ outliers	101464	1948 (3.04-3.00)

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	123	

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 883 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 Frataxin homolog, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	112	Total	C	N	O	S	3	0	0
			883	561	143	177	2			

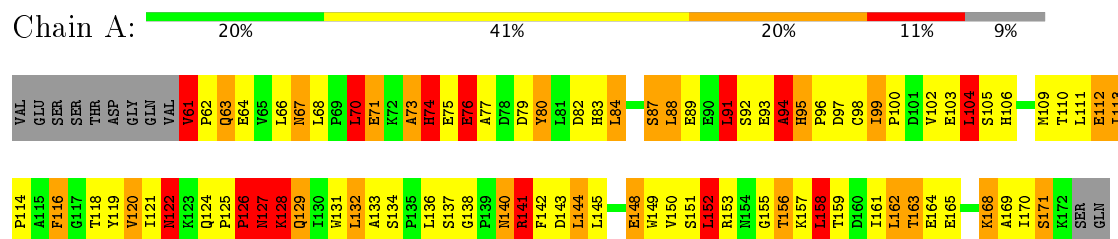
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	73	ALA	TYR	ENGINEERED	UNP Q07540

### 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: Frataxin homolog, mitochondrial



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 21 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	121.22Å 121.22Å 121.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.74 – 3.01 28.57 – 3.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.74-3.01) 100.0 (28.57-3.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.30 (at 3.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.286 , 0.305 0.238 , 0.255	Depositor DCC
$R_{free}$ test set	300 reflections (5.01%)	DCC
Wilson B-factor (Å <sup>2</sup> )	82.4	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 70.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.045 for -l,-k,-h	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	883	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.09% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	2.02	26/903 (2.9%)	1.75	19/1233 (1.5%)

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	A	0	4

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	128	LYS	N-CA	12.85	1.72	1.46
1	A	148	GLU	CG-CD	12.16	1.70	1.51
1	A	128	LYS	CD-CE	11.84	1.80	1.51
1	A	129	GLN	N-CA	10.62	1.67	1.46
1	A	148	GLU	CB-CG	10.44	1.72	1.52
1	A	75	GLU	CD-OE2	9.67	1.36	1.25
1	A	127	ASN	CB-CG	9.50	1.73	1.51
1	A	128	LYS	CB-CG	8.60	1.75	1.52
1	A	128	LYS	CE-NZ	8.38	1.70	1.49
1	A	148	GLU	CD-OE2	7.63	1.34	1.25
1	A	112	GLU	CG-CD	6.82	1.62	1.51
1	A	61	VAL	CA-CB	6.60	1.68	1.54
1	A	61	VAL	CB-CG2	6.46	1.66	1.52
1	A	112	GLU	CB-CG	6.30	1.64	1.52
1	A	128	LYS	CG-CD	6.04	1.73	1.52
1	A	75	GLU	CB-CG	-5.84	1.41	1.52
1	A	74	HIS	N-CA	5.83	1.58	1.46
1	A	76	GLU	CG-CD	5.60	1.60	1.51
1	A	127	ASN	CG-ND2	5.57	1.46	1.32
1	A	80	TYR	CD2-CE2	5.53	1.47	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	148	GLU	CD-OE1	5.43	1.31	1.25
1	A	120	VAL	CA-CB	-5.20	1.43	1.54
1	A	94	ALA	CA-CB	5.17	1.63	1.52
1	A	61	VAL	CA-C	5.09	1.66	1.52
1	A	129	GLN	CA-C	5.04	1.66	1.52
1	A	71	GLU	CB-CG	-5.01	1.42	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	129	GLN	N-CA-C	10.40	139.08	111.00
1	A	74	HIS	CB-CA-C	-8.82	92.75	110.40
1	A	152	LEU	CA-CB-CG	-8.40	95.97	115.30
1	A	128	LYS	CD-CE-NZ	8.39	131.01	111.70
1	A	122	ASN	CB-CA-C	-7.53	95.33	110.40
1	A	152	LEU	CB-CG-CD1	6.87	122.68	111.00
1	A	61	VAL	CB-CA-C	6.70	124.12	111.40
1	A	162	LEU	CB-CG-CD1	-6.58	99.82	111.00
1	A	145	LEU	C-N-CA	-6.55	105.32	121.70
1	A	75	GLU	OE1-CD-OE2	6.48	131.08	123.30
1	A	91	LEU	CA-CB-CG	6.45	130.13	115.30
1	A	70	LEU	CB-CG-CD1	6.11	121.39	111.00
1	A	79	ASP	CB-CG-OD2	-6.04	112.87	118.30
1	A	73	ALA	CA-C-N	5.68	129.69	117.20
1	A	75	GLU	CG-CD-OE1	-5.61	107.08	118.30
1	A	128	LYS	CA-C-N	5.50	129.30	117.20
1	A	104	LEU	CB-CG-CD1	5.49	120.34	111.00
1	A	158	LEU	CB-CG-CD2	-5.43	101.77	111.00
1	A	141	ARG	NE-CZ-NH2	5.40	123.00	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	122	ASN	Peptide
1	A	126	PRO	Peptide
1	A	127	ASN	Peptide
1	A	61	VAL	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	883	0	865	160	3
All	All	883	0	865	160	3

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 92.

All (160) 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:128:LYS:CG	1:A:128:LYS:CB	1.75	1.57
1:A:129:GLN:CA	1:A:129:GLN:N	1.67	1.56
1:A:128:LYS:CD	1:A:128:LYS:CE	1.80	1.55
1:A:128:LYS:NZ	1:A:128:LYS:CE	1.70	1.52
1:A:128:LYS:CA	1:A:128:LYS:N	1.72	1.50
1:A:112:GLU:OE2	1:A:118:THR:HG21	1.43	1.19
1:A:92:SER:HA	1:A:99:ILE:CD1	1.74	1.18
1:A:73:ALA:O	1:A:74:HIS:HB2	1.33	1.12
1:A:92:SER:CA	1:A:99:ILE:HD11	1.82	1.09
1:A:113:ILE:HG22	1:A:113:ILE:O	1.54	1.05
1:A:95:HIS:CE1	1:A:97:ASP:HA	1.91	1.05
1:A:95:HIS:O	1:A:95:HIS:CG	2.11	1.03
1:A:132:LEU:HD22	1:A:133:ALA:N	1.75	1.01
1:A:95:HIS:O	1:A:95:HIS:CD2	2.15	0.99
1:A:129:GLN:HA	1:A:129:GLN:N	1.74	0.98
1:A:95:HIS:HB2	1:A:99:ILE:CG1	1.92	0.98
1:A:95:HIS:HB2	1:A:99:ILE:HG13	1.46	0.96
1:A:73:ALA:O	1:A:74:HIS:CB	2.11	0.96
1:A:76:GLU:N	1:A:76:GLU:OE2	2.03	0.91
1:A:95:HIS:CE1	1:A:97:ASP:CA	2.53	0.91
1:A:128:LYS:HD3	1:A:152:LEU:CB	2.01	0.90
1:A:96:PRO:HD2	1:A:98:CYS:SG	2.12	0.90
1:A:63:GLN:HA	1:A:66:LEU:CD2	2.04	0.88
1:A:126:PRO:C	1:A:128:LYS:H	1.78	0.86
1:A:126:PRO:C	1:A:128:LYS:N	2.34	0.80
1:A:92:SER:HA	1:A:99:ILE:HD11	0.86	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:SER:HB3	1:A:165:GLU:OE2	1.83	0.78
1:A:127:ASN:O	1:A:128:LYS:HG3	1.84	0.78
1:A:66:LEU:H	1:A:66:LEU:HD23	1.47	0.78
1:A:128:LYS:CB	1:A:152:LEU:HD13	2.15	0.76
1:A:112:GLU:HA	1:A:118:THR:HG22	1.67	0.75
1:A:140:ASN:HD22	1:A:141:ARG:H	1.31	0.75
1:A:93:GLU:C	1:A:95:HIS:H	1.89	0.75
1:A:116:PHE:CD2	1:A:116:PHE:N	2.54	0.75
1:A:128:LYS:HD2	1:A:141:ARG:HG2	1.68	0.74
1:A:95:HIS:HB2	1:A:99:ILE:HG12	1.68	0.74
1:A:168:LYS:CD	1:A:168:LYS:N	2.51	0.73
1:A:132:LEU:CD2	1:A:133:ALA:N	2.51	0.72
1:A:132:LEU:HD22	1:A:132:LEU:C	2.10	0.72
1:A:99:ILE:O	1:A:99:ILE:HG13	1.89	0.71
1:A:120:VAL:HG12	1:A:121:ILE:N	2.06	0.71
1:A:84:LEU:HD13	1:A:109:MET:HE1	1.71	0.71
1:A:111:LEU:HB2	1:A:119:TYR:HB2	1.73	0.70
1:A:113:ILE:CG2	1:A:113:ILE:O	2.33	0.70
1:A:67:ASN:H	1:A:67:ASN:HD22	1.38	0.70
1:A:93:GLU:O	1:A:95:HIS:N	2.23	0.70
1:A:128:LYS:HD3	1:A:152:LEU:HB3	1.73	0.70
1:A:151:SER:HB3	1:A:156:THR:H	1.57	0.69
1:A:129:GLN:HG3	1:A:129:GLN:O	1.91	0.69
1:A:88:LEU:HB3	1:A:102:VAL:HG11	1.74	0.69
1:A:128:LYS:CG	1:A:152:LEU:HD13	2.22	0.69
1:A:158:LEU:HG	1:A:162:LEU:HD11	1.74	0.68
1:A:128:LYS:HD2	1:A:141:ARG:CG	2.24	0.68
1:A:67:ASN:N	1:A:67:ASN:HD22	1.90	0.67
1:A:84:LEU:C	1:A:84:LEU:HD22	2.15	0.66
1:A:127:ASN:O	1:A:128:LYS:CG	2.44	0.66
1:A:168:LYS:HD2	1:A:168:LYS:N	2.09	0.65
1:A:63:GLN:HA	1:A:66:LEU:HD21	1.77	0.65
1:A:151:SER:HB3	1:A:156:THR:N	2.12	0.65
1:A:95:HIS:CE1	1:A:97:ASP:C	2.71	0.64
1:A:140:ASN:HD22	1:A:141:ARG:N	1.96	0.64
1:A:99:ILE:O	1:A:99:ILE:CG1	2.45	0.64
1:A:93:GLU:OE1	1:A:93:GLU:N	2.31	0.64
1:A:102:VAL:HG12	1:A:111:LEU:HD21	1.79	0.63
1:A:89:GLU:HB3	1:A:102:VAL:CG2	2.28	0.63
1:A:84:LEU:HD13	1:A:109:MET:CE	2.29	0.63
1:A:132:LEU:CD2	1:A:132:LEU:C	2.67	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:LEU:O	1:A:162:LEU:HD12	1.99	0.62
1:A:170:ILE:HD12	1:A:170:ILE:H	1.65	0.61
1:A:89:GLU:HA	1:A:92:SER:HB2	1.83	0.61
1:A:93:GLU:C	1:A:95:HIS:N	2.54	0.60
1:A:128:LYS:HB3	1:A:152:LEU:HD13	1.82	0.60
1:A:66:LEU:H	1:A:66:LEU:CD2	2.14	0.60
1:A:83:HIS:CD2	1:A:83:HIS:N	2.69	0.59
1:A:67:ASN:ND2	1:A:67:ASN:H	2.00	0.59
1:A:84:LEU:O	1:A:84:LEU:HD22	2.02	0.58
1:A:116:PHE:CE1	1:A:136:LEU:HD22	2.39	0.57
1:A:120:VAL:CG1	1:A:121:ILE:N	2.68	0.56
1:A:122:ASN:HB2	1:A:131:TRP:HB3	1.86	0.56
1:A:80:TYR:CE2	1:A:158:LEU:HD21	2.40	0.56
1:A:159:THR:O	1:A:163:THR:HB	2.06	0.56
1:A:112:GLU:OE2	1:A:118:THR:CG2	2.36	0.56
1:A:128:LYS:HD2	1:A:141:ARG:HD3	1.88	0.56
1:A:66:LEU:N	1:A:66:LEU:HD23	2.21	0.56
1:A:63:GLN:HA	1:A:66:LEU:HD22	1.85	0.55
1:A:61:VAL:HG12	1:A:62:PRO:HD3	1.88	0.55
1:A:96:PRO:CD	1:A:98:CYS:SG	2.93	0.54
1:A:168:LYS:HD3	1:A:168:LYS:N	2.23	0.54
1:A:127:ASN:C	1:A:128:LYS:HG3	2.29	0.54
1:A:128:LYS:N	1:A:128:LYS:CB	2.66	0.53
1:A:128:LYS:HD3	1:A:152:LEU:CG	2.39	0.53
1:A:95:HIS:HE1	1:A:97:ASP:HA	1.63	0.53
1:A:83:HIS:H	1:A:83:HIS:CD2	2.28	0.52
1:A:82:ASP:C	1:A:84:LEU:H	2.13	0.52
1:A:150:VAL:HG13	1:A:151:SER:N	2.23	0.52
1:A:67:ASN:N	1:A:67:ASN:ND2	2.56	0.51
1:A:124:GLN:HG3	1:A:125:PRO:HD2	1.92	0.51
1:A:116:PHE:CE2	1:A:169:ALA:HB2	2.46	0.51
1:A:120:VAL:C	1:A:121:ILE:HG13	2.32	0.50
1:A:128:LYS:CD	1:A:141:ARG:HD3	2.41	0.50
1:A:161:ILE:O	1:A:165:GLU:HG3	2.10	0.50
1:A:128:LYS:C	1:A:129:GLN:HA	2.31	0.50
1:A:102:VAL:HG12	1:A:111:LEU:CD2	2.42	0.49
1:A:170:ILE:O	1:A:171:SER:C	2.51	0.49
1:A:134:SER:O	1:A:138:GLY:N	2.45	0.49
1:A:64:GLU:HG2	1:A:64:GLU:O	2.12	0.49
1:A:128:LYS:CB	1:A:152:LEU:CD1	2.90	0.49
1:A:168:LYS:HD3	1:A:168:LYS:H	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:LEU:O	1:A:91:LEU:HB2	2.13	0.49
1:A:83:HIS:O	1:A:87:SER:HB3	2.13	0.48
1:A:116:PHE:HE2	1:A:169:ALA:HB2	1.78	0.48
1:A:93:GLU:HB2	1:A:94:ALA:H	1.33	0.48
1:A:109:MET:HG2	1:A:110:THR:N	2.27	0.48
1:A:148:GLU:HG3	1:A:157:LYS:NZ	2.29	0.48
1:A:122:ASN:HD22	1:A:124:GLN:HE22	1.60	0.48
1:A:128:LYS:HB3	1:A:152:LEU:CD1	2.43	0.48
1:A:128:LYS:CD	1:A:152:LEU:HD13	2.44	0.48
1:A:116:PHE:CD1	1:A:136:LEU:HD22	2.49	0.48
1:A:80:TYR:CZ	1:A:158:LEU:HD21	2.48	0.47
1:A:88:LEU:HD12	1:A:88:LEU:HA	1.38	0.47
1:A:144:LEU:HD13	1:A:149:TRP:CE2	2.49	0.47
1:A:126:PRO:HA	1:A:128:LYS:CA	2.44	0.47
1:A:89:GLU:HB3	1:A:102:VAL:HG21	1.97	0.46
1:A:82:ASP:C	1:A:84:LEU:N	2.68	0.46
1:A:164:GLU:O	1:A:165:GLU:C	2.53	0.46
1:A:91:LEU:C	1:A:93:GLU:H	2.19	0.46
1:A:150:VAL:HG13	1:A:155:GLY:HA2	1.97	0.46
1:A:148:GLU:HG3	1:A:157:LYS:HZ1	1.81	0.46
1:A:141:ARG:HH11	1:A:141:ARG:HD2	1.64	0.45
1:A:112:GLU:HA	1:A:118:THR:CG2	2.41	0.45
1:A:129:GLN:O	1:A:129:GLN:CG	2.63	0.45
1:A:93:GLU:O	1:A:95:HIS:O	2.35	0.45
1:A:128:LYS:HD2	1:A:141:ARG:CD	2.45	0.44
1:A:129:GLN:CB	1:A:129:GLN:N	2.68	0.43
1:A:142:PHE:HA	1:A:150:VAL:O	2.18	0.43
1:A:111:LEU:HA	1:A:111:LEU:HD23	1.43	0.43
1:A:125:PRO:HA	1:A:126:PRO:HD3	1.68	0.43
1:A:84:LEU:CD1	1:A:109:MET:CE	2.97	0.43
1:A:151:SER:N	1:A:156:THR:O	2.52	0.43
1:A:89:GLU:HB3	1:A:102:VAL:HG22	2.02	0.42
1:A:93:GLU:N	1:A:93:GLU:CD	2.71	0.42
1:A:116:PHE:H	1:A:116:PHE:HD2	1.66	0.42
1:A:122:ASN:HD22	1:A:124:GLN:NE2	2.17	0.42
1:A:140:ASN:HB3	1:A:142:PHE:CE2	2.54	0.42
1:A:70:LEU:HD12	1:A:70:LEU:HA	1.91	0.42
1:A:119:TYR:OH	1:A:165:GLU:HB3	2.20	0.41
1:A:170:ILE:HD12	1:A:170:ILE:N	2.33	0.41
1:A:82:ASP:O	1:A:84:LEU:N	2.53	0.41
1:A:104:LEU:CD2	1:A:104:LEU:C	2.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:162:LEU:HD12	1:A:162:LEU:H	1.85	0.41
1:A:68:LEU:HB2	1:A:70:LEU:HD13	2.02	0.41
1:A:100:PRO:HD2	1:A:114:PRO:HD3	2.02	0.41
1:A:103:GLU:HB3	1:A:110:THR:OG1	2.21	0.41
1:A:89:GLU:C	1:A:91:LEU:H	2.23	0.41
1:A:113:ILE:HG22	1:A:116:PHE:HD2	1.85	0.41
1:A:128:LYS:CD	1:A:152:LEU:CB	2.87	0.41
1:A:113:ILE:HA	1:A:114:PRO:HD3	1.87	0.40
1:A:76:GLU:N	1:A:76:GLU:CD	2.73	0.40
1:A:76:GLU:O	1:A:77:ALA:C	2.59	0.40
1:A:66:LEU:CD2	1:A:66:LEU:N	2.82	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:ASN:CA	1:A:129:GLN:NE2[7_454]	1.85	0.35
1:A:127:ASN:CB	1:A:129:GLN:NE2[7_454]	1.89	0.31
1:A:127:ASN:ND2	1:A:129:GLN:CG[7_454]	2.12	0.08

## 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	110/123 (89%)	79 (72%)	25 (23%)	6 (6%)	<b>2</b> <b>12</b>

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	94	ALA
1	A	126	PRO
1	A	105	SER

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Mol	Chain	Res	Type
1	A	127	ASN
1	A	128	LYS
1	A	74	HIS

### 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	101/111 (91%)	72 (71%)	29 (29%)	<b>0</b> <b>2</b>

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	GLN
1	A	67	ASN
1	A	70	LEU
1	A	71	GLU
1	A	76	GLU
1	A	84	LEU
1	A	87	SER
1	A	88	LEU
1	A	91	LEU
1	A	95	HIS
1	A	99	ILE
1	A	104	LEU
1	A	106	HIS
1	A	113	ILE
1	A	116	PHE
1	A	127	ASN
1	A	128	LYS
1	A	132	LEU
1	A	140	ASN
1	A	141	ARG
1	A	143	ASP
1	A	144	LEU
1	A	152	LEU

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Mol	Chain	Res	Type
1	A	153	ARG
1	A	156	THR
1	A	158	LEU
1	A	163	THR
1	A	168	LYS
1	A	171	SER

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	63	GLN
1	A	67	ASN
1	A	83	HIS
1	A	95	HIS
1	A	124	GLN
1	A	127	ASN

### 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 [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	112/123 (91%)	-0.45	0 100 100	43, 71, 100, 100	1 (0%)

There are no RSRZ outliers to report.

### 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.