



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 06:16 PM GMT

PDB ID : 1F13  
Title : RECOMBINANT HUMAN CELLULAR COAGULATION FACTOR XIII  
Authors : Weiss, M.S.; Hilgenfeld, R.  
Deposited on : 1998-01-16  
Resolution : 2.10 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

---

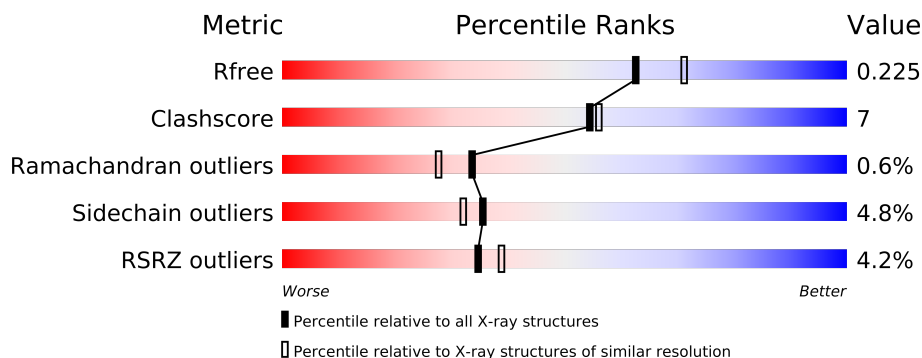
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.10 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))
$R_{free}$	66092	3012 (2.10-2.10)
Clashscore	79885	3649 (2.10-2.10)
Ramachandran outliers	78287	3610 (2.10-2.10)
Sidechain outliers	78261	3611 (2.10-2.10)
RSRZ outliers	66119	3013 (2.10-2.10)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	731	
1	B	731	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12043 atoms, of which 0 are hydrogen and 0 are deuterium.

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 CELLULAR COAGULATION FACTOR XIII ZYMOGEN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	722	Total	C	N	O	S	0	0	0
			5791	3670	1001	1093	27			
1	B	719	Total	C	N	O	S	0	0	0
			5765	3655	994	1089	27			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	651	GLU	GLN	CONFLICT	UNP P00488
B	651	GLU	GLN	CONFLICT	UNP P00488

- Molecule 2 is water.

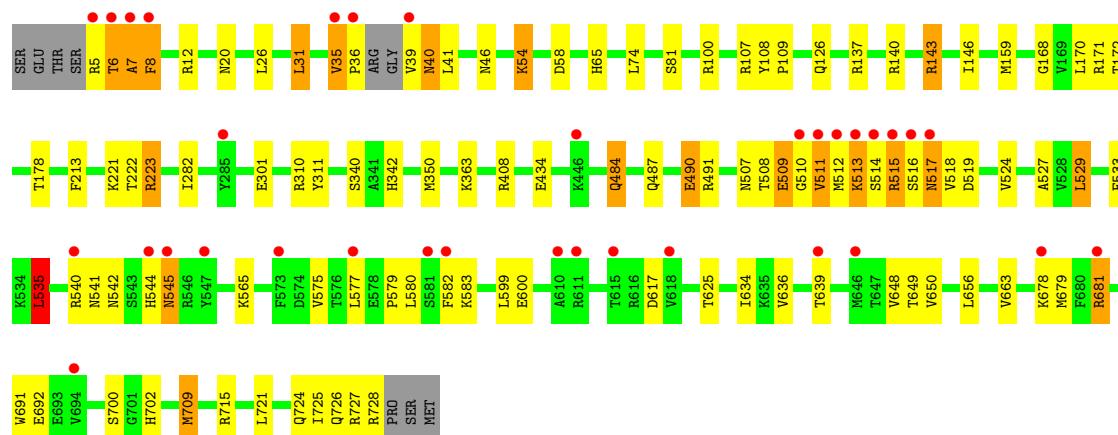
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	263	Total	O	0	0
			263	263		
2	B	224	Total	O	0	0
			224	224		

### 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 errors 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.

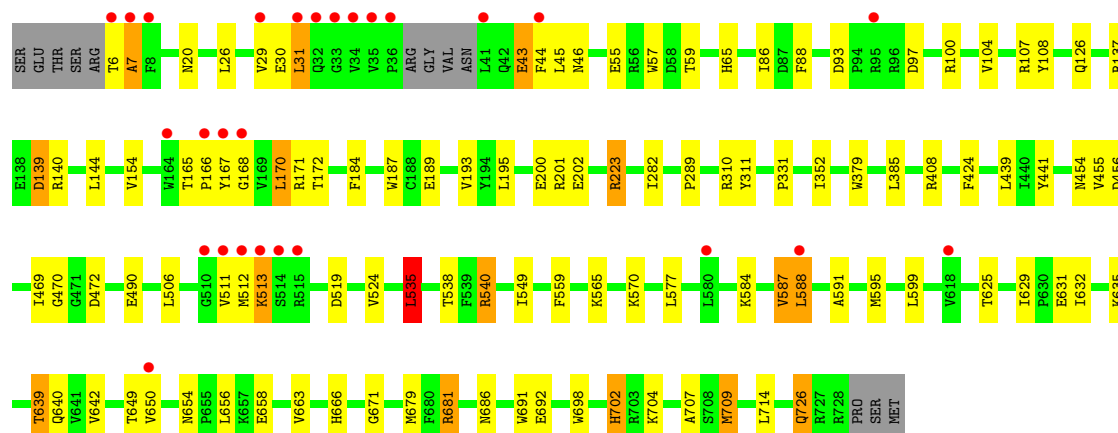
#### • Molecule 1: CELLULAR COAGULATION FACTOR XIII ZYMOGEN

Chain A: 



#### • Molecule 1: CELLULAR COAGULATION FACTOR XIII ZYMOGEN

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	134.59Å 72.78Å 101.05Å 90.00° 106.08° 90.00°	Depositor
Resolution (Å)	40.00 – 2.10 58.24 – 2.10	Depositor EDS
% Data completeness (in resolution range)	82.0 (40.00-2.10) 81.6 (58.24-2.10)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.51 (at 2.10Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, $R_{free}$	0.183 , 0.236 0.177 , 0.225	Depositor DCC
$R_{free}$ test set	1795 reflections (2.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	32.9	Xtriage
Anisotropy	0.395	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 41.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 89656 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	12043	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.63% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.53	0/5926	0.75	1/8041 (0.0%)
1	B	0.51	0/5900	0.73	1/8006 (0.0%)
All	All	0.52	0/11826	0.74	2/16047 (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	A	535	LEU	CA-CB-CG	6.64	130.56	115.30
1	B	535	LEU	CA-CB-CG	6.06	129.23	115.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5791	0	5656	85	0
1	B	5765	0	5628	77	0
2	A	263	0	0	9	0
2	B	224	0	0	8	0
All	All	12043	0	11284	155	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (155) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:538:THR:HG22	1:B:584:LYS:HG2	1.41	0.99
1:B:535:LEU:HB3	1:B:587:VAL:HG22	1.51	0.91
1:A:515:ARG:HA	1:A:515:ARG:NE	1.89	0.87
1:A:40:ASN:HD22	1:A:40:ASN:H	1.18	0.87
1:A:143:ARG:HB2	1:A:143:ARG:HH11	1.48	0.79
1:A:575:VAL:HG13	1:A:583:LYS:HD3	1.67	0.76
1:B:681:ARG:NH1	1:B:681:ARG:HA	2.01	0.76
1:B:649:THR:HG22	1:B:692:GLU:HG2	1.68	0.75
1:A:524:VAL:HG22	1:A:535:LEU:HD22	1.69	0.73
1:A:636:VAL:HG12	1:A:648:VAL:HG22	1.73	0.71
1:B:524:VAL:HG22	1:B:535:LEU:HD22	1.73	0.70
1:A:507:ASN:OD1	1:A:509:GLU:HG3	1.91	0.70
1:B:513:LYS:NZ	1:B:513:LYS:HA	2.07	0.69
1:A:511:VAL:C	1:A:513:LYS:H	1.94	0.69
1:A:40:ASN:ND2	1:A:40:ASN:H	1.90	0.69
1:A:512:MET:C	1:A:514:SER:H	1.96	0.68
1:B:588:LEU:HB3	2:B:2137:HOH:O	1.93	0.67
1:B:642:VAL:HG23	1:B:726:GLN:O	1.94	0.67
1:A:159:MET:SD	2:A:2189:HOH:O	2.53	0.66
1:B:519:ASP:HB2	1:B:540:ARG:HB3	1.78	0.66
1:A:681:ARG:HA	1:A:681:ARG:NH1	2.10	0.65
1:B:331:PRO:HB2	1:B:379:TRP:HB3	1.78	0.65
1:B:513:LYS:HA	1:B:513:LYS:HZ3	1.62	0.64
1:B:635:LYS:HB2	1:B:649:THR:OG1	1.97	0.64
1:A:35:VAL:HB	1:A:36:PRO:C	2.17	0.64
1:A:39:VAL:HG12	1:A:40:ASN:H	1.63	0.63
1:B:202:GLU:HG3	2:B:2212:HOH:O	2.01	0.59
1:B:137:ARG:HD2	2:B:1921:HOH:O	2.02	0.58
1:A:515:ARG:CZ	1:A:515:ARG:HA	2.34	0.58
1:A:529:LEU:HD21	1:A:656:LEU:HD21	1.86	0.57
1:A:727:ARG:O	1:A:728:ARG:HB2	2.05	0.57
1:A:511:VAL:C	1:A:513:LYS:N	2.58	0.56
1:B:663:VAL:HG13	1:B:709:MET:HE2	1.86	0.56
1:A:213:PHE:CD1	1:A:222:THR:HG22	2.40	0.56
1:B:189:GLU:HG3	2:B:2216:HOH:O	2.05	0.56
1:B:107:ARG:HD2	1:B:108:TYR:CZ	2.41	0.56
1:A:484:GLN:HB3	1:A:487:GLN:HE21	1.71	0.56
1:B:43:GLU:O	1:B:43:GLU:HG2	2.05	0.55
1:A:513:LYS:C	1:A:515:ARG:H	2.09	0.55
1:A:39:VAL:HG12	1:A:40:ASN:N	2.22	0.55
1:A:524:VAL:HG22	1:A:535:LEU:CD2	2.36	0.55
1:A:510:GLY:C	1:A:512:MET:H	2.10	0.55

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:223:ARG:NH2	2:A:857:HOH:O	2.40	0.55
1:A:178:THR:HB	2:A:2189:HOH:O	2.07	0.54
1:A:8:PHE:HE2	1:B:559:PHE:CD1	2.26	0.54
1:A:649:THR:HG22	1:A:692:GLU:HG3	1.90	0.54
1:B:570:LYS:HE2	2:B:2264:HOH:O	2.07	0.54
1:B:663:VAL:HG13	1:B:709:MET:CE	2.38	0.54
1:B:44:PHE:CZ	1:B:166:PRO:HD2	2.43	0.54
1:A:282:ILE:HD11	1:B:7:ALA:N	2.23	0.53
2:A:2248:HOH:O	1:B:282:ILE:HD12	2.09	0.52
1:A:508:THR:O	1:A:511:VAL:HG12	2.09	0.52
1:B:441:TYR:HD2	1:B:454:ASN:HD22	1.57	0.52
1:A:515:ARG:HB3	1:A:617:ASP:OD2	2.10	0.52
1:A:517:ASN:ND2	1:A:517:ASN:H	2.08	0.52
1:A:663:VAL:HG13	1:A:709:MET:HE2	1.91	0.52
1:B:26:LEU:HD11	1:B:104:VAL:HG21	1.91	0.52
1:A:575:VAL:CG1	1:A:583:LYS:HD3	2.40	0.51
1:A:663:VAL:HG13	1:A:709:MET:CE	2.40	0.51
1:A:5:ARG:HG3	1:A:5:ARG:HH11	1.76	0.51
1:A:519:ASP:HB2	1:A:540:ARG:HB3	1.93	0.51
1:A:143:ARG:CB	1:A:143:ARG:HH11	2.18	0.51
1:A:35:VAL:HB	1:A:36:PRO:CA	2.41	0.51
1:A:527:ALA:HB2	1:A:533:PHE:HB3	1.93	0.50
1:B:45:LEU:HD22	1:B:88:PHE:HB3	1.93	0.50
1:B:30:GLU:O	1:B:168:GLY:HA3	2.12	0.49
1:B:86:ILE:CD1	1:B:144:LEU:HD12	2.42	0.49
1:A:107:ARG:HD2	1:A:108:TYR:CZ	2.48	0.49
1:A:301:GLU:HG2	2:A:2273:HOH:O	2.13	0.49
1:B:511:VAL:HG12	1:B:512:MET:H	1.77	0.49
1:A:81:SER:HA	1:A:146:ILE:O	2.13	0.48
1:B:107:ARG:HG2	1:B:107:ARG:O	2.13	0.48
1:B:45:LEU:HD22	1:B:88:PHE:CG	2.49	0.48
1:B:629:ILE:HD13	1:B:714:LEU:HD11	1.96	0.48
1:A:221:LYS:HG3	2:A:2266:HOH:O	2.13	0.47
1:B:588:LEU:N	1:B:588:LEU:HD22	2.29	0.47
1:B:385:LEU:HD22	1:B:424:PHE:HB3	1.96	0.47
1:A:7:ALA:N	1:B:282:ILE:HD11	2.30	0.47
1:A:678:LYS:HB2	1:A:691:TRP:CE2	2.49	0.46
1:A:545:ASN:C	1:A:579:PRO:HG3	2.36	0.46
1:B:649:THR:HG22	1:B:692:GLU:CG	2.43	0.46
1:B:681:ARG:HA	1:B:681:ARG:CZ	2.45	0.46
1:A:8:PHE:CE2	1:B:559:PHE:CD1	3.03	0.46
1:B:439:LEU:HB2	1:B:456:ASP:HB3	1.97	0.46

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:40:ASN:N	1:A:40:ASN:HD22	1.98	0.46
1:A:600:GLU:OE1	1:A:715:ARG:HD2	2.16	0.46
1:B:352:ILE:HD11	1:B:439:LEU:HD22	1.98	0.46
1:A:512:MET:C	1:A:514:SER:N	2.67	0.45
1:B:154:VAL:HG21	1:B:184:PHE:CE2	2.51	0.45
1:A:529:LEU:HD21	1:A:656:LEU:CD2	2.46	0.45
1:B:55:GLU:HG3	1:B:57:TRP:CZ2	2.52	0.45
1:A:8:PHE:HE2	1:B:559:PHE:HD1	1.64	0.45
1:B:223:ARG:NH2	2:B:1857:HOH:O	2.49	0.45
1:B:65:HIS:HD2	2:B:1810:HOH:O	1.99	0.45
1:A:517:ASN:OD1	1:A:518:VAL:HG23	2.17	0.45
1:B:93:ASP:O	1:B:97:ASP:HB2	2.17	0.45
1:A:8:PHE:CE2	1:B:559:PHE:HD1	2.35	0.44
1:B:31:LEU:HA	1:B:167:TYR:O	2.16	0.44
1:B:455:VAL:HB	1:B:512:MET:HG2	1.98	0.44
1:A:171:ARG:HG2	1:A:172:THR:O	2.17	0.44
1:A:540:ARG:HD2	1:A:582:PHE:CB	2.47	0.44
1:B:139:ASP:HB3	1:B:140:ARG:H	1.49	0.44
1:B:549:ILE:HD11	1:B:577:LEU:HD12	2.00	0.44
1:B:55:GLU:HG2	2:B:2103:HOH:O	2.17	0.44
1:B:588:LEU:H	1:B:588:LEU:HD22	1.83	0.44
1:A:509:GLU:CD	1:A:510:GLY:H	2.22	0.44
1:B:29:VAL:HG12	1:B:31:LEU:HD12	2.00	0.43
1:B:591:ALA:O	1:B:595:MET:HG2	2.18	0.43
1:B:6:THR:O	1:B:7:ALA:HB3	2.19	0.43
1:A:31:LEU:HD12	1:A:168:GLY:HA3	2.01	0.43
1:B:654:ASN:O	1:B:686:ASN:HA	2.19	0.43
1:A:213:PHE:HD1	1:A:222:THR:HG22	1.83	0.43
1:A:540:ARG:NH1	1:A:582:PHE:HB3	2.34	0.43
1:B:666:HIS:O	1:B:707:ALA:HA	2.19	0.43
1:B:656:LEU:C	1:B:658:GLU:H	2.22	0.43
1:A:137:ARG:HD2	2:A:2315:HOH:O	2.18	0.43
1:B:184:PHE:HB3	1:B:193:VAL:HG21	1.99	0.43
1:A:310:ARG:HA	1:A:311:TYR:HA	1.83	0.43
1:B:565:LYS:HD2	1:B:599:LEU:HD21	2.01	0.43
1:B:632:ILE:HD11	1:B:709:MET:HB2	2.00	0.42
1:B:472:ASP:OD2	1:B:704:LYS:NZ	2.52	0.42
1:A:650:VAL:HB	1:A:691:TRP:HB3	2.01	0.42
1:A:577:LEU:HD23	1:A:577:LEU:H	1.83	0.42
1:B:31:LEU:H	1:B:31:LEU:HD13	1.83	0.42
1:A:12:ARG:HD2	2:A:2251:HOH:O	2.18	0.42
1:A:517:ASN:O	1:A:542:ASN:HB2	2.19	0.42

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:107:ARG:C	1:A:109:PRO:HD3	2.40	0.42
1:A:342:HIS:ND1	1:A:434:GLU:OE2	2.43	0.42
1:A:490:GLU:HG3	1:A:491:ARG:N	2.34	0.42
1:A:65:HIS:HD2	2:A:810:HOH:O	2.01	0.42
1:A:514:SER:C	1:A:516:SER:H	2.23	0.42
1:B:310:ARG:HA	1:B:311:TYR:HA	1.81	0.42
1:A:544:HIS:CD2	1:A:544:HIS:H	2.38	0.41
1:A:700:SER:HA	1:A:725:ILE:HB	2.02	0.41
1:B:511:VAL:HG12	1:B:512:MET:N	2.35	0.41
1:A:515:ARG:HH12	1:A:517:ASN:HB3	1.84	0.41
1:A:580:LEU:N	1:A:580:LEU:HD22	2.34	0.41
1:A:541:ASN:O	1:A:580:LEU:HA	2.21	0.41
1:B:671:GLY:HA2	1:B:698:TRP:NE1	2.36	0.41
1:B:171:ARG:HG2	1:B:172:THR:O	2.20	0.41
1:A:282:ILE:HD12	1:B:7:ALA:HB2	2.03	0.41
1:B:650:VAL:HB	1:B:691:TRP:HB3	2.02	0.41
1:B:187:TRP:CE2	1:B:201:ARG:HD3	2.56	0.41
1:A:143:ARG:CG	1:A:143:ARG:HH11	2.33	0.41
1:A:634:ILE:O	1:A:721:LEU:HD22	2.21	0.41
1:A:565:LYS:HD2	1:A:599:LEU:HD21	2.02	0.41
1:A:727:ARG:NH1	1:A:727:ARG:HB3	2.35	0.41
1:A:54:LYS:HB3	1:A:74:LEU:HD13	2.03	0.41
1:B:639:THR:O	1:B:640:GLN:HB2	2.21	0.41
1:B:702:HIS:C	1:B:702:HIS:ND1	2.74	0.40
1:A:6:THR:O	1:A:7:ALA:HB2	2.20	0.40
1:B:165:THR:CG2	1:B:170:LEU:HD22	2.51	0.40
1:B:200:GLU:HG2	1:B:469:ILE:HD11	2.03	0.40
1:B:31:LEU:CD1	1:B:31:LEU:H	2.34	0.40
1:A:724:GLN:HE21	1:A:724:GLN:HB2	1.70	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	718/731 (98%)	680 (95%)	32 (4%)	6 (1%)	27	20
1	B	715/731 (98%)	679 (95%)	34 (5%)	2 (0%)	50	49
All	All	1433/1462 (98%)	1359 (95%)	66 (5%)	8 (1%)	33	28

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	8	PHE
1	A	35	VAL
1	A	6	THR
1	A	7	ALA
1	B	470	GLY
1	B	7	ALA
1	A	513	LYS
1	A	511	VAL

### 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	636/644 (99%)	603 (95%)	33 (5%)	32	29
1	B	633/644 (98%)	605 (96%)	28 (4%)	39	36
All	All	1269/1288 (98%)	1208 (95%)	61 (5%)	35	32

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	ASN
1	A	26	LEU
1	A	31	LEU
1	A	40	ASN
1	A	41	LEU
1	A	46	ASN
1	A	54	LYS
1	A	58	ASP
1	A	100	ARG

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	126	GLN
1	A	140	ARG
1	A	143	ARG
1	A	170	LEU
1	A	223	ARG
1	A	340	SER
1	A	350	MET
1	A	363	LYS
1	A	408	ARG
1	A	484	GLN
1	A	490	GLU
1	A	509	GLU
1	A	515	ARG
1	A	517	ASN
1	A	529	LEU
1	A	535	LEU
1	A	545	ASN
1	A	625	THR
1	A	639	THR
1	A	679	MET
1	A	681	ARG
1	A	702	HIS
1	A	709	MET
1	A	726	GLN
1	B	20	ASN
1	B	31	LEU
1	B	43	GLU
1	B	46	ASN
1	B	59	THR
1	B	100	ARG
1	B	126	GLN
1	B	139	ASP
1	B	170	LEU
1	B	195	LEU
1	B	223	ARG
1	B	289	PRO
1	B	408	ARG
1	B	490	GLU
1	B	506	LEU
1	B	513	LYS
1	B	535	LEU
1	B	540	ARG

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	B	587	VAL
1	B	588	LEU
1	B	625	THR
1	B	631	GLU
1	B	639	THR
1	B	679	MET
1	B	681	ARG
1	B	702	HIS
1	B	709	MET
1	B	726	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	65	HIS
1	A	126	GLN
1	A	484	GLN
1	A	487	GLN
1	A	526	ASN
1	A	544	HIS
1	A	640	GLN
1	A	724	GLN
1	B	65	HIS
1	B	110	GLN
1	B	126	GLN
1	B	267	ASN
1	B	436	ASN
1	B	507	ASN
1	B	526	ASN
1	B	724	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

There are no ligands in this entry.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 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	722/731 (98%)	0.12	34 (4%) 30 33	16, 34, 72, 96	0
1	B	719/731 (98%)	-0.01	27 (3%) 38 43	19, 37, 68, 92	0
All	All	1441/1462 (98%)	0.05	61 (4%) 35 39	16, 35, 71, 96	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	34	VAL	10.2
1	B	33	GLY	10.0
1	B	36	PRO	9.0
1	B	35	VAL	8.3
1	B	6	THR	8.2
1	B	512	MET	7.4
1	B	7	ALA	6.9
1	A	6	THR	6.5
1	B	31	LEU	6.3
1	A	516	SER	6.0
1	A	512	MET	5.4
1	B	167	TYR	5.1
1	B	41	LEU	5.0
1	A	7	ALA	4.9
1	B	511	VAL	4.8
1	A	514	SER	4.7
1	A	513	LYS	4.4
1	A	515	ARG	4.2
1	A	577	LEU	4.1
1	A	5	ARG	4.1
1	B	514	SER	4.0
1	A	582	PHE	4.0
1	A	544	HIS	3.9
1	A	39	VAL	3.7

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	511	VAL	3.7
1	B	32	GLN	3.7
1	A	8	PHE	3.6
1	B	164	TRP	3.6
1	B	8	PHE	3.6
1	A	547	TYR	3.6
1	B	513	LYS	3.3
1	A	681	ARG	3.3
1	A	540	ARG	3.1
1	B	515	ARG	3.1
1	A	639	THR	3.0
1	A	694	VAL	2.9
1	A	611	ARG	2.8
1	B	168	GLY	2.7
1	A	615	THR	2.7
1	B	44	PHE	2.6
1	B	650	VAL	2.5
1	A	285	TYR	2.5
1	A	35	VAL	2.5
1	B	166	PRO	2.4
1	A	678	LYS	2.4
1	B	580	LEU	2.3
1	A	545	ASN	2.3
1	B	588	LEU	2.3
1	A	36	PRO	2.3
1	A	517	ASN	2.3
1	A	581	SER	2.2
1	B	29	VAL	2.2
1	A	618	VAL	2.2
1	A	510	GLY	2.2
1	A	610	ALA	2.2
1	B	510	GLY	2.1
1	B	618	VAL	2.1
1	A	573	PHE	2.1
1	A	446	LYS	2.1
1	A	646	MET	2.0
1	B	95	ARG	2.0

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

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



### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 6.4 Ligands ⓘ

There are no ligands in this entry.

### 6.5 Other polymers ⓘ

There are no such residues in this entry.