



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2017 – 11:08 pm GMT

PDB ID : 1RKC  
Title : Human vinculin head (1-258) in complex with talin's vinculin binding site 3 (residues 1944-1969)  
Authors : Izard, T.; Evans, G.; Borgon, R.A.; Rush, C.L.; Bricogne, G.; Bois, P.R.  
Deposited on : 2003-11-21  
Resolution : 2.70 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) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

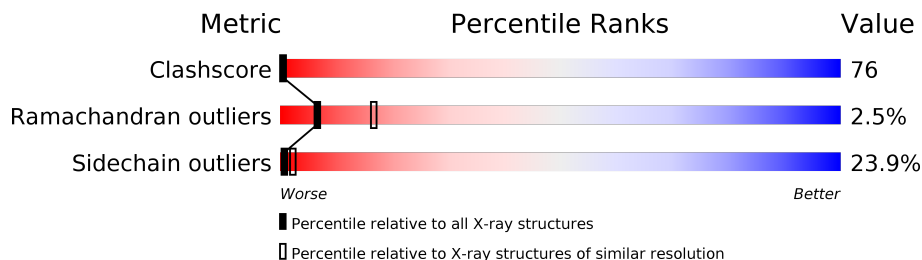
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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))
Clashscore	112137	2590 (2.70-2.70)
Ramachandran outliers	110173	2550 (2.70-2.70)
Sidechain outliers	110143	2550 (2.70-2.70)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	262	
2	B	26	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2223 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 Vinculin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	0	0
			2020	1272	342	392	14			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	HIS	-	EXPRESSION TAG	UNP P18206
A	-2	HIS	-	EXPRESSION TAG	UNP P18206
A	-1	HIS	-	EXPRESSION TAG	UNP P18206
A	0	HIS	-	EXPRESSION TAG	UNP P18206

- Molecule 2 is a protein called Talin.

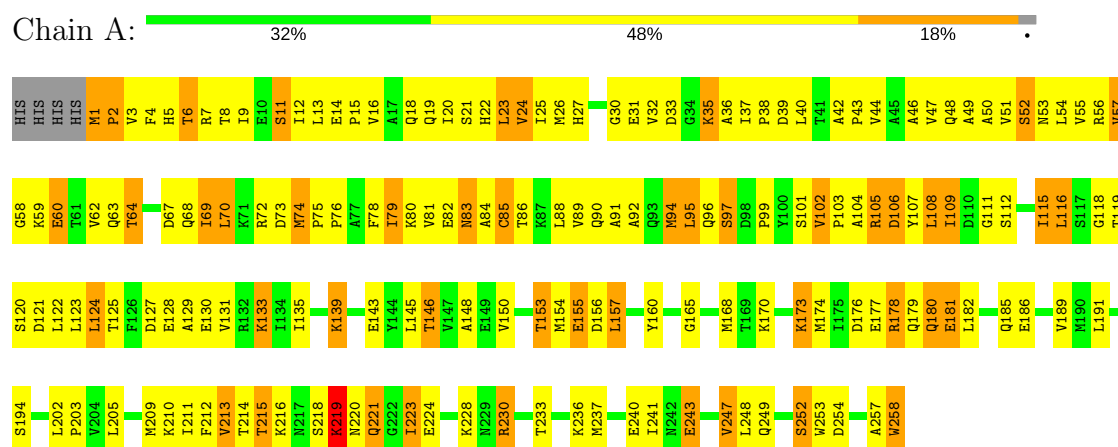
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	26	Total	C	N	O	0	0	0
			203	129	36	38			

### 3 Residue-property plots

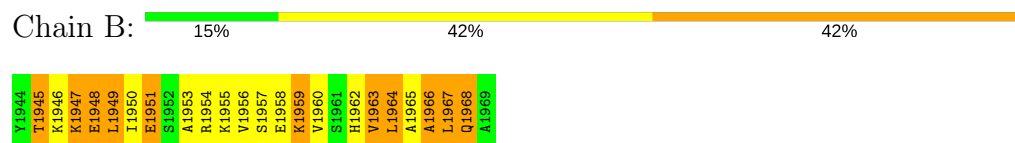
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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.

Note EDS was not executed.

#### • Molecule 1: Vinculin



#### • Molecule 2: Talin



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	I 41 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	155.68Å 155.68Å 155.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.70	Depositor
% Data completeness (in resolution range)	99.7 (15.00-2.70)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	BUSTER-TNT	Depositor
R, $R_{free}$	0.244 , 0.335	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2223	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	113.0	wwPDB-VP

## 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	0.60	0/2047	0.82	3/2773 (0.1%)
2	B	0.42	0/204	0.67	0/272
All	All	0.58	0/2251	0.81	3/3045 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1	MET	C-N-CD	-6.70	105.85	120.60
1	A	102	VAL	C-N-CD	-6.54	106.20	120.60
1	A	165	GLY	C-N-CD	-5.67	108.14	120.60

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	2020	0	2074	323	0
2	B	203	0	219	52	0
All	All	2223	0	2293	344	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 76.

All (344) 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:26:MET:HB3	1:A:31:GLU:OE2	1.38	1.18
1:A:81:VAL:HG23	1:A:118:GLY:HA3	1.24	1.17
1:A:153:THR:HG22	1:A:156:ASP:H	1.07	1.10
1:A:37:ILE:HA	1:A:40:LEU:HD12	1.13	1.09
1:A:54:LEU:HD21	2:B:1953:ALA:HB2	1.36	1.08
1:A:43:PRO:HB2	2:B:1963:VAL:HG23	1.35	1.03
1:A:75:PRO:HA	1:A:78:PHE:HD1	1.24	1.01
1:A:74:MET:HG3	1:A:125:THR:HG21	1.43	0.97
1:A:12:ILE:HG22	1:A:13:LEU:HD23	1.47	0.96
1:A:26:MET:CB	1:A:31:GLU:OE2	2.14	0.95
1:A:178:ARG:HH11	1:A:178:ARG:HG2	1.27	0.95
1:A:88:LEU:HD22	1:A:108:LEU:HD22	1.50	0.93
1:A:35:LYS:NZ	1:A:99:PRO:HA	1.84	0.93
1:A:88:LEU:HD11	1:A:112:SER:HB3	1.51	0.93
1:A:84:ALA:HB1	1:A:115:ILE:HG12	1.50	0.91
1:A:75:PRO:HA	1:A:78:PHE:CD1	2.06	0.91
1:A:23:LEU:HD22	2:B:1967:LEU:HB2	1.52	0.90
1:A:115:ILE:HG21	2:B:1960:VAL:HG21	1.53	0.90
1:A:202:LEU:HB3	1:A:203:PRO:HD3	1.54	0.90
1:A:88:LEU:HD22	1:A:108:LEU:HA	1.53	0.89
1:A:101:SER:HB3	1:A:104:ALA:HB3	1.55	0.88
1:A:153:THR:HG22	1:A:156:ASP:N	1.89	0.88
1:A:88:LEU:HD11	1:A:112:SER:CB	2.04	0.87
1:A:88:LEU:HD22	1:A:108:LEU:CD2	2.04	0.87
1:A:69:ILE:HD13	1:A:257:ALA:CB	2.04	0.87
1:A:88:LEU:CD2	1:A:108:LEU:HA	2.05	0.86
1:A:84:ALA:CB	1:A:115:ILE:HG12	2.05	0.86
1:A:44:VAL:CG1	1:A:89:VAL:HG22	2.05	0.86
1:A:211:ILE:O	1:A:215:THR:HB	1.76	0.85
1:A:178:ARG:O	1:A:178:ARG:HD3	1.75	0.85
1:A:91:ALA:HA	1:A:94:MET:CG	2.08	0.84
1:A:46:ALA:HA	1:A:49:ALA:HB3	1.60	0.83
1:A:91:ALA:HA	1:A:94:MET:HG2	1.61	0.82
1:A:43:PRO:HB2	2:B:1963:VAL:CG2	2.08	0.82
1:A:23:LEU:CD2	2:B:1967:LEU:HB2	2.10	0.81
1:A:94:MET:HB2	1:A:104:ALA:HB2	1.62	0.81
1:A:81:VAL:HG23	1:A:118:GLY:CA	2.08	0.81
1:A:88:LEU:HD11	1:A:112:SER:CA	2.11	0.80
1:A:249:GLN:HG2	1:A:253:TRP:CZ3	2.17	0.79
1:A:35:LYS:HZ2	1:A:95:LEU:HD13	1.48	0.79
1:A:55:VAL:HA	1:A:78:PHE:HE2	1.46	0.79
1:A:69:ILE:HD13	1:A:257:ALA:HB1	1.66	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:MET:HG3	1:A:104:ALA:HB1	1.65	0.78
1:A:68:GLN:O	1:A:72:ARG:HG3	1.83	0.77
1:A:4:PHE:CZ	1:A:13:LEU:HB2	2.19	0.77
1:A:37:ILE:HA	1:A:40:LEU:CD1	2.06	0.77
1:A:54:LEU:CD2	2:B:1953:ALA:HB2	2.12	0.77
1:A:4:PHE:CE2	1:A:13:LEU:HB2	2.20	0.76
1:A:170:LYS:O	1:A:174:MET:HG3	1.84	0.76
1:A:249:GLN:HG2	1:A:253:TRP:HZ3	1.48	0.76
1:A:55:VAL:O	1:A:59:LYS:HG2	1.85	0.76
1:A:105:ARG:HH21	2:B:1967:LEU:HA	1.50	0.76
1:A:35:LYS:HD2	1:A:37:ILE:HG12	1.67	0.76
1:A:74:MET:CG	1:A:125:THR:HG21	2.15	0.75
1:A:233:THR:O	1:A:237:MET:HG3	1.86	0.75
1:A:215:THR:HG23	1:A:221:GLN:CG	2.16	0.75
1:A:6:THR:HG21	1:A:181:GLU:HG3	1.69	0.75
1:A:178:ARG:HG2	1:A:178:ARG:NH1	1.91	0.74
1:A:74:MET:HG3	1:A:125:THR:CG2	2.15	0.73
1:A:55:VAL:HA	1:A:78:PHE:CE2	2.22	0.73
1:A:78:PHE:O	1:A:82:GLU:HG2	1.89	0.73
1:A:88:LEU:HD21	1:A:112:SER:H	1.52	0.73
1:A:54:LEU:HD11	2:B:1953:ALA:CB	2.18	0.72
1:A:24:VAL:HG12	1:A:25:ILE:N	2.04	0.72
1:A:154:MET:HE2	1:A:214:THR:HG23	1.72	0.72
1:A:6:THR:CG2	1:A:181:GLU:HG3	2.21	0.71
1:A:178:ARG:HD3	1:A:178:ARG:C	2.10	0.70
1:A:219:LYS:HD2	1:A:220:ASN:HB2	1.74	0.70
2:B:1966:ALA:C	2:B:1967:LEU:HD23	2.12	0.70
1:A:191:LEU:HD21	1:A:247:VAL:HG22	1.74	0.70
1:A:215:THR:HG23	1:A:221:GLN:HG3	1.73	0.70
1:A:91:ALA:HA	1:A:94:MET:SD	2.32	0.69
2:B:1967:LEU:HD23	2:B:1967:LEU:N	2.07	0.69
1:A:37:ILE:CG2	1:A:96:GLN:HG3	2.22	0.69
1:A:133:LYS:HE3	1:A:174:MET:CE	2.23	0.69
1:A:9:ILE:CD1	1:A:123:LEU:HB3	2.22	0.69
1:A:40:LEU:HA	1:A:43:PRO:CG	2.23	0.68
1:A:46:ALA:HA	1:A:49:ALA:CB	2.23	0.68
1:A:12:ILE:HG22	1:A:13:LEU:CD2	2.22	0.68
1:A:35:LYS:HZ3	1:A:99:PRO:HA	1.58	0.68
1:A:35:LYS:HD3	1:A:95:LEU:HB3	1.75	0.68
1:A:35:LYS:HZ1	1:A:99:PRO:HA	1.59	0.68
1:A:67:ASP:HB3	1:A:70:LEU:HB2	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:PRO:O	1:A:79:ILE:HB	1.94	0.67
1:A:95:LEU:HD23	1:A:104:ALA:CB	2.24	0.67
1:A:55:VAL:HG22	1:A:78:PHE:CD2	2.30	0.67
1:A:55:VAL:HG22	1:A:78:PHE:HD2	1.59	0.67
1:A:154:MET:CE	1:A:214:THR:HG23	2.25	0.67
1:A:42:ALA:N	1:A:43:PRO:HD2	2.10	0.66
1:A:128:GLU:HG2	1:A:252:SER:OG	1.95	0.66
2:B:1947:LYS:O	2:B:1950:ILE:N	2.24	0.66
1:A:133:LYS:HE3	1:A:174:MET:HE2	1.78	0.65
1:A:35:LYS:CD	1:A:95:LEU:HD13	2.27	0.64
1:A:116:LEU:O	1:A:119:THR:N	2.30	0.64
1:A:91:ALA:HB3	1:A:108:LEU:CD2	2.28	0.64
1:A:101:SER:HB3	1:A:104:ALA:CB	2.26	0.64
1:A:35:LYS:HZ2	1:A:95:LEU:CD1	2.11	0.64
1:A:4:PHE:HE2	1:A:13:LEU:HD12	1.63	0.64
1:A:89:VAL:O	1:A:92:ALA:HB3	1.97	0.63
1:A:237:MET:O	1:A:241:ILE:HG13	1.99	0.63
1:A:50:ALA:HA	1:A:53:ASN:ND2	2.13	0.63
1:A:35:LYS:NZ	1:A:95:LEU:HD13	2.13	0.63
1:A:9:ILE:HD13	1:A:123:LEU:HB3	1.79	0.63
1:A:37:ILE:HD13	1:A:96:GLN:HA	1.81	0.63
1:A:54:LEU:HD11	2:B:1953:ALA:HB1	1.80	0.62
1:A:44:VAL:HG13	1:A:89:VAL:HG22	1.79	0.61
1:A:53:ASN:O	1:A:57:VAL:HG23	1.99	0.61
1:A:60:GLU:O	1:A:64:THR:HB	2.00	0.61
1:A:102:VAL:HB	1:A:103:PRO:HD3	1.82	0.61
1:A:14:GLU:N	1:A:15:PRO:HD2	2.14	0.61
1:A:215:THR:HG22	1:A:223:ILE:HG12	1.82	0.61
1:A:148:ALA:HB2	1:A:160:TYR:OH	2.01	0.61
1:A:131:VAL:HG13	1:A:248:LEU:HD22	1.81	0.61
1:A:128:GLU:CG	1:A:252:SER:OG	2.48	0.60
1:A:88:LEU:HD21	1:A:112:SER:N	2.16	0.60
1:A:191:LEU:HD21	1:A:247:VAL:CG2	2.31	0.60
1:A:58:GLY:C	1:A:78:PHE:HZ	2.05	0.60
1:A:46:ALA:HB1	2:B:1959:LYS:NZ	2.16	0.60
1:A:153:THR:HG23	1:A:155:GLU:H	1.67	0.60
1:A:35:LYS:NZ	1:A:95:LEU:HD22	2.17	0.60
1:A:218:SER:O	1:A:221:GLN:HB2	2.03	0.59
1:A:37:ILE:HG21	1:A:96:GLN:CG	2.32	0.59
2:B:1963:VAL:HG12	2:B:1964:LEU:HD23	1.83	0.59
1:A:40:LEU:HA	1:A:43:PRO:HG3	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:PRO:N	1:A:76:PRO:HD2	2.17	0.59
1:A:35:LYS:HZ2	1:A:95:LEU:CG	2.16	0.59
2:B:1945:THR:HA	2:B:1948:GLU:OE2	2.03	0.58
1:A:36:ALA:O	1:A:40:LEU:HG	2.03	0.58
1:A:81:VAL:HG13	1:A:115:ILE:HD13	1.84	0.58
2:B:1945:THR:HG23	2:B:1948:GLU:OE2	2.03	0.58
1:A:74:MET:CB	1:A:125:THR:HG21	2.34	0.58
1:A:223:ILE:HG22	1:A:224:GLU:N	2.17	0.58
1:A:37:ILE:HG21	1:A:96:GLN:HG3	1.83	0.58
1:A:94:MET:SD	1:A:104:ALA:HA	2.44	0.58
1:A:257:ALA:O	1:A:258:TRP:HB2	2.03	0.58
1:A:95:LEU:HD23	1:A:104:ALA:HB3	1.85	0.58
1:A:75:PRO:O	1:A:78:PHE:HB2	2.04	0.57
1:A:11:SER:OG	2:B:1954:ARG:NH2	2.35	0.57
1:A:178:ARG:HH11	1:A:178:ARG:CG	2.09	0.57
2:B:1960:VAL:O	2:B:1964:LEU:HG	2.05	0.57
1:A:44:VAL:HG11	1:A:89:VAL:HG22	1.87	0.57
1:A:88:LEU:CD2	1:A:111:GLY:HA3	2.33	0.57
1:A:202:LEU:HB3	1:A:203:PRO:CD	2.32	0.57
1:A:88:LEU:HD11	1:A:112:SER:N	2.20	0.57
1:A:16:VAL:HA	1:A:19:GLN:HB2	1.87	0.56
1:A:205:LEU:HG	1:A:209:MET:HE2	1.87	0.56
1:A:153:THR:CG2	1:A:155:GLU:H	2.19	0.56
2:B:1965:ALA:O	2:B:1968:GLN:HG3	2.05	0.56
1:A:5:HIS:HB2	1:A:124:LEU:HD21	1.87	0.56
1:A:54:LEU:HD11	2:B:1953:ALA:HB2	1.87	0.56
1:A:215:THR:HG23	1:A:221:GLN:HB3	1.88	0.56
1:A:219:LYS:HD2	1:A:219:LYS:C	2.25	0.55
1:A:94:MET:O	1:A:95:LEU:HD23	2.06	0.55
1:A:131:VAL:O	1:A:135:ILE:HD12	2.05	0.55
1:A:91:ALA:CA	1:A:94:MET:HG2	2.33	0.55
1:A:39:ASP:O	1:A:43:PRO:HG3	2.06	0.55
1:A:94:MET:HG3	1:A:104:ALA:CB	2.33	0.55
1:A:185:GLN:O	1:A:189:VAL:HG23	2.06	0.55
1:A:88:LEU:HD22	1:A:108:LEU:HD23	1.87	0.55
1:A:19:GLN:O	1:A:23:LEU:HB2	2.06	0.54
2:B:1947:LYS:O	2:B:1949:LEU:N	2.40	0.54
1:A:20:ILE:CG1	2:B:1964:LEU:HD13	2.38	0.54
1:A:75:PRO:HB2	1:A:76:PRO:HD3	1.89	0.54
1:A:44:VAL:HG11	1:A:89:VAL:HG13	1.90	0.54
1:A:179:GLN:HA	1:A:182:LEU:HD12	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:GLY:O	1:A:62:VAL:HG23	2.08	0.54
1:A:173:LYS:O	1:A:177:GLU:HG3	2.06	0.54
1:A:22:HIS:C	1:A:24:VAL:H	2.11	0.54
1:A:37:ILE:CA	1:A:40:LEU:HD12	2.09	0.53
1:A:88:LEU:HD23	1:A:107:TYR:O	2.08	0.53
1:A:35:LYS:HZ2	1:A:95:LEU:CB	2.22	0.53
1:A:133:LYS:NZ	2:B:1946:LYS:NZ	2.56	0.53
1:A:88:LEU:HD23	1:A:111:GLY:HA3	1.90	0.53
1:A:54:LEU:CG	2:B:1953:ALA:HB2	2.38	0.53
1:A:51:VAL:HG23	2:B:1956:VAL:HG13	1.90	0.53
1:A:218:SER:HB3	1:A:221:GLN:NE2	2.24	0.53
2:B:1948:GLU:HA	2:B:1951:GLU:CD	2.29	0.52
2:B:1954:ARG:O	2:B:1958:GLU:HG3	2.08	0.52
1:A:202:LEU:CB	1:A:203:PRO:HD3	2.31	0.52
1:A:35:LYS:HD3	1:A:95:LEU:CD1	2.40	0.52
1:A:35:LYS:HD3	1:A:95:LEU:HD13	1.92	0.52
1:A:95:LEU:CD2	1:A:104:ALA:HB3	2.39	0.52
1:A:91:ALA:HB3	1:A:108:LEU:HD23	1.90	0.51
1:A:218:SER:O	1:A:221:GLN:NE2	2.28	0.51
1:A:153:THR:HG23	1:A:154:MET:N	2.25	0.51
1:A:143:GLU:O	1:A:146:THR:HB	2.11	0.51
1:A:176:ASP:O	1:A:179:GLN:HG3	2.11	0.51
1:A:109:ILE:HG23	1:A:109:ILE:O	2.09	0.51
1:A:3:VAL:HG13	1:A:3:VAL:O	2.11	0.51
1:A:102:VAL:N	1:A:103:PRO:HD2	2.25	0.51
1:A:13:LEU:N	1:A:13:LEU:HD23	2.26	0.50
1:A:191:LEU:CD2	1:A:247:VAL:HG22	2.39	0.50
1:A:91:ALA:CB	1:A:108:LEU:HD23	2.41	0.50
1:A:154:MET:HE2	1:A:214:THR:CG2	2.39	0.50
1:A:104:ALA:C	1:A:106:ASP:H	2.14	0.50
1:A:215:THR:HG23	1:A:221:GLN:CB	2.41	0.50
1:A:35:LYS:NZ	1:A:95:LEU:HA	2.26	0.50
2:B:1945:THR:O	2:B:1948:GLU:HB2	2.11	0.50
1:A:205:LEU:HG	1:A:209:MET:CE	2.42	0.50
1:A:133:LYS:HZ3	2:B:1946:LYS:NZ	2.10	0.50
1:A:72:ARG:O	1:A:73:ASP:OD1	2.30	0.50
1:A:50:ALA:HB1	2:B:1956:VAL:HG22	1.94	0.50
1:A:1:MET:HB3	1:A:2:PRO:CD	2.42	0.49
1:A:168:MET:HG3	1:A:202:LEU:HD22	1.93	0.49
1:A:218:SER:C	1:A:221:GLN:HE21	2.13	0.49
1:A:48:GLN:HA	1:A:48:GLN:OE1	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:69:ILE:HD13	1:A:257:ALA:HB3	1.89	0.49
1:A:105:ARG:HH21	2:B:1967:LEU:CA	2.23	0.49
1:A:154:MET:CE	1:A:213:VAL:HG12	2.42	0.49
1:A:52:SER:O	1:A:56:ARG:HG3	2.12	0.49
1:A:50:ALA:CB	2:B:1956:VAL:HG22	2.42	0.49
1:A:116:LEU:HD23	1:A:116:LEU:N	2.27	0.49
1:A:91:ALA:HB3	1:A:108:LEU:HD21	1.94	0.49
1:A:20:ILE:O	1:A:20:ILE:HG22	2.12	0.49
1:A:24:VAL:HG12	1:A:25:ILE:HG13	1.95	0.48
1:A:58:GLY:O	1:A:78:PHE:HZ	1.95	0.48
1:A:6:THR:OG1	1:A:9:ILE:HG13	2.12	0.48
2:B:1945:THR:HA	2:B:1948:GLU:CD	2.33	0.48
1:A:219:LYS:HD2	1:A:220:ASN:CB	2.42	0.48
2:B:1947:LYS:HA	2:B:1950:ILE:HD13	1.96	0.48
1:A:154:MET:CE	1:A:214:THR:CG2	2.92	0.48
1:A:32:VAL:CG1	1:A:33:ASP:N	2.76	0.48
1:A:82:GLU:OE1	1:A:82:GLU:HA	2.13	0.48
1:A:43:PRO:O	1:A:47:VAL:HB	2.13	0.48
1:A:75:PRO:CD	1:A:76:PRO:HD2	2.44	0.48
1:A:95:LEU:HD23	1:A:104:ALA:HB1	1.93	0.48
1:A:94:MET:HE1	1:A:107:TYR:CD1	2.48	0.48
1:A:16:VAL:HG21	2:B:1957:SER:O	2.12	0.48
1:A:40:LEU:C	1:A:43:PRO:HD2	2.34	0.48
1:A:148:ALA:HB2	1:A:160:TYR:CZ	2.49	0.48
1:A:42:ALA:N	1:A:43:PRO:CD	2.77	0.48
1:A:127:ASP:O	1:A:131:VAL:HG23	2.14	0.47
1:A:30:GLY:O	1:A:31:GLU:C	2.52	0.47
1:A:212:PHE:O	1:A:216:LYS:HB2	2.14	0.47
1:A:105:ARG:NH2	2:B:1967:LEU:HA	2.26	0.47
1:A:81:VAL:HG22	1:A:115:ILE:O	2.14	0.47
1:A:44:VAL:HG12	1:A:44:VAL:O	2.14	0.47
1:A:35:LYS:HZ2	1:A:95:LEU:HD22	1.78	0.47
1:A:249:GLN:O	1:A:253:TRP:HB2	2.14	0.47
1:A:94:MET:HB2	1:A:104:ALA:CB	2.40	0.47
1:A:88:LEU:CD1	1:A:112:SER:HB3	2.35	0.47
1:A:35:LYS:HG3	1:A:95:LEU:HD13	1.96	0.47
1:A:64:THR:HG23	1:A:64:THR:O	2.15	0.47
1:A:153:THR:O	1:A:156:ASP:HB2	2.15	0.47
1:A:62:VAL:HG23	1:A:74:MET:HE3	1.96	0.47
1:A:153:THR:CG2	1:A:155:GLU:N	2.78	0.46
1:A:153:THR:CG2	1:A:154:MET:N	2.76	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:179:GLN:NE2	1:A:180:GLN:HG3	2.30	0.46
1:A:191:LEU:CD2	1:A:247:VAL:CG2	2.94	0.46
1:A:55:VAL:HG12	1:A:56:ARG:N	2.28	0.46
1:A:102:VAL:N	1:A:103:PRO:CD	2.78	0.46
1:A:16:VAL:HG13	2:B:1964:LEU:HD11	1.97	0.46
1:A:24:VAL:HA	1:A:27:HIS:HD2	1.80	0.46
1:A:94:MET:CE	1:A:107:TYR:CD1	2.99	0.46
1:A:4:PHE:CE2	1:A:13:LEU:HD12	2.47	0.46
1:A:243:GLU:O	1:A:247:VAL:HG13	2.16	0.46
1:A:218:SER:CB	1:A:221:GLN:NE2	2.79	0.46
1:A:37:ILE:N	1:A:38:PRO:CD	2.78	0.46
1:A:62:VAL:CG2	1:A:74:MET:HE3	2.46	0.46
1:A:35:LYS:NZ	1:A:95:LEU:CA	2.79	0.46
1:A:6:THR:HG22	1:A:181:GLU:O	2.15	0.45
1:A:70:LEU:HD23	1:A:129:ALA:HB2	1.98	0.45
2:B:1962:HIS:O	2:B:1965:ALA:HB3	2.16	0.45
1:A:84:ALA:CB	1:A:115:ILE:HA	2.45	0.45
1:A:94:MET:SD	1:A:107:TYR:CD1	3.09	0.45
1:A:9:ILE:HD13	1:A:123:LEU:CB	2.46	0.45
2:B:1954:ARG:HG2	2:B:1958:GLU:OE2	2.14	0.45
1:A:54:LEU:HA	1:A:54:LEU:HD23	1.71	0.45
1:A:1:MET:HA	1:A:2:PRO:HD3	1.83	0.45
1:A:20:ILE:HG12	2:B:1964:LEU:HD13	1.97	0.45
1:A:115:ILE:HG22	1:A:116:LEU:HD23	1.97	0.45
1:A:46:ALA:HB1	2:B:1959:LYS:HZ2	1.82	0.45
1:A:55:VAL:CG2	1:A:78:PHE:HD2	2.28	0.45
1:A:135:ILE:CG2	1:A:139:LYS:HD2	2.47	0.45
1:A:84:ALA:HB3	1:A:115:ILE:HG12	1.94	0.44
1:A:59:LYS:O	1:A:63:GLN:HG3	2.16	0.44
1:A:102:VAL:HB	1:A:103:PRO:CD	2.46	0.44
1:A:112:SER:O	1:A:115:ILE:HB	2.17	0.44
1:A:75:PRO:HB2	1:A:76:PRO:CD	2.48	0.44
1:A:35:LYS:CG	1:A:95:LEU:HD13	2.48	0.44
1:A:157:LEU:O	1:A:157:LEU:HD12	2.18	0.44
1:A:69:ILE:HG23	1:A:257:ALA:HB2	1.98	0.44
1:A:7:ARG:O	1:A:11:SER:HB3	2.17	0.44
1:A:16:VAL:HG12	1:A:16:VAL:O	2.17	0.44
1:A:79:ILE:O	1:A:83:ASN:ND2	2.51	0.44
1:A:97:SER:C	1:A:99:PRO:HD3	2.37	0.44
1:A:249:GLN:HG2	1:A:253:TRP:CE3	2.52	0.44
1:A:54:LEU:CD1	2:B:1953:ALA:HB2	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:VAL:HG13	1:A:248:LEU:CD2	2.48	0.44
1:A:210:LYS:O	1:A:214:THR:OG1	2.29	0.43
1:A:230:ARG:HE	1:A:230:ARG:HB3	1.76	0.43
1:A:8:THR:O	1:A:8:THR:HG22	2.19	0.43
1:A:179:GLN:C	1:A:181:GLU:H	2.22	0.43
1:A:35:LYS:HE3	1:A:99:PRO:HB3	1.99	0.43
1:A:94:MET:CG	1:A:104:ALA:CB	2.96	0.43
1:A:50:ALA:HA	1:A:53:ASN:CG	2.39	0.43
1:A:133:LYS:NZ	2:B:1946:LYS:HZ1	2.16	0.43
1:A:4:PHE:CZ	1:A:13:LEU:CB	2.97	0.43
1:A:21:SER:O	1:A:25:ILE:HD12	2.19	0.42
1:A:54:LEU:CD2	2:B:1949:LEU:HD12	2.49	0.42
1:A:86:THR:O	1:A:90:GLN:HB2	2.18	0.42
1:A:35:LYS:CE	1:A:95:LEU:HD13	2.48	0.42
1:A:213:VAL:HG13	1:A:213:VAL:O	2.19	0.42
2:B:1949:LEU:HA	2:B:1949:LEU:HD13	1.92	0.42
1:A:35:LYS:HZ2	1:A:95:LEU:CD2	2.33	0.42
1:A:50:ALA:C	2:B:1956:VAL:HG22	2.39	0.42
1:A:112:SER:O	1:A:115:ILE:N	2.49	0.42
1:A:74:MET:HB2	1:A:125:THR:HG21	2.01	0.42
1:A:186:GLU:HA	1:A:186:GLU:OE1	2.18	0.42
1:A:48:GLN:OE1	1:A:85:CYS:SG	2.77	0.42
1:A:47:VAL:HG23	2:B:1959:LYS:HB3	2.01	0.42
1:A:106:ASP:O	1:A:109:ILE:HG22	2.20	0.42
1:A:88:LEU:C	1:A:90:GLN:N	2.71	0.42
1:A:96:GLN:O	1:A:99:PRO:HD3	2.20	0.42
1:A:194:SER:OG	1:A:243:GLU:HG2	2.20	0.42
2:B:1967:LEU:CD2	2:B:1967:LEU:N	2.77	0.42
1:A:74:MET:HE3	1:A:74:MET:HB3	1.90	0.42
1:A:75:PRO:N	1:A:76:PRO:CD	2.80	0.41
1:A:135:ILE:HG22	1:A:139:LYS:HD2	2.01	0.41
1:A:35:LYS:HG3	1:A:95:LEU:CD1	2.50	0.41
1:A:4:PHE:CE2	1:A:13:LEU:CB	2.99	0.41
1:A:92:ALA:O	1:A:96:GLN:HB2	2.20	0.41
2:B:1965:ALA:O	2:B:1967:LEU:N	2.52	0.41
1:A:35:LYS:HZ1	1:A:95:LEU:HA	1.85	0.41
1:A:75:PRO:HD2	1:A:76:PRO:HD2	2.01	0.41
1:A:84:ALA:CB	1:A:115:ILE:CA	2.98	0.41
1:A:236:LYS:O	1:A:240:GLU:HG2	2.20	0.41
2:B:1960:VAL:CG1	2:B:1960:VAL:O	2.67	0.41
1:A:202:LEU:CB	1:A:203:PRO:CD	2.96	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:GLN:C	1:A:92:ALA:N	2.73	0.41
1:A:48:GLN:O	1:A:48:GLN:HG3	2.21	0.41
1:A:94:MET:HE3	1:A:94:MET:HB3	1.76	0.41
1:A:128:GLU:HG3	1:A:252:SER:OG	2.19	0.41
1:A:35:LYS:NZ	1:A:95:LEU:O	2.53	0.41
1:A:84:ALA:HB1	1:A:111:GLY:O	2.20	0.41
1:A:80:LYS:NZ	1:A:121:ASP:OD2	2.47	0.40
1:A:154:MET:CE	1:A:213:VAL:CG1	3.00	0.40
1:A:154:MET:HE1	1:A:213:VAL:HG12	2.03	0.40
1:A:38:PRO:C	1:A:40:LEU:H	2.24	0.40
1:A:85:CYS:SG	1:A:85:CYS:O	2.79	0.40
2:B:1945:THR:HG23	2:B:1948:GLU:CD	2.41	0.40
1:A:27:HIS:CE1	1:A:31:GLU:OE1	2.75	0.40
1:A:35:LYS:HZ1	1:A:95:LEU:CA	2.34	0.40
1:A:94:MET:CB	1:A:104:ALA:HB2	2.43	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	256/262 (98%)	214 (84%)	37 (14%)	5 (2%)	9	22
2	B	24/26 (92%)	19 (79%)	3 (12%)	2 (8%)	1	1
All	All	280/288 (97%)	233 (83%)	40 (14%)	7 (2%)	6	17

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	PRO
2	B	1948	GLU
2	B	1966	ALA

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Mol	Chain	Res	Type
1	A	23	LEU
1	A	180	GLN
1	A	219	LYS
1	A	97	SER

### 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	229/233 (98%)	179 (78%)	50 (22%)	1	3
2	B	22/22 (100%)	12 (54%)	10 (46%)	0	0
All	All	251/255 (98%)	191 (76%)	60 (24%)	1	2

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	THR
1	A	11	SER
1	A	18	GLN
1	A	24	VAL
1	A	35	LYS
1	A	52	SER
1	A	57	VAL
1	A	60	GLU
1	A	64	THR
1	A	69	ILE
1	A	70	LEU
1	A	74	MET
1	A	79	ILE
1	A	83	ASN
1	A	85	CYS
1	A	94	MET
1	A	95	LEU
1	A	105	ARG
1	A	106	ASP

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Mol	Chain	Res	Type
1	A	108	LEU
1	A	109	ILE
1	A	115	ILE
1	A	116	LEU
1	A	120	SER
1	A	122	LEU
1	A	124	LEU
1	A	130	GLU
1	A	133	LYS
1	A	139	LYS
1	A	145	LEU
1	A	146	THR
1	A	150	VAL
1	A	153	THR
1	A	155	GLU
1	A	157	LEU
1	A	173	LYS
1	A	178	ARG
1	A	181	GLU
1	A	213	VAL
1	A	215	THR
1	A	219	LYS
1	A	221	GLN
1	A	223	ILE
1	A	228	LYS
1	A	230	ARG
1	A	243	GLU
1	A	247	VAL
1	A	252	SER
1	A	254	ASP
1	A	258	TRP
2	B	1945	THR
2	B	1947	LYS
2	B	1949	LEU
2	B	1951	GLU
2	B	1955	LYS
2	B	1959	LYS
2	B	1963	VAL
2	B	1964	LEU
2	B	1967	LEU
2	B	1968	GLN

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	5	HIS
1	A	22	HIS
1	A	27	HIS
1	A	63	GLN
1	A	83	ASN
1	A	221	GLN

### 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 [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

### 6.4 Ligands

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers

EDS was not executed - this section is therefore empty.