



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 11, 2021 – 04:33 PM EDT

PDB ID : 2O94
Title : The 97H/F mutant Structure of a glutamine-rich domain from histone deacetylase 4
Authors : Guo, L.; Han, A.; Bates, D.L.; Chen, L.
Deposited on : 2006-12-13
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

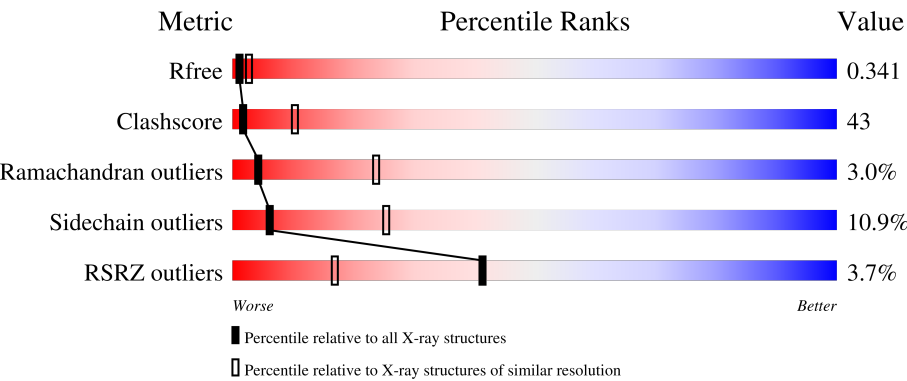
MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.23.2
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.23.2

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 3.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	112	
1	B	112	
1	C	112	
1	D	112	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2328 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 Histone deacetylase 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	68	Total	C	N	O	S	0	0	0
			582	361	111	108	2			
1	B	68	Total	C	N	O	S	0	0	0
			582	361	111	108	2			
1	C	68	Total	C	N	O	S	0	0	0
			582	361	111	108	2			
1	D	68	Total	C	N	O	S	0	0	0
			582	361	111	108	2			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	42	GLY	-	cloning artifact	UNP P56524
A	43	SER	-	cloning artifact	UNP P56524
A	44	SER	-	cloning artifact	UNP P56524
A	45	HIS	-	expression tag	UNP P56524
A	46	HIS	-	expression tag	UNP P56524
A	47	HIS	-	expression tag	UNP P56524
A	48	HIS	-	expression tag	UNP P56524
A	49	HIS	-	expression tag	UNP P56524
A	50	HIS	-	expression tag	UNP P56524
A	51	SER	-	cloning artifact	UNP P56524
A	52	SER	-	cloning artifact	UNP P56524
A	53	GLY	-	cloning artifact	UNP P56524
A	54	LEU	-	cloning artifact	UNP P56524
A	55	VAL	-	cloning artifact	UNP P56524
A	56	PRO	-	cloning artifact	UNP P56524
A	57	ARG	-	cloning artifact	UNP P56524
A	58	GLY	-	cloning artifact	UNP P56524
A	59	SER	-	cloning artifact	UNP P56524
A	60	HIS	-	cloning artifact	UNP P56524
A	61	MET	-	cloning artifact	UNP P56524
A	97	PHE	HIS	engineered mutation	UNP P56524

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Chain	Residue	Modelled	Actual	Comment	Reference
B	42	GLY	-	cloning artifact	UNP P56524
B	43	SER	-	cloning artifact	UNP P56524
B	44	SER	-	cloning artifact	UNP P56524
B	45	HIS	-	expression tag	UNP P56524
B	46	HIS	-	expression tag	UNP P56524
B	47	HIS	-	expression tag	UNP P56524
B	48	HIS	-	expression tag	UNP P56524
B	49	HIS	-	expression tag	UNP P56524
B	50	HIS	-	expression tag	UNP P56524
B	51	SER	-	cloning artifact	UNP P56524
B	52	SER	-	cloning artifact	UNP P56524
B	53	GLY	-	cloning artifact	UNP P56524
B	54	LEU	-	cloning artifact	UNP P56524
B	55	VAL	-	cloning artifact	UNP P56524
B	56	PRO	-	cloning artifact	UNP P56524
B	57	ARG	-	cloning artifact	UNP P56524
B	58	GLY	-	cloning artifact	UNP P56524
B	59	SER	-	cloning artifact	UNP P56524
B	60	HIS	-	cloning artifact	UNP P56524
B	61	MET	-	cloning artifact	UNP P56524
B	97	PHE	HIS	engineered mutation	UNP P56524
C	42	GLY	-	cloning artifact	UNP P56524
C	43	SER	-	cloning artifact	UNP P56524
C	44	SER	-	cloning artifact	UNP P56524
C	45	HIS	-	expression tag	UNP P56524
C	46	HIS	-	expression tag	UNP P56524
C	47	HIS	-	expression tag	UNP P56524
C	48	HIS	-	expression tag	UNP P56524
C	49	HIS	-	expression tag	UNP P56524
C	50	HIS	-	expression tag	UNP P56524
C	51	SER	-	cloning artifact	UNP P56524
C	52	SER	-	cloning artifact	UNP P56524
C	53	GLY	-	cloning artifact	UNP P56524
C	54	LEU	-	cloning artifact	UNP P56524
C	55	VAL	-	cloning artifact	UNP P56524
C	56	PRO	-	cloning artifact	UNP P56524
C	57	ARG	-	cloning artifact	UNP P56524
C	58	GLY	-	cloning artifact	UNP P56524
C	59	SER	-	cloning artifact	UNP P56524
C	60	HIS	-	cloning artifact	UNP P56524
C	61	MET	-	cloning artifact	UNP P56524
C	97	PHE	HIS	engineered mutation	UNP P56524

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Chain	Residue	Modelled	Actual	Comment	Reference
D	42	GLY	-	cloning artifact	UNP P56524
D	43	SER	-	cloning artifact	UNP P56524
D	44	SER	-	cloning artifact	UNP P56524
D	45	HIS	-	expression tag	UNP P56524
D	46	HIS	-	expression tag	UNP P56524
D	47	HIS	-	expression tag	UNP P56524
D	48	HIS	-	expression tag	UNP P56524
D	49	HIS	-	expression tag	UNP P56524
D	50	HIS	-	expression tag	UNP P56524
D	51	SER	-	cloning artifact	UNP P56524
D	52	SER	-	cloning artifact	UNP P56524
D	53	GLY	-	cloning artifact	UNP P56524
D	54	LEU	-	cloning artifact	UNP P56524
D	55	VAL	-	cloning artifact	UNP P56524
D	56	PRO	-	cloning artifact	UNP P56524
D	57	ARG	-	cloning artifact	UNP P56524
D	58	GLY	-	cloning artifact	UNP P56524
D	59	SER	-	cloning artifact	UNP P56524
D	60	HIS	-	cloning artifact	UNP P56524
D	61	MET	-	cloning artifact	UNP P56524
D	97	PHE	HIS	engineered mutation	UNP P56524

- Molecule 1: Histone deacetylase 4



I112	K113	Q114	Q115	Q116	E117	M118	L119	A120	M121	K122	H123	Q124	E125	E126	L127	L128	E129	HIS	GLN	ARG	LYS	LEU	GLU	ARG	HIS	ARG	GLN	GLU	GLN	GLN	GLU	LEU	GLU	LYS	GLN	HIS	ARG	GLU	GLN	LYS	LEU	GLN
------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	189.88Å 61.02Å 60.78Å 90.00° 108.61° 90.00°	Depositor
Resolution (Å)	30.00 – 3.00 29.99 – 2.94	Depositor EDS
% Data completeness (in resolution range)	75.1 (30.00-3.00) 84.2 (29.99-2.94)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 2.95Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.305 , 0.327 0.313 , 0.341	Depositor DCC
R_{free} test set	1348 reflections (9.79%)	wwPDB-VP
Wilson B-factor (Å ²)	81.3	Xtriage
Anisotropy	0.460	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 85.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	0.460 for -h-k-l,l,k 0.469 for -h+k-l,-l,-k 0.468 for -h-2*l,-k,l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	2328	wwPDB-VP
Average B, all atoms (Å ²)	108.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.99% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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	0.43	0/588	0.66	0/784
1	B	0.45	0/588	0.66	0/784
1	C	0.42	0/588	0.66	0/784
1	D	0.46	0/588	0.64	0/784
All	All	0.44	0/2352	0.65	0/3136

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	582	0	586	67	0
1	B	582	0	586	59	0
1	C	582	0	586	59	0
1	D	582	0	586	73	0
All	All	2328	0	2344	203	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 43.

All (203) 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:121:MET:HA	1:D:124:GLN:HG2	1.28	1.11
1:A:121:MET:HA	1:A:124:GLN:HG2	1.41	1.02
1:B:121:MET:HA	1:B:124:GLN:HG2	1.51	0.93
1:A:97:PHE:HE1	1:B:97:PHE:HE1	1.20	0.89
1:A:75:LEU:HD13	1:C:111:HIS:NE2	1.91	0.86
1:D:123:HIS:HA	1:D:126:GLU:HB2	1.57	0.85
1:B:123:HIS:HA	1:B:126:GLU:HB2	1.58	0.85
1:B:111:HIS:CE1	1:D:75:LEU:HD13	2.11	0.85
1:B:63:GLU:HB3	1:B:64:PRO:HD3	1.57	0.84
1:C:97:PHE:HE1	1:D:97:PHE:HE1	1.21	0.82
1:D:62:ALA:HB3	1:D:66:LEU:HD13	1.60	0.82
1:D:114:GLN:HA	1:D:114:GLN:HE21	1.44	0.82
1:A:67:ARG:HH11	1:A:67:ARG:HB2	1.43	0.81
1:B:67:ARG:HG3	1:B:67:ARG:HH11	1.45	0.81
1:A:97:PHE:HE1	1:B:97:PHE:CE1	1.97	0.81
1:C:121:MET:N	1:C:124:GLN:HE21	1.80	0.80
1:D:68:GLU:HG2	1:D:72:GLN:NE2	1.98	0.79
1:C:123:HIS:HA	1:C:126:GLU:HB2	1.67	0.76
1:C:97:PHE:CE1	1:D:97:PHE:HE1	2.04	0.76
1:A:111:HIS:NE2	1:C:75:LEU:HB2	2.02	0.75
1:C:97:PHE:HE1	1:D:97:PHE:CE1	2.04	0.74
1:A:112:ILE:HG12	1:B:112:ILE:HG12	1.69	0.73
1:A:97:PHE:CE1	1:B:97:PHE:HE1	2.06	0.73
1:D:123:HIS:HA	1:D:126:GLU:CB	2.19	0.71
1:B:75:LEU:HB2	1:D:111:HIS:NE2	2.05	0.71
1:A:122:LYS:O	1:A:122:LYS:HD3	1.91	0.70
1:D:117:GLU:O	1:D:120:ALA:HB3	1.92	0.70
1:C:121:MET:HA	1:C:124:GLN:HG2	1.74	0.69
1:A:75:LEU:HD13	1:C:111:HIS:CD2	2.29	0.67
1:B:117:GLU:HA	1:B:120:ALA:HB3	1.77	0.66
1:C:123:HIS:HA	1:C:126:GLU:CB	2.26	0.65
1:A:123:HIS:HA	1:A:126:GLU:HB3	1.78	0.65
1:A:118:MET:SD	1:A:119:LEU:HD23	2.38	0.64
1:B:75:LEU:HD13	1:D:111:HIS:CE1	2.34	0.63
1:A:112:ILE:HG12	1:B:112:ILE:CG1	2.29	0.62
1:A:75:LEU:HD13	1:C:111:HIS:CE1	2.34	0.62
1:A:111:HIS:NE2	1:C:75:LEU:HD13	2.15	0.62
1:D:68:GLU:HG2	1:D:72:GLN:HE22	1.66	0.60
1:B:121:MET:HA	1:B:124:GLN:CG	2.29	0.60
1:C:74:GLU:O	1:C:77:ALA:HB3	2.02	0.60
1:A:74:GLU:O	1:A:77:ALA:HB3	2.02	0.59
1:C:83:GLN:O	1:C:83:GLN:HG2	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:117:GLU:O	1:C:120:ALA:HB3	2.03	0.59
1:B:118:MET:HE2	1:D:68:GLU:HA	1.85	0.58
1:C:63:GLU:HB3	1:C:64:PRO:CD	2.32	0.58
1:D:63:GLU:HB3	1:D:64:PRO:CD	2.33	0.58
1:B:67:ARG:HH11	1:B:67:ARG:CG	2.15	0.58
1:B:82:GLN:HG2	1:D:104:HIS:CD2	2.37	0.58
1:A:97:PHE:CE1	1:B:97:PHE:CE1	2.84	0.57
1:A:113:LYS:O	1:A:116:GLN:HB3	2.04	0.57
1:A:121:MET:CA	1:A:124:GLN:HE21	2.18	0.57
1:C:112:ILE:CG1	1:D:112:ILE:HG12	2.35	0.57
1:C:112:ILE:HG12	1:D:112:ILE:HG12	1.87	0.56
1:D:118:MET:SD	1:D:119:LEU:N	2.79	0.56
1:D:125:GLN:O	1:D:125:GLN:HG2	2.06	0.56
1:B:75:LEU:HD13	1:D:111:HIS:NE2	2.21	0.56
1:C:75:LEU:O	1:C:78:LEU:N	2.34	0.56
1:D:94:GLN:O	1:D:97:PHE:HB3	2.06	0.56
1:A:118:MET:SD	1:A:119:LEU:N	2.79	0.56
1:D:113:LYS:O	1:D:116:GLN:HB3	2.05	0.56
1:D:83:GLN:C	1:D:83:GLN:HE21	2.10	0.55
1:A:111:HIS:CD2	1:C:75:LEU:HD13	2.42	0.55
1:B:115:GLN:HA	1:D:71:LEU:CD1	2.36	0.55
1:C:112:ILE:HD11	1:D:112:ILE:HG12	1.89	0.55
1:D:74:GLU:O	1:D:77:ALA:HB3	2.07	0.55
1:B:117:GLU:O	1:B:120:ALA:HB3	2.08	0.54
1:D:67:ARG:HG2	1:D:67:ARG:HH11	1.73	0.54
1:A:94:GLN:O	1:A:97:PHE:HB3	2.08	0.54
1:B:68:GLU:HG3	1:D:118:MET:CE	2.38	0.53
1:B:76:LEU:HD23	1:B:76:LEU:O	2.08	0.53
1:C:114:GLN:C	1:C:116:GLN:H	2.10	0.53
1:A:68:GLU:HG2	1:A:72:GLN:OE1	2.07	0.53
1:C:97:PHE:CE1	1:D:97:PHE:CE1	2.87	0.53
1:A:117:GLU:O	1:A:120:ALA:HB3	2.09	0.53
1:B:74:GLU:O	1:B:77:ALA:HB3	2.09	0.52
1:B:82:GLN:HG2	1:D:104:HIS:CG	2.44	0.52
1:C:63:GLU:HB3	1:C:64:PRO:HD3	1.91	0.52
1:A:122:LYS:HD3	1:A:122:LYS:C	2.29	0.52
1:A:118:MET:HE2	1:C:68:GLU:HA	1.92	0.52
1:C:126:GLU:O	1:C:126:GLU:HG3	2.10	0.52
1:A:111:HIS:CE1	1:C:75:LEU:HD13	2.46	0.51
1:A:128:LEU:HD12	1:A:128:LEU:H	1.76	0.51
1:B:68:GLU:HG3	1:D:118:MET:HE2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:GLU:O	1:A:126:GLU:OE1	2.29	0.51
1:D:121:MET:CA	1:D:124:GLN:HG2	2.21	0.51
1:D:126:GLU:OE2	1:D:129:GLU:OE2	2.29	0.51
1:B:75:LEU:O	1:B:78:LEU:N	2.43	0.51
1:C:117:GLU:HA	1:C:120:ALA:HB3	1.93	0.50
1:A:63:GLU:HB3	1:A:64:PRO:CD	2.41	0.50
1:A:118:MET:HE1	1:C:68:GLU:HB2	1.94	0.50
1:B:126:GLU:OE2	1:B:126:GLU:HA	2.11	0.50
1:B:94:GLN:O	1:B:97:PHE:HB3	2.12	0.49
1:D:119:LEU:HA	1:D:122:LYS:HG2	1.94	0.49
1:C:64:PRO:HG2	1:C:65:ALA:H	1.77	0.49
1:D:121:MET:N	1:D:124:GLN:HE21	2.10	0.49
1:D:114:GLN:NE2	1:D:114:GLN:CA	2.76	0.49
1:B:63:GLU:CB	1:B:64:PRO:HD3	2.38	0.49
1:C:112:ILE:CD1	1:D:112:ILE:HG12	2.43	0.48
1:D:62:ALA:HB3	1:D:66:LEU:CD1	2.40	0.48
1:B:67:ARG:HG3	1:B:67:ARG:NH1	2.22	0.48
1:D:126:GLU:O	1:D:126:GLU:HG3	2.11	0.48
1:B:75:LEU:C	1:B:77:ALA:H	2.17	0.48
1:C:67:ARG:HD2	1:C:70:GLN:OE1	2.13	0.48
1:D:121:MET:C	1:D:123:HIS:H	2.16	0.48
1:A:83:GLN:O	1:A:83:GLN:NE2	2.47	0.48
1:A:112:ILE:HG12	1:B:112:ILE:CD1	2.44	0.48
1:A:64:PRO:C	1:A:66:LEU:H	2.17	0.47
1:A:84:ILE:O	1:A:85:GLN:C	2.52	0.47
1:D:114:GLN:HA	1:D:114:GLN:NE2	2.20	0.47
1:D:63:GLU:CB	1:D:64:PRO:CD	2.93	0.47
1:C:114:GLN:C	1:C:116:GLN:N	2.68	0.47
1:C:121:MET:CA	1:C:124:GLN:HE21	2.28	0.47
1:B:118:MET:SD	1:B:119:LEU:N	2.88	0.47
1:C:118:MET:SD	1:C:119:LEU:N	2.88	0.47
1:D:117:GLU:HA	1:D:120:ALA:HB3	1.97	0.47
1:B:75:LEU:C	1:B:77:ALA:N	2.68	0.47
1:B:123:HIS:HA	1:B:126:GLU:CB	2.38	0.47
1:C:121:MET:CA	1:C:124:GLN:HG2	2.43	0.47
1:D:114:GLN:HE21	1:D:114:GLN:CA	2.11	0.46
1:B:100:LEU:HD12	1:B:100:LEU:HA	1.65	0.46
1:D:67:ARG:HG2	1:D:67:ARG:NH1	2.30	0.46
1:A:117:GLU:HA	1:A:120:ALA:HB3	1.97	0.46
1:A:94:GLN:HG2	1:D:90:ILE:HG21	1.98	0.46
1:B:63:GLU:O	1:B:66:LEU:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:121:MET:HA	1:D:124:GLN:CG	2.21	0.46
1:A:83:GLN:HE21	1:A:83:GLN:C	2.20	0.46
1:B:98:GLU:O	1:B:102:ARG:HG3	2.16	0.46
1:B:117:GLU:HA	1:B:120:ALA:CB	2.45	0.46
1:D:83:GLN:C	1:D:83:GLN:NE2	2.69	0.46
1:B:116:GLN:O	1:B:120:ALA:N	2.47	0.45
1:A:68:GLU:CB	1:C:118:MET:HE1	2.46	0.45
1:A:90:ILE:HD13	1:A:90:ILE:HA	1.85	0.45
1:D:121:MET:CA	1:D:124:GLN:HE21	2.29	0.45
1:A:103:GLN:O	1:A:104:HIS:C	2.54	0.45
1:B:117:GLU:CA	1:B:120:ALA:HB3	2.46	0.45
1:B:63:GLU:HB3	1:B:64:PRO:CD	2.38	0.45
1:B:104:HIS:CD2	1:D:82:GLN:HG2	2.52	0.45
1:D:100:LEU:HD12	1:D:100:LEU:HA	1.73	0.45
1:A:93:PHE:HE2	1:D:90:ILE:HD11	1.81	0.45
1:C:64:PRO:C	1:C:66:LEU:H	2.20	0.45
1:C:64:PRO:C	1:C:66:LEU:N	2.70	0.45
1:C:98:GLU:O	1:C:102:ARG:HG3	2.17	0.45
1:C:103:GLN:O	1:C:104:HIS:C	2.55	0.45
1:B:121:MET:CA	1:B:124:GLN:HG2	2.34	0.45
1:C:118:MET:SD	1:C:119:LEU:HD23	2.57	0.45
1:C:101:SER:O	1:C:105:GLU:HB2	2.17	0.45
1:A:71:LEU:CD1	1:C:115:GLN:HA	2.47	0.44
1:A:84:ILE:O	1:A:87:GLN:N	2.50	0.44
1:A:90:ILE:HD11	1:D:93:PHE:HE2	1.81	0.44
1:C:94:GLN:O	1:C:97:PHE:HB3	2.17	0.44
1:A:125:GLN:HA	1:A:128:LEU:HD13	1.98	0.44
1:A:68:GLU:HA	1:C:118:MET:HE2	1.99	0.44
1:D:124:GLN:HG3	1:D:125:GLN:N	2.33	0.44
1:A:68:GLU:O	1:A:72:GLN:HG3	2.19	0.43
1:A:128:LEU:HD12	1:A:128:LEU:N	2.33	0.43
1:B:67:ARG:CG	1:B:67:ARG:NH1	2.75	0.43
1:D:101:SER:O	1:D:105:GLU:HB2	2.18	0.43
1:D:114:GLN:C	1:D:116:GLN:N	2.71	0.43
1:D:103:GLN:O	1:D:104:HIS:C	2.56	0.43
1:A:90:ILE:HG21	1:D:94:GLN:HG2	2.01	0.43
1:C:64:PRO:O	1:C:67:ARG:N	2.51	0.43
1:A:68:GLU:HA	1:C:118:MET:CE	2.49	0.43
1:A:83:GLN:NE2	1:A:83:GLN:C	2.72	0.43
1:A:100:LEU:HD12	1:A:100:LEU:HA	1.66	0.43
1:B:103:GLN:O	1:B:104:HIS:C	2.56	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:PHE:O	1:A:94:GLN:C	2.56	0.43
1:A:98:GLU:O	1:A:102:ARG:HG3	2.18	0.43
1:B:101:SER:O	1:B:105:GLU:HB2	2.19	0.43
1:D:84:ILE:O	1:D:85:GLN:C	2.57	0.43
1:C:128:LEU:HD12	1:C:128:LEU:N	2.34	0.43
1:B:83:GLN:O	1:B:83:GLN:NE2	2.52	0.42
1:A:107:GLN:O	1:A:110:GLU:HB3	2.19	0.42
1:A:112:ILE:HG12	1:B:112:ILE:HD11	2.01	0.42
1:A:119:LEU:HD11	1:B:116:GLN:HG3	2.01	0.42
1:B:111:HIS:ND1	1:D:75:LEU:HD13	2.31	0.42
1:D:64:PRO:C	1:D:66:LEU:N	2.73	0.42
1:A:64:PRO:C	1:A:66:LEU:N	2.73	0.42
1:D:114:GLN:C	1:D:116:GLN:H	2.23	0.42
1:B:93:PHE:HE2	1:C:90:ILE:HD11	1.84	0.42
1:A:68:GLU:HB2	1:C:118:MET:HE1	2.02	0.42
1:C:128:LEU:HD12	1:C:128:LEU:H	1.83	0.42
1:D:109:HIS:O	1:D:112:ILE:HB	2.20	0.42
1:D:110:GLU:O	1:D:114:GLN:HG2	2.20	0.42
1:A:95:ARG:C	1:A:97:PHE:N	2.74	0.41
1:A:120:ALA:O	1:A:124:GLN:N	2.50	0.41
1:A:123:HIS:HD2	1:B:123:HIS:CD2	2.38	0.41
1:C:100:LEU:HD12	1:C:100:LEU:HA	1.65	0.41
1:D:119:LEU:HA	1:D:119:LEU:HD23	1.91	0.41
1:C:117:GLU:HA	1:C:120:ALA:CB	2.50	0.41
1:C:123:HIS:HD2	1:D:123:HIS:O	2.03	0.41
1:C:93:PHE:O	1:C:94:GLN:C	2.59	0.41
1:A:104:HIS:CD2	1:C:82:GLN:HG2	2.55	0.41
1:B:90:ILE:HG21	1:C:94:GLN:HG2	2.03	0.41
1:A:112:ILE:CG1	1:B:112:ILE:HG12	2.43	0.41
1:A:118:MET:SD	1:A:118:MET:C	2.99	0.41
1:B:114:GLN:C	1:B:116:GLN:N	2.72	0.41
1:A:121:MET:CB	1:A:124:GLN:HE21	2.34	0.41
1:D:84:ILE:O	1:D:87:GLN:N	2.54	0.40
1:B:68:GLU:HA	1:D:118:MET:HE2	2.03	0.40
1:D:120:ALA:O	1:D:124:GLN:N	2.50	0.40
1:D:118:MET:C	1:D:120:ALA:N	2.73	0.40
1:C:116:GLN:HG3	1:D:119:LEU:HD11	2.03	0.40
1:B:64:PRO:C	1:B:66:LEU:N	2.74	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	66/112 (59%)	52 (79%)	12 (18%)	2 (3%)	4	24
1	B	66/112 (59%)	53 (80%)	11 (17%)	2 (3%)	4	24
1	C	66/112 (59%)	52 (79%)	13 (20%)	1 (2%)	10	42
1	D	66/112 (59%)	50 (76%)	13 (20%)	3 (4%)	2	14
All	All	264/448 (59%)	207 (78%)	49 (19%)	8 (3%)	4	24

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	120	ALA
1	D	63	GLU
1	A	63	GLU
1	B	63	GLU
1	C	63	GLU
1	D	123	HIS
1	A	118	MET
1	D	120	ALA

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	62/103 (60%)	54 (87%)	8 (13%)	4	19
1	B	62/103 (60%)	58 (94%)	4 (6%)	17	50

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	62/103 (60%)	55 (89%)	7 (11%)	6	24
1	D	62/103 (60%)	54 (87%)	8 (13%)	4	19
All	All	248/412 (60%)	221 (89%)	27 (11%)	6	25

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	ARG
1	A	78	LEU
1	A	83	GLN
1	A	109	HIS
1	A	115	GLN
1	A	118	MET
1	A	122	LYS
1	A	126	GLU
1	B	83	GLN
1	B	109	HIS
1	B	123	HIS
1	B	127	LEU
1	C	63	GLU
1	C	67	ARG
1	C	74	GLU
1	C	109	HIS
1	C	118	MET
1	C	123	HIS
1	C	127	LEU
1	D	83	GLN
1	D	109	HIS
1	D	114	GLN
1	D	115	GLN
1	D	118	MET
1	D	123	HIS
1	D	125	GLN
1	D	127	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	83	GLN
1	A	85	GLN

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Mol	Chain	Res	Type
1	A	87	GLN
1	A	96	GLN
1	A	103	GLN
1	A	114	GLN
1	A	115	GLN
1	A	124	GLN
1	B	69	GLN
1	B	73	GLN
1	B	83	GLN
1	B	87	GLN
1	B	103	GLN
1	B	109	HIS
1	B	114	GLN
1	B	123	HIS
1	B	125	GLN
1	C	72	GLN
1	C	80	GLN
1	C	82	GLN
1	C	87	GLN
1	C	103	GLN
1	C	114	GLN
1	C	115	GLN
1	C	123	HIS
1	C	124	GLN
1	D	72	GLN
1	D	83	GLN
1	D	85	GLN
1	D	87	GLN
1	D	96	GLN
1	D	103	GLN
1	D	114	GLN
1	D	124	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th 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 ²)	Q<0.9
1	A	68/112 (60%)	0.16	0 100 100	56, 102, 162, 193	0
1	B	68/112 (60%)	0.33	4 (5%) 22 7	61, 104, 168, 189	0
1	C	68/112 (60%)	0.29	3 (4%) 34 13	56, 102, 166, 187	0
1	D	68/112 (60%)	0.10	3 (4%) 34 13	60, 106, 160, 186	0
All	All	272/448 (60%)	0.22	10 (3%) 41 17	56, 103, 169, 193	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	124	GLN	4.6
1	C	62	ALA	4.4
1	B	128	LEU	3.9
1	B	125	GLN	3.7
1	B	127	LEU	3.4
1	C	127	LEU	3.3
1	D	127	LEU	2.5
1	B	129	GLU	2.3
1	D	123	HIS	2.2
1	C	129	GLU	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 monosaccharides 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.