



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 07:04 PM UTC

PDB ID : 2VXP / pdb\_00002vxp  
Title : The fourth FAS1 domain structure of human Bigh3  
Authors : Yoo, J.-H.; Cho, H.-S.  
Deposited on : 2008-07-08  
Resolution : 2.50 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  
Xtrriage (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

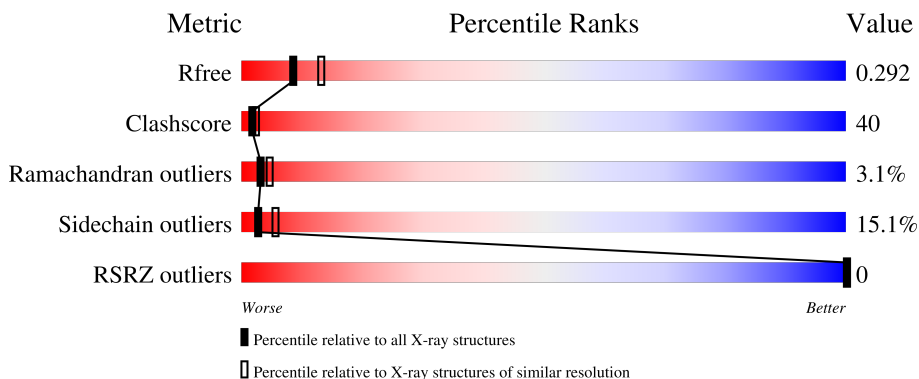
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	132	 37% 52% 9% ..
1	B	132	 34% 41% 22% ..

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2010 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 TRANSFORMING GROWTH FACTOR-BETA-INDUCED PROTEIN IG-H3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	131	989	623	174	188	4	0	0	0
1	B	131	989	623	174	188	4	0	0	0

- Molecule 2 is water.

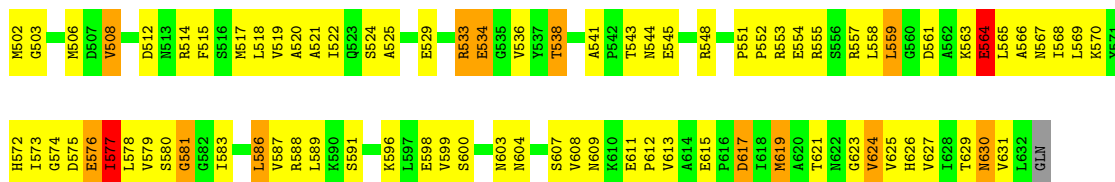
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	17	17	17	0	0
2	B	15	15	15	0	0

### 3 Residue-property plots

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.

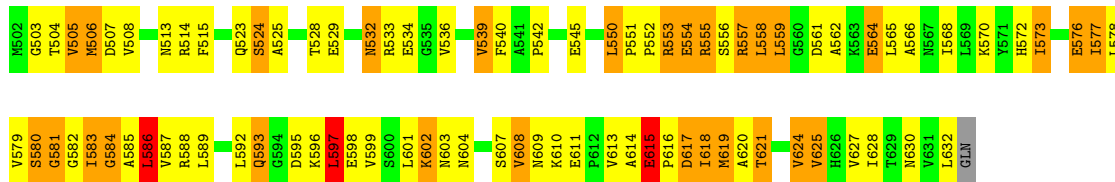
- Molecule 1: TRANSFORMING GROWTH FACTOR-BETA-INDUCED PROTEIN IG-H3

Chain A: 



- Molecule 1: TRANSFORMING GROWTH FACTOR-BETA-INDUCED PROTEIN IG-H3

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	62.93Å 62.93Å 143.29Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.50 50.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.0 (50.00-2.50) 98.1 (50.00-2.50)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.58 (at 2.51Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.270 , 0.285 0.273 , 0.292	Depositor DCC
$R_{free}$ test set	1081 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	36.7	Xtrriage
Anisotropy	0.164	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 25.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.40$ , $\langle L^2 \rangle = 0.23$	Xtrriage
Estimated twinning fraction	0.428 for h,-h-k,-l	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	2010	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.11% 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	0.58	0/1000	1.25	13/1354 (1.0%)
1	B	0.59	0/1000	1.46	27/1354 (2.0%)
All	All	0.58	0/2000	1.36	40/2708 (1.5%)

There are no bond length outliers.

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	508	VAL	N-CA-C	-9.55	98.64	111.44
1	B	505	VAL	N-CA-C	-8.98	101.46	110.62
1	B	615	GLU	CA-C-N	8.66	130.67	119.84
1	B	615	GLU	C-N-CA	8.66	130.67	119.84
1	B	532	ASN	N-CA-C	-7.92	103.19	113.17
1	B	524	SER	N-CA-C	-7.46	103.98	113.23
1	B	551	PRO	N-CA-C	-7.39	103.37	110.47
1	B	534	GLU	N-CA-C	7.36	120.36	110.35
1	B	617	ASP	N-CA-C	7.18	121.44	111.74
1	B	593	GLN	N-CA-C	-6.74	104.02	111.36
1	A	573	ILE	N-CA-C	6.57	117.30	107.51
1	B	625	VAL	N-CA-C	6.57	117.37	108.17
1	A	581	GLY	N-CA-C	-6.51	97.75	113.18
1	B	586	LEU	N-CA-C	-6.45	98.86	108.99
1	B	559	LEU	N-CA-C	6.22	120.61	112.89
1	B	507	ASP	N-CA-C	-6.13	105.76	113.18
1	B	597	LEU	CA-CB-CG	6.13	137.74	116.30
1	B	608	VAL	CB-CA-C	-6.06	103.98	111.32
1	B	552	PRO	N-CA-C	5.95	121.80	113.53
1	B	528	THR	N-CA-C	5.57	117.81	111.02
1	B	551	PRO	CA-C-N	5.57	125.23	119.05
1	B	551	PRO	C-N-CA	5.57	125.23	119.05
1	B	581	GLY	N-CA-C	-5.52	100.10	113.18
1	A	552	PRO	N-CA-C	5.50	123.81	112.47
1	B	576	GLU	N-CA-C	5.50	116.63	108.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	630	ASN	N-CA-C	-5.44	100.25	108.52
1	A	596	LYS	N-CA-C	5.37	117.46	110.43
1	A	587	VAL	N-CA-C	5.33	116.98	108.87
1	B	508	VAL	N-CA-C	-5.29	105.69	110.82
1	A	619	MET	N-CA-C	5.27	117.33	110.53
1	B	525	ALA	N-CA-C	-5.26	106.37	112.89
1	A	577	ILE	N-CA-C	-5.21	100.87	108.17
1	B	621	THR	N-CA-C	-5.18	106.47	112.89
1	A	525	ALA	N-CA-C	-5.17	106.97	113.28
1	B	555	ARG	N-CA-C	-5.17	105.65	111.28
1	A	520	ALA	N-CA-C	-5.13	104.95	111.11
1	B	506	MET	N-CA-C	5.11	117.59	111.71
1	A	591	SER	N-CA-C	-5.05	103.73	110.55
1	A	541	ALA	N-CA-C	5.01	116.65	109.04
1	B	573	ILE	CB-CA-C	-5.01	105.26	111.32

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	989	0	1028	64	0
1	B	989	0	1028	98	0
2	A	17	0	0	4	0
2	B	15	0	0	3	0
All	All	2010	0	2056	162	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 40.

All (162) 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:578:LEU:HD11	1:A:583:ILE:HD11	1.16	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:505:VAL:HG11	1:B:539:VAL:HG11	1.38	1.04
1:A:557:ARG:O	1:A:564:GLU:HG3	1.62	0.97
1:B:503:GLY:HA2	1:B:621:THR:H	1.31	0.92
1:A:578:LEU:HD11	1:A:583:ILE:CD1	2.03	0.88
1:B:539:VAL:HB	1:B:573:ILE:HD13	1.58	0.86
1:B:599:VAL:HA	1:B:607:SER:O	1.81	0.80
1:B:618:ILE:HG12	1:B:625:VAL:O	1.80	0.80
1:B:504:THR:HG23	1:B:532:ASN:HA	1.62	0.80
1:A:518:LEU:O	1:A:522:ILE:HG12	1.83	0.79
1:A:536:VAL:HB	1:A:577:ILE:HD12	1.63	0.79
1:B:553:ARG:HH11	1:B:553:ARG:HG3	1.48	0.78
1:A:617:ASP:HB3	1:A:626:HIS:CE1	2.19	0.77
1:A:538:THR:HG23	1:A:574:GLY:HA3	1.66	0.76
1:B:503:GLY:HA2	1:B:621:THR:N	2.01	0.76
1:B:557:ARG:HG3	1:B:557:ARG:HH11	1.51	0.76
1:A:577:ILE:H	1:A:577:ILE:HD13	1.50	0.75
1:B:504:THR:HG21	1:B:532:ASN:OD1	1.87	0.74
1:B:583:ILE:HB	1:B:587:VAL:HB	1.70	0.74
1:B:582:GLY:O	1:B:583:ILE:HG12	1.88	0.73
1:B:578:LEU:HD21	1:B:583:ILE:HG21	1.71	0.73
1:B:586:LEU:C	1:B:586:LEU:HD13	2.14	0.72
1:A:551:PRO:HG2	1:A:554:GLU:HB3	1.72	0.72
1:B:529:GLU:O	1:B:533:ARG:HB2	1.90	0.71
1:A:538:THR:HG21	1:A:589:LEU:HD13	1.71	0.71
1:B:588:ARG:HB3	1:B:596:LYS:HG2	1.71	0.71
1:B:553:ARG:HH11	1:B:553:ARG:CG	2.02	0.71
1:A:538:THR:HG21	1:A:589:LEU:CD1	2.21	0.70
1:B:555:ARG:HG2	1:B:559:LEU:HD22	1.73	0.70
1:B:555:ARG:O	1:B:559:LEU:HB2	1.92	0.70
1:A:604:ASN:ND2	2:A:2012:HOH:O	2.26	0.69
1:B:539:VAL:HB	1:B:573:ILE:CD1	2.23	0.69
1:B:566:ALA:O	1:B:570:LYS:HG2	1.94	0.68
1:B:611:GLU:OE1	1:B:630:ASN:ND2	2.25	0.68
1:A:553:ARG:O	1:A:557:ARG:HG3	1.95	0.66
1:A:555:ARG:O	1:A:559:LEU:HG	1.94	0.66
1:B:619:MET:HA	1:B:619:MET:HE2	1.78	0.66
1:A:577:ILE:HD13	1:A:577:ILE:N	2.11	0.66
1:B:576:GLU:CG	1:B:589:LEU:HD11	2.26	0.65
1:B:503:GLY:H	1:B:620:ALA:HA	1.61	0.65
1:A:608:VAL:O	1:A:608:VAL:HG13	1.97	0.64
1:B:570:LYS:HB2	1:B:592:LEU:HB2	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:578:LEU:HD11	1:B:587:VAL:HG11	1.79	0.64
1:A:617:ASP:HB3	1:A:626:HIS:HE1	1.62	0.64
1:A:588:ARG:HD2	1:A:598:GLU:OE1	1.97	0.63
1:A:561:ASP:HB3	1:A:564:GLU:HB2	1.81	0.62
1:B:503:GLY:CA	1:B:621:THR:HG23	2.30	0.62
1:B:505:VAL:HG11	1:B:539:VAL:CG1	2.22	0.62
1:B:578:LEU:HD21	1:B:583:ILE:CG2	2.30	0.62
1:A:568:ILE:HG23	1:A:572:HIS:HE1	1.65	0.61
1:A:517:MET:H	1:A:544:ASN:HD21	1.50	0.59
1:B:523:GLN:O	1:B:523:GLN:NE2	2.33	0.59
1:B:540:PHE:CE1	1:B:597:LEU:HD21	2.37	0.59
1:B:503:GLY:HA2	1:B:621:THR:HG23	1.85	0.59
1:A:545:GLU:HA	1:A:548:ARG:HB2	1.86	0.56
1:A:566:ALA:O	1:A:570:LYS:HB2	2.05	0.56
1:A:630:ASN:HA	2:A:2007:HOH:O	2.06	0.55
1:B:504:THR:CG2	1:B:532:ASN:HA	2.34	0.55
1:B:554:GLU:HG3	1:B:558:LEU:HD23	1.89	0.55
1:A:538:THR:CG2	1:A:574:GLY:HA3	2.35	0.55
1:A:506:MET:HE1	1:A:519:VAL:HG13	1.87	0.55
1:A:514:ARG:NH2	2:A:2004:HOH:O	2.40	0.55
1:B:614:ALA:O	1:B:615:GLU:O	2.25	0.55
1:A:577:ILE:HG22	1:A:623:GLY:HA2	1.89	0.55
1:A:538:THR:HG22	1:A:576:GLU:O	2.07	0.54
1:B:580:SER:O	1:B:581:GLY:C	2.49	0.54
1:B:613:VAL:HG11	1:B:616:PRO:HB3	1.89	0.54
1:B:586:LEU:HD13	1:B:587:VAL:N	2.23	0.53
1:B:582:GLY:C	1:B:583:ILE:HG23	2.33	0.53
1:B:618:ILE:N	1:B:618:ILE:HD13	2.22	0.53
1:B:586:LEU:C	1:B:586:LEU:CD1	2.81	0.53
1:B:593:GLN:NE2	1:B:595:ASP:OD2	2.39	0.52
1:A:579:VAL:HG23	1:A:581:GLY:O	2.09	0.52
1:B:578:LEU:HD23	1:B:579:VAL:N	2.25	0.52
1:A:536:VAL:CB	1:A:577:ILE:HD12	2.37	0.52
1:B:505:VAL:HG13	1:B:625:VAL:HG22	1.92	0.51
1:A:603:ASN:C	1:A:604:ASN:HD22	2.19	0.51
1:A:617:ASP:HB2	1:A:624:VAL:HG21	1.93	0.51
1:B:568:ILE:HG12	1:B:572:HIS:HE1	1.77	0.50
1:B:611:GLU:CD	1:B:630:ASN:HB2	2.36	0.50
1:A:536:VAL:HB	1:A:577:ILE:CD1	2.36	0.50
1:B:576:GLU:CD	1:B:589:LEU:HD11	2.36	0.50
1:A:521:ALA:HB1	1:A:569:LEU:HD11	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:609:ASN:HD22	1:B:609:ASN:N	2.09	0.50
1:B:553:ARG:CG	1:B:553:ARG:NH1	2.70	0.50
1:A:503:GLY:H	1:A:621:THR:HG23	1.76	0.49
1:B:554:GLU:HG3	1:B:558:LEU:CD2	2.42	0.49
1:B:557:ARG:NH2	2:B:2006:HOH:O	2.45	0.49
1:B:597:LEU:CD1	1:B:608:VAL:CG1	2.91	0.49
1:B:597:LEU:HD12	1:B:608:VAL:CG1	2.42	0.49
1:B:608:VAL:O	1:B:609:ASN:HB2	2.13	0.49
1:A:529:GLU:O	1:A:533:ARG:N	2.46	0.49
1:A:603:ASN:C	1:A:604:ASN:ND2	2.70	0.48
1:B:562:ALA:O	1:B:565:LEU:HB3	2.13	0.48
1:A:572:HIS:NE2	1:A:631:VAL:HG13	2.28	0.48
1:A:617:ASP:HB2	1:A:619:MET:HE1	1.96	0.48
1:A:561:ASP:O	1:A:564:GLU:HB2	2.14	0.48
1:A:543:THR:OG1	1:A:629:THR:HG22	2.13	0.48
1:B:557:ARG:HG3	1:B:557:ARG:NH1	2.24	0.48
1:A:512:ASP:HB3	1:A:515:PHE:CD2	2.49	0.47
1:B:561:ASP:OD2	1:B:564:GLU:HB2	2.15	0.47
1:A:586:LEU:C	1:A:586:LEU:HD22	2.40	0.47
1:B:503:GLY:O	1:B:504:THR:HB	2.14	0.47
1:B:582:GLY:O	1:B:583:ILE:HG23	2.14	0.47
1:A:538:THR:HG21	1:A:589:LEU:HD11	1.96	0.47
1:B:580:SER:HB2	1:B:617:ASP:OD2	2.14	0.47
1:B:504:THR:HG22	1:B:506:MET:N	2.28	0.47
1:B:615:GLU:HA	1:B:616:PRO:HD2	1.71	0.47
1:A:613:VAL:HG13	1:A:627:VAL:O	2.15	0.47
1:A:608:VAL:O	1:A:609:ASN:HB2	2.16	0.46
1:B:539:VAL:HG22	1:B:539:VAL:O	2.15	0.46
1:B:588:ARG:O	1:B:596:LYS:HE3	2.15	0.46
1:A:577:ILE:CG2	1:A:623:GLY:HA2	2.45	0.46
1:A:580:SER:O	1:A:581:GLY:C	2.56	0.46
1:B:577:ILE:HD13	1:B:624:VAL:CG1	2.46	0.46
1:A:538:THR:CG2	1:A:589:LEU:HD13	2.44	0.46
1:B:557:ARG:HH11	1:B:557:ARG:CG	2.26	0.46
1:B:577:ILE:HD13	1:B:624:VAL:HG11	1.97	0.46
1:B:582:GLY:O	1:B:583:ILE:CG1	2.61	0.46
1:B:602:LYS:O	1:B:603:ASN:C	2.59	0.45
1:B:515:PHE:HZ	1:B:627:VAL:HG22	1.82	0.45
1:B:542:PRO:HA	1:B:628:ILE:HG13	1.97	0.45
1:B:587:VAL:HG12	1:B:599:VAL:HB	1.97	0.45
1:A:502:MET:HB3	1:A:621:THR:HG23	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:584:GLY:HA2	1:B:601:LEU:HD23	1.98	0.45
1:B:554:GLU:OE2	1:B:554:GLU:HA	2.16	0.44
1:A:533:ARG:HB3	1:A:534:GLU:H	1.62	0.44
1:A:583:ILE:HG21	1:A:600:SER:HA	2.00	0.44
1:B:503:GLY:HA2	1:B:621:THR:OG1	2.18	0.44
1:A:565:LEU:O	1:A:566:ALA:C	2.61	0.44
1:B:604:ASN:HB3	2:B:2013:HOH:O	2.16	0.44
1:B:576:GLU:HG3	1:B:589:LEU:HD11	1.97	0.44
1:B:583:ILE:O	1:B:584:GLY:C	2.61	0.43
1:B:598:GLU:O	1:B:608:VAL:HA	2.17	0.43
1:B:614:ALA:O	1:B:627:VAL:HB	2.17	0.43
1:B:568:ILE:O	1:B:572:HIS:ND1	2.50	0.43
1:A:613:VAL:HG12	1:A:615:GLU:H	1.84	0.43
1:B:587:VAL:CG1	1:B:599:VAL:HB	2.49	0.43
1:A:564:GLU:O	1:A:567:ASN:HB3	2.18	0.43
1:A:536:VAL:CG1	1:A:577:ILE:HD12	2.49	0.42
1:B:515:PHE:CZ	1:B:627:VAL:HG22	2.54	0.42
1:B:578:LEU:HD23	1:B:579:VAL:H	1.84	0.42
1:B:589:LEU:O	1:B:596:LYS:HA	2.19	0.42
1:A:599:VAL:HA	1:A:607:SER:O	2.19	0.42
1:B:503:GLY:HA3	1:B:621:THR:HG23	2.01	0.42
1:B:540:PHE:HE1	1:B:597:LEU:HD21	1.82	0.42
1:B:602:LYS:HA	1:B:602:LYS:HD2	1.34	0.42
1:B:550:LEU:HD12	1:B:550:LEU:HA	1.89	0.42
1:B:609:ASN:N	1:B:609:ASN:ND2	2.65	0.42
1:A:563:LYS:HA	1:A:563:LYS:HD2	1.61	0.42
1:A:551:PRO:HG2	1:A:554:GLU:CB	2.48	0.42
1:A:564:GLU:OE2	1:A:564:GLU:HA	2.20	0.41
1:B:572:HIS:NE2	1:B:632:LEU:HB2	2.34	0.41
1:B:580:SER:C	1:B:581:GLY:O	2.58	0.41
1:A:503:GLY:HA2	2:A:2001:HOH:O	2.21	0.41
1:B:610:LYS:HD3	2:B:2015:HOH:O	2.21	0.41
1:A:611:GLU:OE2	1:A:612:PRO:HD2	2.21	0.41
1:B:561:ASP:O	1:B:562:ALA:C	2.64	0.41
1:B:589:LEU:O	1:B:596:LYS:HG3	2.22	0.40
1:B:583:ILE:HD12	1:B:587:VAL:HG12	2.03	0.40
1:A:608:VAL:O	1:A:608:VAL:CG1	2.66	0.40
1:B:588:ARG:HG2	1:B:598:GLU:OE1	2.22	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	129/132 (98%)	118 (92%)	8 (6%)	3 (2%)	5	8
1	B	129/132 (98%)	108 (84%)	16 (12%)	5 (4%)	2	3
All	All	258/264 (98%)	226 (88%)	24 (9%)	8 (3%)	3	5

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	617	ASP
1	B	583	ILE
1	B	585	ALA
1	B	584	GLY
1	B	615	GLU
1	A	533	ARG
1	A	564	GLU
1	B	580	SER

### 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	109/110 (99%)	96 (88%)	13 (12%)	5	11
1	B	109/110 (99%)	89 (82%)	20 (18%)	2	3
All	All	218/220 (99%)	185 (85%)	33 (15%)	3	5

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	508	VAL
1	A	524	SER
1	A	534	GLU
1	A	538	THR
1	A	558	LEU
1	A	559	LEU
1	A	564	GLU
1	A	575	ASP
1	A	576	GLU
1	A	577	ILE
1	A	586	LEU
1	A	624	VAL
1	A	625	VAL
1	B	513	ASN
1	B	514	ARG
1	B	524	SER
1	B	536	VAL
1	B	539	VAL
1	B	545	GLU
1	B	550	LEU
1	B	553	ARG
1	B	554	GLU
1	B	556	SER
1	B	557	ARG
1	B	558	LEU
1	B	564	GLU
1	B	577	ILE
1	B	586	LEU
1	B	597	LEU
1	B	602	LYS
1	B	618	ILE
1	B	619	MET
1	B	624	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	532	ASN
1	A	544	ASN
1	A	603	ASN
1	B	513	ASN
1	B	567	ASN
1	B	603	ASN

*Continued on next page...*

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Mol	Chain	Res	Type
1	B	609	ASN

### 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 [i](#)

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

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	131/132 (99%)	-1.27	0 <a href="#">100</a> <a href="#">100</a>	11, 31, 42, 46	0
1	B	131/132 (99%)	-1.22	0 <a href="#">100</a> <a href="#">100</a>	14, 38, 51, 59	0
All	All	262/264 (99%)	-1.24	0 <a href="#">100</a> <a href="#">100</a>	11, 33, 48, 59	0

There are no RSRZ outliers to report.

### 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.