



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 23, 2024 – 01:13 PM EDT

PDB ID : 4LSD
Title : Myokine structure
Authors : Schumacher, M.A.; Ohashi, T.; Shah, R.S.; Chinnam, N.; Erickson, H.
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

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>.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.36.1
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) : 2.36.1

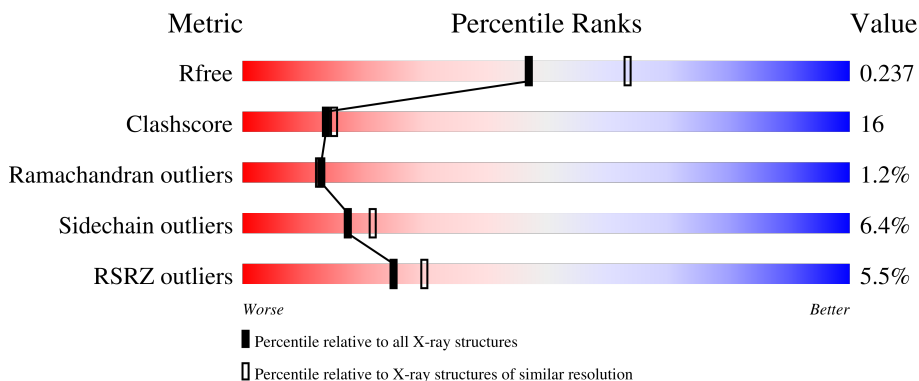
1 Overall quality at a glance

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.



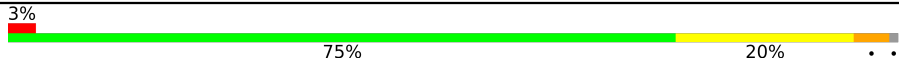


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6980 (2.30-2.26)
Clashscore	141614	7711 (2.30-2.26)
Ramachandran outliers	138981	7597 (2.30-2.26)
Sidechain outliers	138945	7598 (2.30-2.26)
RSRZ outliers	127900	6849 (2.30-2.26)

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 ≥ 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	 3% 64% 31% . .
1	B	99	 66% 27% .. 5%
1	C	99	 4% 67% 26% . .
1	D	99	 75% 18% . 5%
1	E	99	 63% 27% 5% 5%

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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

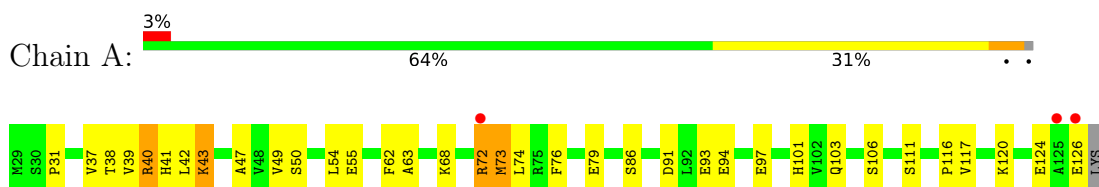
- Molecule 2 is water.

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

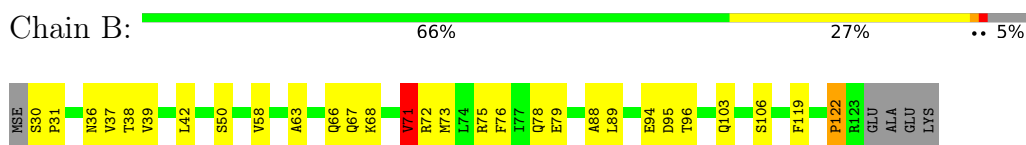
3 Residue-property plots [i](#)

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.

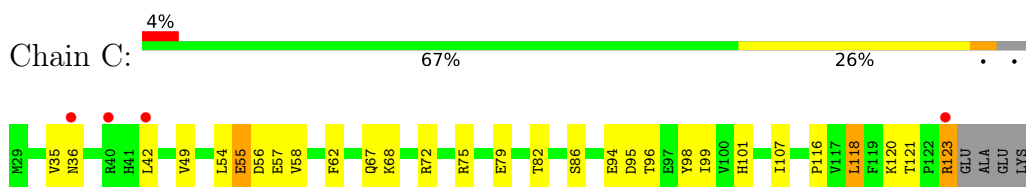
- Molecule 1: Fibronectin type III domain-containing protein 5



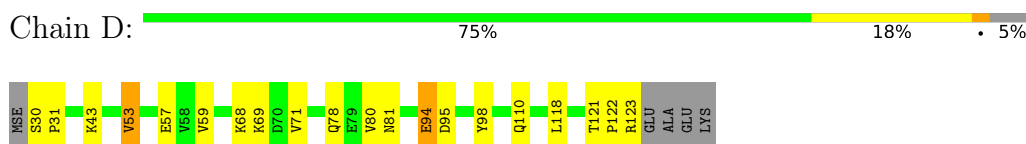
- Molecule 1: Fibronectin type III domain-containing protein 5



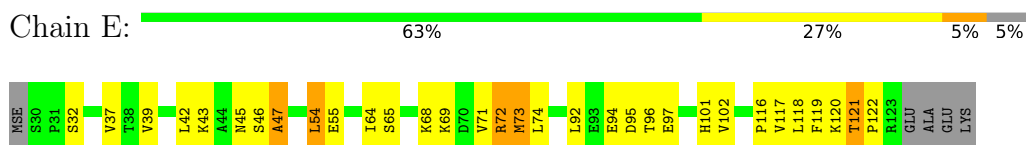
- Molecule 1: Fibronectin type III domain-containing protein 5



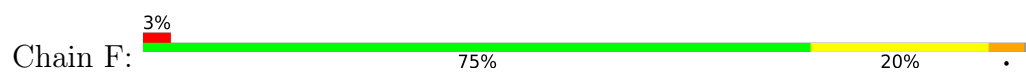
- Molecule 1: Fibronectin type III domain-containing protein 5



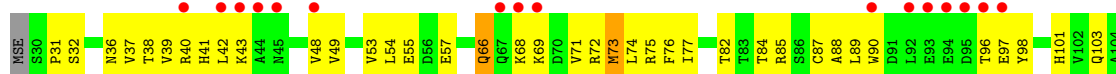
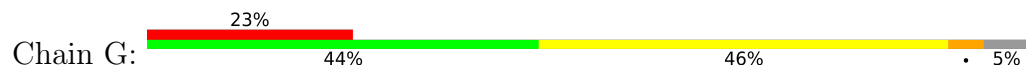
- Molecule 1: Fibronectin type III domain-containing protein 5



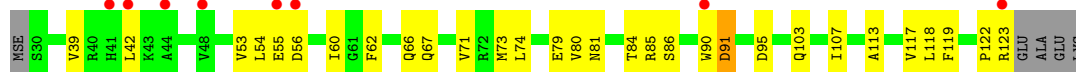
- Molecule 1: Fibronectin type III domain-containing protein 5



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	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) 91.6 (66.52-2.28)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	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 ²)	48.5	Xtrriage
Anisotropy	0.171	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 46.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	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 ²)	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.*

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	0/780	0.71	0/1061
1	B	0.44	0/753	0.77	1/1025 (0.1%)
1	C	0.42	0/761	0.72	0/1035
1	D	0.44	0/753	0.77	0/1025
1	E	0.47	0/753	0.74	0/1025
1	F	0.43	0/785	0.80	1/1067 (0.1%)
1	G	0.39	0/753	0.65	0/1025
1	H	0.37	0/753	0.67	0/1025
All	All	0.42	0/6091	0.73	2/8288 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	90	TRP	N-CA-C	5.47	125.78	111.00
1	B	71	VAL	CB-CA-C	-5.12	101.68	111.40

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	29	0
1	B	740	0	740	19	0
1	C	748	0	749	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	740	0	740	13	0
1	E	740	0	740	26	0
1	F	772	0	770	30	0
1	G	740	0	740	50	0
1	H	740	0	740	26	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	190	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 (190) 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: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:A:91:ASP:HB2	1:F:127:LYS:HZ2	1.48	0.77
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: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:73:MSE:HE2	1:E:74:LEU:C	2.15	0.67
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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: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:91:ASP:CG	1:F:127:LYS:HZ3	2.04	0.60
1:F:59:VAL:CG2	1:F:85:ARG:HH21	2.15	0.60
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: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:HE2	1:A:74:LEU:C	2.26	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:A:73:MSE:CE	1:A:74:LEU:C	2.75	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:68:LYS:HG3	1:D:98:TYR:CE1	2.42	0.54
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: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:43:LYS:C	1:G:122:PRO:HD2	2.29	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: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: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
1:G:101:HIS:CE1	1:G:116:PRO:HB3	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:42:LEU:C	1:G:43:LYS:HD2	2.36	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:D:57:GLU:H	1:D:57:GLU:CD	2.20	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: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: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: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:A:91:ASP:CB	1:F:127:LYS:HZ2	2.23	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: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: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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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%)	15	16
1	B	92/99 (93%)	85 (92%)	6 (6%)	1 (1%)	14	14
1	C	93/99 (94%)	86 (92%)	6 (6%)	1 (1%)	14	14
1	D	92/99 (93%)	88 (96%)	4 (4%)	0	100	100
1	E	92/99 (93%)	82 (89%)	7 (8%)	3 (3%)	4	2
1	F	96/99 (97%)	92 (96%)	3 (3%)	1 (1%)	15	16
1	G	92/99 (93%)	80 (87%)	11 (12%)	1 (1%)	14	14
1	H	92/99 (93%)	88 (96%)	3 (3%)	1 (1%)	14	14
All	All	745/792 (94%)	693 (93%)	43 (6%)	9 (1%)	13	12

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

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Mol	Chain	Res	Type
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%)	80 (92%)	7 (8%)	12	14
1	B	85/87 (98%)	82 (96%)	3 (4%)	36	48
1	C	86/87 (99%)	77 (90%)	9 (10%)	7	7
1	D	85/87 (98%)	79 (93%)	6 (7%)	14	17
1	E	85/87 (98%)	80 (94%)	5 (6%)	19	24
1	F	88/87 (101%)	85 (97%)	3 (3%)	37	49
1	G	85/87 (98%)	78 (92%)	7 (8%)	11	13
1	H	85/87 (98%)	81 (95%)	4 (5%)	26	34
All	All	686/696 (99%)	642 (94%)	44 (6%)	17	21

All (44) 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	126	GLU
1	B	42	LEU
1	B	71	VAL
1	B	94	GLU

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Mol	Chain	Res	Type
1	C	54	LEU
1	C	55	GLU
1	C	57	GLU
1	C	67	GLN
1	C	82	THR
1	C	94	GLU
1	C	107	ILE
1	C	118	LEU
1	C	123	ARG
1	D	53	VAL
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	42	LEU
1	F	64	ILE
1	F	124	GLU
1	G	36	ASN
1	G	66	GLN
1	G	73	MSE
1	G	85	ARG
1	G	103	GLN
1	G	107	ILE
1	G	119	PHE
1	H	56	ASP
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 (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	ASN
1	A	101	HIS
1	A	103	GLN
1	A	110	GLN

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Mol	Chain	Res	Type
1	B	66	GLN
1	B	67	GLN
1	B	103	GLN
1	B	108	GLN
1	C	101	HIS
1	C	103	GLN
1	C	110	GLN
1	D	36	ASN
1	D	45	ASN
1	D	66	GLN
1	D	110	GLN
1	E	45	ASN
1	E	67	GLN
1	E	108	GLN
1	F	66	GLN
1	F	81	ASN
1	F	101	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 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

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, 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	96/99 (96%)	0.41	3 (3%) 49 54	37, 49, 69, 94	0
1	B	93/99 (93%)	0.30	0 100 100	30, 40, 54, 65	0
1	C	93/99 (93%)	0.32	4 (4%) 35 40	32, 49, 68, 79	0
1	D	93/99 (93%)	0.26	0 100 100	26, 38, 56, 69	0
1	E	93/99 (93%)	0.52	0 100 100	26, 43, 82, 91	0
1	F	97/99 (97%)	0.30	3 (3%) 49 54	29, 39, 60, 85	0
1	G	93/99 (93%)	1.19	23 (24%) 0 0	40, 72, 94, 105	0
1	H	93/99 (93%)	0.59	8 (8%) 10 13	41, 59, 77, 96	0
All	All	751/792 (94%)	0.48	41 (5%) 25 30	26, 47, 83, 105	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	123	ARG	4.3
1	G	90	TRP	4.3
1	F	125	ALA	4.2
1	H	90	TRP	4.1
1	G	69	LYS	4.0
1	G	119	PHE	3.8
1	G	68	LYS	3.5
1	G	92	LEU	3.5
1	F	127	LYS	3.5
1	A	125	ALA	3.5
1	H	44	ALA	3.4
1	G	95	ASP	3.4
1	G	40	ARG	3.4
1	G	118	LEU	3.3
1	G	48	VAL	3.1
1	H	123	ARG	3.1

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Mol	Chain	Res	Type	RSRZ
1	G	42	LEU	3.1
1	G	122	PRO	3.0
1	A	126	GLU	3.0
1	G	45	ASN	3.0
1	F	126	GLU	2.8
1	C	42	LEU	2.8
1	G	96	THR	2.8
1	C	40	ARG	2.7
1	G	44	ALA	2.7
1	G	43	LYS	2.7
1	C	36	ASN	2.7
1	G	93	GLU	2.7
1	G	94	GLU	2.6
1	G	120	LYS	2.6
1	H	55	GLU	2.5
1	H	56	ASP	2.5
1	H	42	LEU	2.3
1	G	97	GLU	2.3
1	G	116	PRO	2.2
1	G	67	GLN	2.2
1	C	123	ARG	2.2
1	H	48	VAL	2.1
1	G	117	VAL	2.1
1	H	41	HIS	2.1
1	A	72	ARG	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no 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.