



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:08 PM UTC

PDB ID : 4LSD / pdb\_00004lsd  
Title : Myokine structure  
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Deposited on : 2013-07-22  
Resolution : 2.28 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

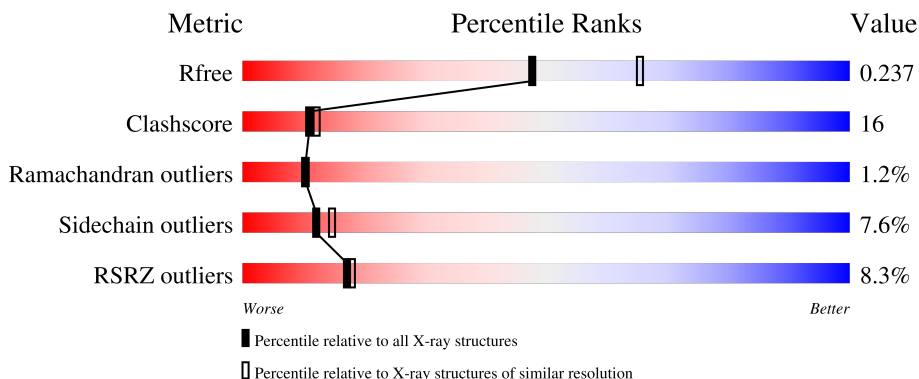
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.28 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.



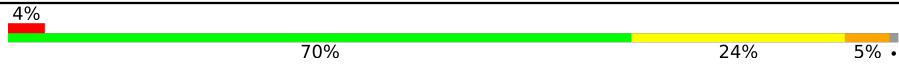

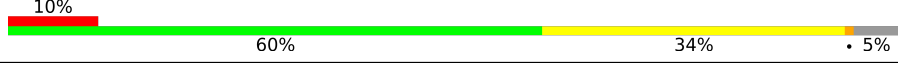
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	9078 (2.30-2.26)
Clashscore	190562	9802 (2.30-2.26)
Ramachandran outliers	187476	9690 (2.30-2.26)
Sidechain outliers	187428	9691 (2.30-2.26)
RSRZ outliers	180081	9085 (2.30-2.26)

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  $\geq 3$ , 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	99	
1	B	99	
1	C	99	
1	D	99	
1	E	99	

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Mol	Chain	Length	Quality of chain
1	F	99	
1	G	99	
1	H	99	

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 6225 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 Fibronectin type III domain-containing protein 5.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	98	767	483	132	149	1	2	0	0	0
1	B	94	740	467	128	143	1	1	0	0	0
1	C	95	748	472	129	144	1	2	0	0	0
1	D	94	740	467	128	143	1	1	0	0	0
1	E	94	740	467	128	143	1	1	0	0	0
1	F	98	772	486	133	151	1	1	0	0	0
1	G	94	740	467	128	143	1	1	0	0	0
1	H	94	740	467	128	143	1	1	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	29	MSE	-	initiating methionine	UNP Q8NAU1
B	29	MSE	-	initiating methionine	UNP Q8NAU1
C	29	MSE	-	initiating methionine	UNP Q8NAU1
D	29	MSE	-	initiating methionine	UNP Q8NAU1
E	29	MSE	-	initiating methionine	UNP Q8NAU1
F	29	MSE	-	initiating methionine	UNP Q8NAU1
G	29	MSE	-	initiating methionine	UNP Q8NAU1
H	29	MSE	-	initiating methionine	UNP Q8NAU1

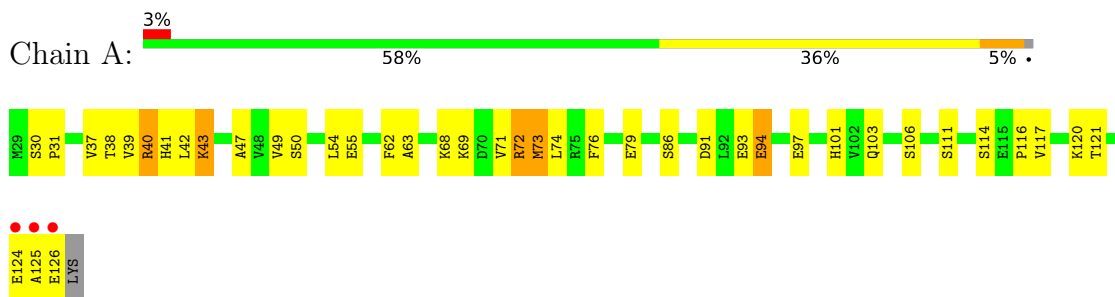
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	16	Total O 16 16	0	0
2	B	50	Total O 50 50	0	0
2	C	30	Total O 30 30	0	0
2	D	47	Total O 47 47	0	0
2	E	30	Total O 30 30	0	0
2	F	42	Total O 42 42	0	0
2	G	6	Total O 6 6	0	0
2	H	17	Total O 17 17	0	0

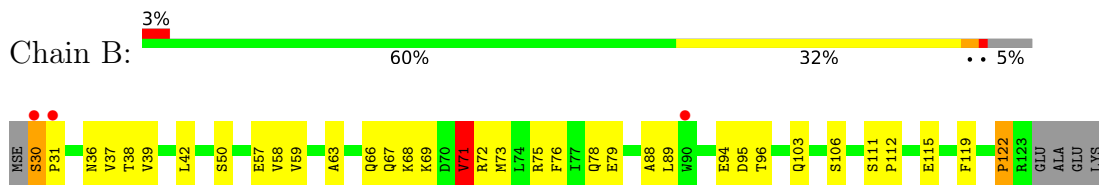
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

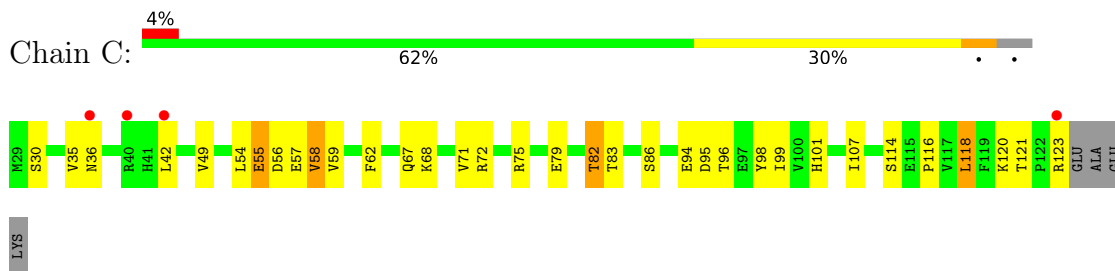
- Molecule 1: Fibronectin type III domain-containing protein 5



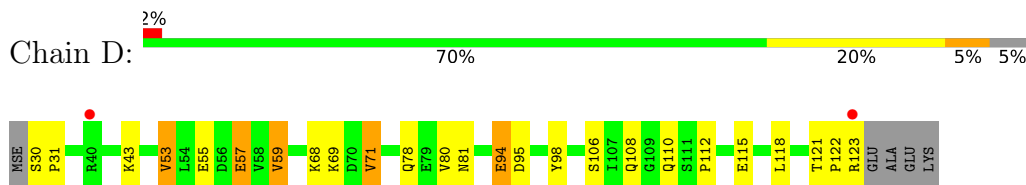
- Molecule 1: Fibronectin type III domain-containing protein 5



- Molecule 1: Fibronectin type III domain-containing protein 5

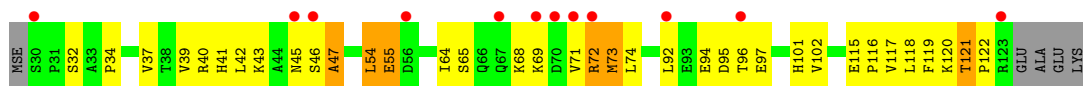


- Molecule 1: Fibronectin type III domain-containing protein 5



- Molecule 1: Fibronectin type III domain-containing protein 5

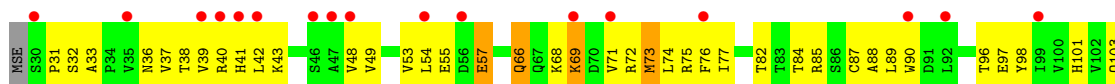




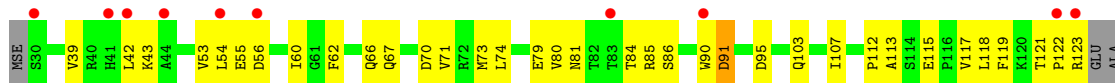
- Molecule 1: Fibronectin type III domain-containing protein 5



- Molecule 1: Fibronectin type III domain-containing protein 5



- Molecule 1: Fibronectin type III domain-containing protein 5



GLU  
LYS

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	93.20Å 93.20Å 284.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	66.52 – 2.28 66.52 – 2.28	Depositor EDS
% Data completeness (in resolution range)	91.6 (66.52-2.28) 91.7 (66.52-2.28)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.67 (at 2.27Å)	Xtrriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.224 , 0.236 0.225 , 0.237	Depositor DCC
$R_{free}$ test set	6327 reflections (11.79%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	48.5	Xtrriage
Anisotropy	0.171	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 45.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6225	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% 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	0.52	0/780	1.05	7/1061 (0.7%)
1	B	0.50	0/753	1.15	11/1025 (1.1%)
1	C	0.48	0/761	1.01	7/1035 (0.7%)
1	D	0.51	0/753	1.06	4/1025 (0.4%)
1	E	0.51	0/753	1.05	3/1025 (0.3%)
1	F	0.49	0/785	1.10	3/1067 (0.3%)
1	G	0.42	0/753	0.97	3/1025 (0.3%)
1	H	0.41	0/753	1.03	5/1025 (0.5%)
All	All	0.48	0/6091	1.05	43/8288 (0.5%)

There are no bond length outliers.

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	90	TRP	N-CA-C	9.23	125.78	113.97
1	A	30	SER	N-CA-C	8.99	116.31	108.22
1	D	71	VAL	N-CA-C	8.82	122.12	109.51
1	E	115	GLU	N-CA-C	-7.64	100.12	109.83
1	F	112	PRO	N-CA-C	-7.59	100.59	111.22
1	B	30	SER	CA-C-N	7.37	127.82	119.93
1	B	30	SER	C-N-CA	7.37	127.82	119.93
1	A	125	ALA	N-CA-C	7.32	120.87	109.96
1	F	115	GLU	N-CA-C	-7.22	100.66	109.83
1	D	112	PRO	N-CA-C	-7.14	100.70	111.41
1	C	114	SER	N-CA-C	-7.12	100.53	110.35
1	H	121	THR	CA-C-N	7.01	127.18	120.31
1	H	121	THR	C-N-CA	7.01	127.18	120.31
1	B	59	VAL	N-CA-C	6.74	117.79	108.89
1	B	112	PRO	N-CA-C	-6.68	101.18	111.13
1	A	114	SER	N-CA-C	-6.65	101.94	110.19
1	B	69	LYS	N-CA-C	-6.56	104.13	111.28
1	H	112	PRO	N-CA-C	-6.46	102.18	111.22
1	A	69	LYS	N-CA-C	-6.43	103.55	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	57	GLU	N-CA-C	6.39	119.06	111.71
1	G	33	ALA	CA-C-N	6.22	126.24	119.89
1	G	33	ALA	C-N-CA	6.22	126.24	119.89
1	H	115	GLU	N-CA-C	-6.02	101.77	110.08
1	E	55	GLU	N-CA-C	6.02	123.62	110.80
1	C	59	VAL	N-CA-C	5.90	116.03	108.12
1	A	121	THR	CA-C-N	5.88	125.85	120.21
1	A	121	THR	C-N-CA	5.88	125.85	120.21
1	B	71	VAL	CB-CA-C	-5.86	101.68	111.29
1	A	71	VAL	N-CA-C	5.85	117.87	109.51
1	D	115	GLU	N-CA-C	-5.81	102.45	109.83
1	D	59	VAL	N-CA-C	5.71	116.43	108.89
1	C	30	SER	CA-C-N	5.70	126.02	119.92
1	C	30	SER	C-N-CA	5.70	126.02	119.92
1	B	111	SER	CA-C-N	-5.55	115.02	120.52
1	B	111	SER	C-N-CA	-5.55	115.02	120.52
1	H	70	ASP	N-CA-C	5.45	119.51	112.86
1	C	83	THR	N-CA-C	-5.44	105.27	112.94
1	B	71	VAL	N-CA-C	5.41	120.59	109.34
1	E	34	PRO	N-CA-C	-5.40	103.32	111.41
1	C	82	THR	CB-CA-C	-5.29	99.83	110.30
1	B	115	GLU	N-CA-C	-5.12	103.32	109.83
1	C	58	VAL	CB-CA-C	-5.12	106.56	111.06
1	G	121	THR	N-CA-C	5.08	117.06	110.40

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	767	0	762	32	0
1	B	740	0	740	19	0
1	C	748	0	749	19	0
1	D	740	0	740	14	0
1	E	740	0	740	27	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	772	0	770	32	0
1	G	740	0	740	50	0
1	H	740	0	740	27	0
2	A	16	0	0	0	0
2	B	50	0	0	1	0
2	C	30	0	0	3	0
2	D	47	0	0	5	0
2	E	30	0	0	0	0
2	F	42	0	0	2	0
2	G	6	0	0	1	0
2	H	17	0	0	2	0
All	All	6225	0	5981	197	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 16.

All (197) 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:C:99:ILE:HG12	1:C:118:LEU:HD12	1.50	0.92
1:A:91:ASP:HB2	1:F:127:LYS:HZ2	1.34	0.91
1:H:84:THR:HG22	1:H:86:SER:H	1.41	0.86
1:E:73:MSE:HE3	1:F:79:GLU:HB3	1.64	0.78
1:G:68:LYS:HB2	1:G:71:VAL:CG1	2.18	0.73
1:G:66:GLN:HE21	1:G:89:LEU:HD23	1.53	0.73
1:E:73:MSE:HE2	1:E:74:LEU:C	2.15	0.71
1:G:107:ILE:HD13	1:G:107:ILE:H	1.55	0.71
1:A:31:PRO:HG3	1:A:106:SER:HB2	1.73	0.70
1:E:73:MSE:HE2	1:E:74:LEU:O	1.90	0.70
1:B:88:ALA:CB	1:F:127:LYS:HE3	2.21	0.70
1:A:73:MSE:HE2	1:A:74:LEU:N	2.07	0.69
1:G:101:HIS:NE2	1:G:116:PRO:HB3	2.08	0.69
1:A:91:ASP:HB2	1:F:127:LYS:NZ	2.08	0.69
1:G:43:LYS:O	1:G:122:PRO:HD2	1.92	0.69
1:E:101:HIS:NE2	1:E:116:PRO:HB3	2.10	0.67
1:G:68:LYS:HB2	1:G:71:VAL:HG12	1.77	0.67
1:B:88:ALA:HB2	1:F:127:LYS:HE3	1.77	0.66
1:C:95:ASP:HA	1:C:121:THR:O	1.94	0.66
1:G:90:TRP:CD1	1:H:90:TRP:CG	2.83	0.66
1:F:30:SER:N	2:F:221:HOH:O	2.29	0.66
1:E:43:LYS:O	1:E:121:THR:HG22	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:31:PRO:HG3	1:G:106:SER:OG	1.97	0.65
1:G:66:GLN:NE2	1:G:89:LEU:HD23	2.13	0.64
1:G:42:LEU:H	1:G:42:LEU:HD12	1.64	0.63
1:A:91:ASP:CB	1:F:127:LYS:HZ2	2.10	0.63
1:A:39:VAL:HG23	1:A:117:VAL:HG11	1.81	0.63
1:E:72:ARG:HG3	1:F:81:ASN:ND2	2.14	0.62
1:D:53:VAL:HG13	2:D:239:HOH:O	1.99	0.62
1:F:64:ILE:CD1	1:F:102:VAL:HG13	2.29	0.62
1:E:39:VAL:HG21	1:E:119:PHE:CD1	2.34	0.62
1:E:37:VAL:O	1:E:117:VAL:HG21	2.00	0.61
1:E:64:ILE:HD13	1:E:102:VAL:HG22	1.81	0.61
1:G:68:LYS:HB2	1:G:71:VAL:HG11	1.83	0.61
1:D:68:LYS:HD3	2:D:214:HOH:O	1.99	0.61
1:G:73:MSE:HE1	1:H:79:GLU:HB3	1.83	0.61
1:B:36:ASN:O	1:B:38:THR:HG23	2.01	0.61
1:G:75:ARG:NH1	1:H:90:TRP:CH2	2.69	0.61
1:A:73:MSE:HE2	1:A:74:LEU:C	2.26	0.60
1:F:59:VAL:CG2	1:F:85:ARG:HH21	2.15	0.60
1:A:73:MSE:CE	1:A:74:LEU:C	2.75	0.59
1:C:35:VAL:HG12	1:C:36:ASN:H	1.67	0.59
1:A:76:PHE:HB2	1:B:78:GLN:HG3	1.84	0.59
1:C:42:LEU:HD11	1:C:120:LYS:O	2.02	0.59
1:G:73:MSE:CE	1:H:79:GLU:HB3	2.33	0.58
1:A:73:MSE:HE3	1:B:79:GLU:HB3	1.85	0.58
1:D:57:GLU:OE1	2:D:233:HOH:O	2.16	0.58
1:G:42:LEU:O	1:G:43:LYS:HD2	2.04	0.58
1:G:43:LYS:C	1:G:122:PRO:HD2	2.29	0.57
1:G:90:TRP:CD1	1:H:90:TRP:CD1	2.92	0.57
1:G:75:ARG:HD3	1:G:77:ILE:HD11	1.87	0.57
1:H:42:LEU:HD21	1:H:119:PHE:CD1	2.40	0.56
1:A:73:MSE:CE	1:A:74:LEU:O	2.54	0.56
1:F:39:VAL:HG23	1:F:47:ALA:HB1	1.88	0.56
1:E:101:HIS:CE1	1:E:116:PRO:HB3	2.41	0.55
1:G:90:TRP:HD1	1:H:90:TRP:CD1	2.24	0.55
1:D:43:LYS:O	1:D:122:PRO:HD2	2.06	0.55
1:G:107:ILE:HD13	1:G:107:ILE:N	2.22	0.55
1:F:59:VAL:HG22	1:F:85:ARG:HH21	1.72	0.55
1:E:69:LYS:HB2	1:E:97:GLU:HB2	1.89	0.55
1:F:55:GLU:O	1:F:56:ASP:HB2	2.06	0.55
1:G:32:SER:HB2	1:G:54:LEU:HD23	1.88	0.54
1:D:68:LYS:HG3	1:D:98:TYR:CE1	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:101:HIS:HD2	2:F:201:HOH:O	1.91	0.54
1:C:101:HIS:CE1	1:C:116:PRO:HG3	2.43	0.54
1:D:95:ASP:HA	1:D:121:THR:O	2.08	0.54
1:H:67:GLN:HB2	1:H:74:LEU:HD23	1.89	0.54
1:A:91:ASP:CG	1:F:127:LYS:HZ3	2.16	0.53
1:G:73:MSE:HE1	1:H:79:GLU:CB	2.38	0.53
1:H:60:ILE:HD11	1:H:107:ILE:HD13	1.90	0.53
1:B:63:ALA:HB1	1:B:76:PHE:CZ	2.44	0.53
1:B:66:GLN:HE21	1:B:89:LEU:HD22	1.74	0.53
1:G:98:TYR:HB2	1:G:119:PHE:CE2	2.43	0.53
1:B:88:ALA:HB3	1:F:127:LYS:HE3	1.90	0.53
1:E:68:LYS:HB2	1:E:71:VAL:CG1	2.39	0.53
1:A:91:ASP:CB	1:F:127:LYS:NZ	2.72	0.52
1:G:105:ILE:HG12	1:G:110:GLN:HG2	1.90	0.52
1:A:68:LYS:NZ	1:A:93:GLU:HG3	2.23	0.52
1:F:74:LEU:HD21	1:F:76:PHE:CE1	2.45	0.52
1:D:78:GLN:HB3	2:D:238:HOH:O	2.08	0.52
1:F:39:VAL:CG2	1:F:47:ALA:HB1	2.40	0.51
1:H:39:VAL:HG23	1:H:117:VAL:HG11	1.91	0.51
1:G:40:ARG:O	1:G:48:VAL:HG12	2.11	0.51
1:F:123:ARG:HH11	1:F:123:ARG:HG3	1.76	0.51
1:G:82:THR:OG1	1:G:84:THR:HG22	2.10	0.51
1:G:74:LEU:HD13	1:G:76:PHE:CZ	2.46	0.51
1:E:97:GLU:HG2	1:E:120:LYS:HE3	1.93	0.51
1:G:89:LEU:HD12	1:G:89:LEU:N	2.27	0.50
1:G:69:LYS:HB2	1:G:97:GLU:HB2	1.93	0.50
1:C:101:HIS:HE1	1:C:116:PRO:HG3	1.75	0.50
1:C:35:VAL:HG12	1:C:36:ASN:N	2.27	0.50
1:E:32:SER:CB	1:E:54:LEU:HD23	2.41	0.50
1:G:42:LEU:C	1:G:43:LYS:HD2	2.36	0.50
1:A:63:ALA:HB3	1:A:103:GLN:HE21	1.77	0.50
1:C:49:VAL:O	1:C:86:SER:HA	2.12	0.50
1:G:53:VAL:HG12	1:G:57:GLU:HB2	1.94	0.50
1:A:31:PRO:HG3	1:A:106:SER:CB	2.40	0.50
1:B:72:ARG:HB3	2:B:201:HOH:O	2.12	0.50
1:B:63:ALA:HB3	1:B:103:GLN:HB3	1.94	0.49
1:C:56:ASP:O	2:C:221:HOH:O	2.20	0.49
1:D:57:GLU:H	1:D:57:GLU:CD	2.20	0.49
1:A:101:HIS:CE1	1:A:116:PRO:HB3	2.47	0.49
1:B:66:GLN:NE2	1:B:89:LEU:HD22	2.28	0.49
1:G:107:ILE:H	1:G:107:ILE:CD1	2.22	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:101:HIS:CE1	1:G:116:PRO:HB3	2.48	0.49
1:H:62:PHE:CE2	1:H:79:GLU:HB2	2.48	0.48
1:E:68:LYS:HE3	1:E:71:VAL:HB	1.96	0.48
1:H:42:LEU:HD21	1:H:119:PHE:HD1	1.78	0.47
1:A:94:GLU:OE1	1:A:124:GLU:CB	2.62	0.47
1:A:40:ARG:HD2	1:F:90:TRP:CZ3	2.50	0.47
1:C:68:LYS:HE2	1:C:96:THR:HG21	1.96	0.47
1:H:53:VAL:HG12	1:H:54:LEU:N	2.30	0.47
1:H:80:VAL:O	1:H:81:ASN:HB2	2.15	0.47
1:A:42:LEU:HD12	1:A:47:ALA:HB2	1.96	0.47
1:D:94:GLU:OE2	1:D:123:ARG:HA	2.15	0.47
1:A:73:MSE:HE1	1:A:74:LEU:O	2.15	0.46
1:D:78:GLN:NE2	2:D:238:HOH:O	2.47	0.46
1:E:97:GLU:CG	1:E:120:LYS:HE3	2.45	0.46
1:E:45:ASN:O	1:E:92:LEU:HB2	2.15	0.46
1:F:64:ILE:HD11	1:F:102:VAL:HG13	1.97	0.46
1:A:37:VAL:HA	1:A:50:SER:O	2.15	0.46
1:F:64:ILE:HD12	1:F:102:VAL:HG13	1.95	0.46
1:C:68:LYS:HG2	2:C:203:HOH:O	2.16	0.46
1:G:54:LEU:HD22	1:G:54:LEU:N	2.32	0.45
1:C:55:GLU:HB2	1:C:58:VAL:HG22	1.97	0.45
1:E:64:ILE:CD1	1:E:102:VAL:HG22	2.45	0.45
1:H:118:LEU:HD12	1:H:118:LEU:N	2.32	0.45
1:E:42:LEU:HD21	1:E:120:LYS:O	2.17	0.45
1:D:53:VAL:HG11	1:D:59:VAL:HG21	1.99	0.44
1:E:73:MSE:CE	1:E:74:LEU:O	2.63	0.44
1:E:94:GLU:O	1:E:96:THR:HG23	2.16	0.44
1:C:55:GLU:O	1:C:58:VAL:HG22	2.18	0.44
1:F:123:ARG:HG3	1:F:123:ARG:NH1	2.32	0.44
1:A:103:GLN:HG3	1:A:111:SER:O	2.17	0.44
1:C:99:ILE:HG12	1:C:118:LEU:CD1	2.34	0.44
1:F:101:HIS:CE1	1:F:116:PRO:HB3	2.52	0.44
1:G:31:PRO:HB2	1:G:111:SER:HB3	2.00	0.44
1:G:88:ALA:HB3	1:H:91:ASP:OD2	2.17	0.44
1:A:73:MSE:HE2	1:A:74:LEU:O	2.17	0.43
1:B:39:VAL:HG21	1:B:119:PHE:CD2	2.53	0.43
1:F:59:VAL:HG22	1:F:85:ARG:NH2	2.33	0.43
1:H:84:THR:HG22	1:H:85:ARG:N	2.33	0.43
1:B:37:VAL:HA	1:B:50:SER:O	2.18	0.43
1:A:73:MSE:HE1	1:A:74:LEU:C	2.43	0.43
1:E:32:SER:HB2	1:E:54:LEU:HD23	1.98	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:37:VAL:HB	1:G:117:VAL:CG2	2.49	0.43
1:C:68:LYS:HA	1:C:98:TYR:HA	2.01	0.43
1:A:62:PHE:CE2	1:A:79:GLU:HB2	2.54	0.43
1:B:30:SER:HA	1:B:31:PRO:HD3	1.80	0.43
1:A:62:PHE:CD1	1:A:62:PHE:C	2.96	0.42
1:E:46:SER:O	1:E:47:ALA:HB2	2.19	0.42
1:E:118:LEU:HD12	1:E:118:LEU:N	2.34	0.42
1:H:43:LYS:C	1:H:122:PRO:HD2	2.44	0.42
1:E:45:ASN:HA	1:E:92:LEU:O	2.19	0.42
1:H:122:PRO:HA	2:H:206:HOH:O	2.19	0.42
1:B:67:GLN:NE2	1:B:71:VAL:HG13	2.34	0.42
1:E:95:ASP:HA	1:E:121:THR:O	2.18	0.42
1:G:40:ARG:HD3	1:G:41:HIS:CE1	2.54	0.42
1:C:35:VAL:HG11	2:G:203:HOH:O	2.20	0.42
1:G:96:THR:O	1:G:120:LYS:HA	2.18	0.42
1:A:41:HIS:O	1:A:43:LYS:HE2	2.19	0.42
1:B:58:VAL:O	1:B:106:SER:HB2	2.20	0.42
1:F:74:LEU:HD21	1:F:76:PHE:HE1	1.84	0.42
1:H:53:VAL:HG12	1:H:54:LEU:H	1.85	0.42
1:A:49:VAL:O	1:A:86:SER:HA	2.20	0.42
1:A:97:GLU:HG2	1:A:120:LYS:HB2	2.02	0.42
1:H:60:ILE:O	1:H:81:ASN:N	2.53	0.42
1:A:91:ASP:CG	1:F:127:LYS:NZ	2.77	0.42
1:A:72:ARG:HH11	1:A:72:ARG:HB2	1.85	0.42
1:G:31:PRO:HG2	1:G:109:GLY:O	2.20	0.42
1:B:73:MSE:SE	1:B:75:ARG:HB2	2.71	0.41
1:F:124:GLU:O	1:F:126:GLU:HG2	2.20	0.41
1:D:30:SER:N	1:D:31:PRO:HD2	2.35	0.41
1:G:41:HIS:HB2	1:G:48:VAL:HG12	2.02	0.41
1:D:80:VAL:O	1:D:81:ASN:HB2	2.20	0.41
1:B:68:LYS:HD2	1:B:96:THR:HG21	2.02	0.41
1:C:123:ARG:N	1:C:123:ARG:HD2	2.36	0.41
1:G:38:THR:HG22	1:G:39:VAL:N	2.35	0.41
1:G:54:LEU:HA	1:G:54:LEU:HD13	1.86	0.41
1:G:72:ARG:HG3	1:G:72:ARG:O	2.21	0.41
1:D:106:SER:C	1:D:108:GLN:H	2.29	0.41
1:G:43:LYS:O	1:G:121:THR:HG22	2.21	0.41
1:B:95:ASP:N	1:B:122:PRO:O	2.47	0.41
1:C:75:ARG:NH2	2:C:212:HOH:O	2.53	0.41
1:F:68:LYS:HE3	1:F:96:THR:HG23	2.03	0.41
1:G:32:SER:CB	1:G:54:LEU:HD23	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:49:VAL:HG12	1:G:87:CYS:O	2.21	0.41
1:G:75:ARG:NH1	1:H:90:TRP:CZ3	2.82	0.41
1:G:121:THR:HA	1:G:122:PRO:HD3	1.91	0.41
1:H:66:GLN:NE2	2:H:210:HOH:O	2.32	0.41
1:H:71:VAL:HG12	1:H:73:MSE:H	1.85	0.41
1:H:103:GLN:HG2	1:H:113:ALA:HA	2.03	0.41
1:C:62:PHE:CE2	1:C:79:GLU:HB2	2.55	0.41
1:F:60:ILE:HD12	1:F:60:ILE:C	2.46	0.40
1:E:40:ARG:O	1:E:41:HIS:C	2.63	0.40
1:F:80:VAL:O	1:F:81:ASN:HB2	2.21	0.40
1:G:98:TYR:CB	1:G:119:PHE:CE2	3.05	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	96/99 (97%)	92 (96%)	3 (3%)	1 (1%)	12	13
1	B	92/99 (93%)	85 (92%)	6 (6%)	1 (1%)	11	11
1	C	93/99 (94%)	86 (92%)	6 (6%)	1 (1%)	11	11
1	D	92/99 (93%)	88 (96%)	4 (4%)	0	100	100
1	E	92/99 (93%)	82 (89%)	7 (8%)	3 (3%)	3	1
1	F	96/99 (97%)	92 (96%)	3 (3%)	1 (1%)	12	13
1	G	92/99 (93%)	80 (87%)	11 (12%)	1 (1%)	11	11
1	H	92/99 (93%)	88 (96%)	3 (3%)	1 (1%)	11	11
All	All	745/792 (94%)	693 (93%)	43 (6%)	9 (1%)	10	10

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	122	PRO
1	E	55	GLU
1	G	55	GLU
1	E	47	ALA
1	H	55	GLU
1	F	81	ASN
1	A	72	ARG
1	C	72	ARG
1	E	122	PRO

### 5.3.2 Protein sidechains [i](#)

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	87/87 (100%)	79 (91%)	8 (9%)	8	10
1	B	85/87 (98%)	82 (96%)	3 (4%)	32	45
1	C	86/87 (99%)	77 (90%)	9 (10%)	6	7
1	D	85/87 (98%)	77 (91%)	8 (9%)	8	9
1	E	85/87 (98%)	80 (94%)	5 (6%)	18	23
1	F	88/87 (101%)	82 (93%)	6 (7%)	14	18
1	G	85/87 (98%)	76 (89%)	9 (11%)	6	7
1	H	85/87 (98%)	81 (95%)	4 (5%)	23	33
All	All	686/696 (99%)	634 (92%)	52 (8%)	12	15

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	THR
1	A	40	ARG
1	A	43	LYS
1	A	54	LEU
1	A	55	GLU
1	A	73	MSE
1	A	94	GLU

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	126	GLU
1	B	42	LEU
1	B	71	VAL
1	B	94	GLU
1	C	54	LEU
1	C	55	GLU
1	C	57	GLU
1	C	67	GLN
1	C	71	VAL
1	C	82	THR
1	C	94	GLU
1	C	107	ILE
1	C	118	LEU
1	D	53	VAL
1	D	55	GLU
1	D	57	GLU
1	D	69	LYS
1	D	71	VAL
1	D	94	GLU
1	D	110	GLN
1	D	118	LEU
1	E	54	LEU
1	E	65	SER
1	E	72	ARG
1	E	73	MSE
1	E	121	THR
1	F	39	VAL
1	F	42	LEU
1	F	53	VAL
1	F	64	ILE
1	F	69	LYS
1	F	124	GLU
1	G	36	ASN
1	G	57	GLU
1	G	66	GLN
1	G	69	LYS
1	G	73	MSE
1	G	85	ARG
1	G	103	GLN
1	G	107	ILE
1	G	119	PHE
1	H	56	ASP

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Mol	Chain	Res	Type
1	H	91	ASP
1	H	95	ASP
1	H	123	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	ASN
1	A	45	ASN
1	A	101	HIS
1	A	103	GLN
1	A	108	GLN
1	A	110	GLN
1	B	41	HIS
1	B	66	GLN
1	B	67	GLN
1	B	103	GLN
1	B	108	GLN
1	C	45	ASN
1	C	101	HIS
1	C	103	GLN
1	C	110	GLN
1	D	36	ASN
1	D	45	ASN
1	D	110	GLN
1	E	66	GLN
1	E	67	GLN
1	E	78	GLN
1	E	103	GLN
1	E	108	GLN
1	E	110	GLN
1	F	101	HIS
1	F	103	GLN
1	F	108	GLN
1	G	41	HIS
1	G	66	GLN
1	G	67	GLN
1	G	78	GLN
1	H	101	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 oligosaccharides 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	96/99 (96%)	0.37	3 (3%) 51 53	37, 49, 69, 94	0
1	B	93/99 (93%)	0.05	3 (3%) 50 52	30, 40, 54, 65	0
1	C	93/99 (93%)	0.42	4 (4%) 40 41	32, 49, 68, 79	0
1	D	93/99 (93%)	-0.18	2 (2%) 62 64	26, 38, 56, 69	0
1	E	93/99 (93%)	0.53	12 (12%) 7 8	26, 43, 82, 91	0
1	F	97/99 (97%)	0.00	4 (4%) 41 43	29, 39, 60, 85	0
1	G	93/99 (93%)	1.44	24 (25%) 1 1	40, 72, 94, 105	0
1	H	93/99 (93%)	0.82	10 (10%) 11 12	41, 59, 77, 96	0
All	All	751/792 (94%)	0.43	62 (8%) 17 18	26, 47, 83, 105	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	90	TRP	5.2
1	F	125	ALA	5.1
1	G	42	LEU	5.0
1	A	125	ALA	4.7
1	F	127	LYS	4.2
1	E	30	SER	4.1
1	E	71	VAL	3.8
1	G	123	ARG	3.7
1	H	30	SER	3.7
1	G	117	VAL	3.6
1	G	118	LEU	3.6
1	G	48	VAL	3.5
1	H	123	ARG	3.5
1	F	126	GLU	3.4
1	C	42	LEU	3.2
1	G	40	ARG	3.1

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	H	54	LEU	3.1
1	E	72	ARG	3.0
1	G	116	PRO	3.0
1	G	35	VAL	3.0
1	H	90	TRP	3.0
1	E	70	ASP	2.9
1	G	39	VAL	2.8
1	G	56	ASP	2.8
1	A	126	GLU	2.8
1	C	123	ARG	2.7
1	E	123	ARG	2.7
1	G	108	GLN	2.6
1	G	30	SER	2.6
1	H	41	HIS	2.6
1	E	45	ASN	2.6
1	G	69	LYS	2.6
1	D	123	ARG	2.5
1	G	71	VAL	2.4
1	B	31	PRO	2.4
1	D	40	ARG	2.4
1	G	41	HIS	2.4
1	H	56	ASP	2.4
1	H	42	LEU	2.4
1	E	69	LYS	2.3
1	G	47	ALA	2.3
1	G	46	SER	2.3
1	G	119	PHE	2.3
1	H	44	ALA	2.3
1	A	124	GLU	2.3
1	E	56	ASP	2.2
1	E	46	SER	2.2
1	B	90	TRP	2.2
1	C	36	ASN	2.2
1	E	92	LEU	2.2
1	F	123	ARG	2.2
1	E	96	THR	2.2
1	G	54	LEU	2.1
1	G	92	LEU	2.1
1	G	99	ILE	2.1
1	G	76	PHE	2.1
1	G	115	GLU	2.1
1	C	40	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	H	122	PRO	2.1
1	B	30	SER	2.1
1	E	67	GLN	2.0
1	H	83	THR	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 oligosaccharides 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.