



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 03:32 AM UTC

PDB ID : 1PDG / pdb_00001pdg
Title : CRYSTAL STRUCTURE OF HUMAN PLATELET-DERIVED GROWTH FACTOR BB
Authors : Oefner, C.; Darcy, A.D.; Winkler, F.K.; Eggimann, B.; Hosnag, M.
Deposited on : 1992-07-14
Resolution : 3.00 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	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

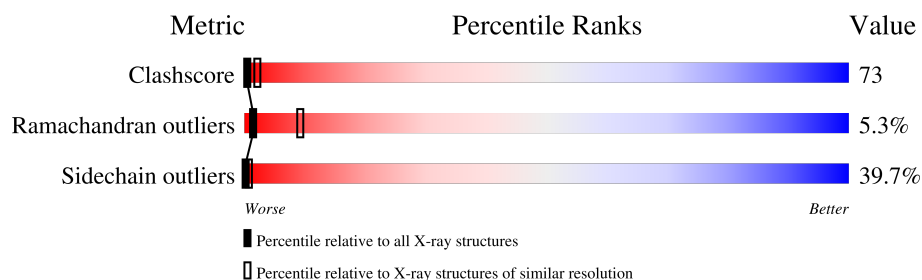
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	109	
1	B	109	
1	C	109	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2035 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 PLATELET-DERIVED GROWTH FACTOR BB.

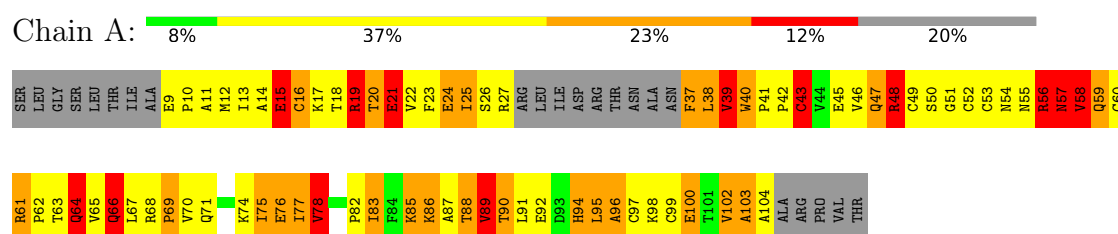
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	87	Total	C	N	O	S	0	0	0
			692	433	129	121	9			
1	B	83	Total	C	N	O	S	0	0	0
			659	412	122	116	9			
1	C	86	Total	C	N	O	S	0	0	0
			684	430	125	120	9			

3 Residue-property plots

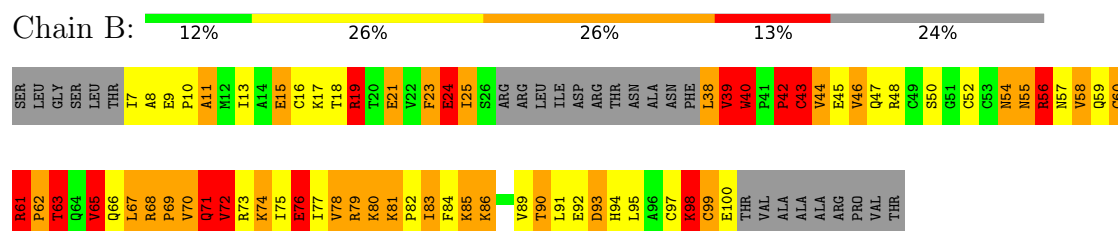
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

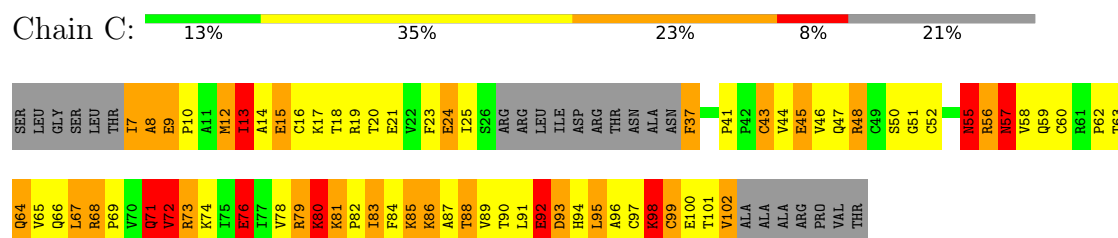
• Molecule 1: PLATELET-DERIVED GROWTH FACTOR BB



• Molecule 1: PLATELET-DERIVED GROWTH FACTOR BB



• Molecule 1: PLATELET-DERIVED GROWTH FACTOR BB



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	147.30Å 31.80Å 90.10Å 90.00° 98.70° 90.00°	Depositor
Resolution (Å)	6.00 – 3.00	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-3.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.209 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2035	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

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	1.54	4/702 (0.6%)	2.61	58/947 (6.1%)
1	B	1.40	1/668 (0.1%)	2.40	48/901 (5.3%)
1	C	1.38	3/694 (0.4%)	2.34	27/937 (2.9%)
All	All	1.44	8/2064 (0.4%)	2.45	133/2785 (4.8%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	2
All	All	0	3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	88	THR	CA-CB	7.43	1.62	1.52
1	A	45	GLU	CA-CB	-6.68	1.43	1.53
1	B	44	VAL	N-CA	6.23	1.53	1.46
1	A	88	THR	N-CA	-5.77	1.39	1.46
1	C	45	GLU	CA-CB	-5.71	1.44	1.53
1	A	59	GLN	N-CA	5.62	1.52	1.45
1	C	73	ARG	CA-CB	-5.20	1.45	1.53
1	C	72	VAL	N-CA	5.10	1.52	1.46

All (133) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	73	ARG	CA-CB-CG	17.08	148.26	114.10
1	B	44	VAL	CB-CA-C	14.36	125.23	111.74
1	C	93	ASP	CA-CB-CG	14.01	126.61	112.60

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	37	PHE	CA-CB-CG	13.17	126.97	113.80
1	C	45	GLU	CA-CB-CG	13.12	140.34	114.10
1	A	45	GLU	CA-CB-CG	11.82	137.75	114.10
1	A	87	ALA	CA-C-N	9.58	136.50	122.41
1	A	87	ALA	C-N-CA	9.58	136.50	122.41
1	B	57	ASN	CA-CB-CG	9.44	122.04	112.60
1	C	72	VAL	N-CA-CB	-8.60	99.58	111.19
1	A	89	VAL	CA-C-O	-8.47	111.38	121.80
1	A	39	VAL	N-CA-CB	-8.46	102.50	112.40
1	B	65	VAL	CB-CA-C	8.40	125.06	111.29
1	B	43	CYS	N-CA-CB	-8.16	99.11	110.38
1	A	19	ARG	NE-CZ-NH1	-8.04	113.46	121.50
1	B	57	ASN	CB-CA-C	7.91	124.29	110.85
1	C	98	LYS	N-CA-CB	7.83	123.73	110.49
1	A	16	CYS	CA-C-O	-7.83	112.50	121.66
1	B	19	ARG	CA-C-O	-7.82	112.30	121.44
1	B	44	VAL	CA-C-O	-7.79	115.44	120.66
1	A	89	VAL	N-CA-C	-7.37	99.21	109.45
1	C	57	ASN	N-CA-C	-7.32	103.94	114.12
1	A	97	CYS	CA-CB-SG	7.30	131.18	114.40
1	C	73	ARG	CB-CG-CD	7.25	127.97	111.30
1	A	21	GLU	CB-CG-CD	7.17	124.78	112.60
1	C	24	GLU	CA-C-O	-7.15	112.86	120.80
1	A	25	ILE	N-CA-CB	7.14	120.46	111.46
1	A	25	ILE	CA-C-O	7.11	128.14	120.53
1	A	47	GLN	CA-C-N	6.93	134.03	122.87
1	A	47	GLN	C-N-CA	6.93	134.03	122.87
1	B	72	VAL	N-CA-CB	-6.91	100.11	111.44
1	B	56	ARG	CD-NE-CZ	6.85	134.00	124.40
1	C	55	ASN	CA-C-N	6.77	134.47	121.54
1	C	55	ASN	C-N-CA	6.77	134.47	121.54
1	A	57	ASN	CA-CB-CG	6.74	119.34	112.60
1	B	19	ARG	CD-NE-CZ	-6.63	115.12	124.40
1	C	76	GLU	N-CA-CB	-6.62	100.55	111.74
1	B	23	PHE	CA-CB-CG	6.61	120.41	113.80
1	B	45	GLU	CB-CG-CD	6.60	123.83	112.60
1	B	43	CYS	CA-CB-SG	-6.54	99.36	114.40
1	A	100	GLU	CA-C-N	6.53	131.19	121.72
1	A	100	GLU	C-N-CA	6.53	131.19	121.72
1	A	78	VAL	N-CA-C	6.50	117.24	107.75
1	A	43	CYS	CA-C-N	6.42	131.42	121.85
1	A	43	CYS	C-N-CA	6.42	131.42	121.85

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	37	PHE	CA-CB-CG	-6.40	107.40	113.80
1	B	67	LEU	CB-CA-C	6.40	121.28	109.70
1	B	61	ARG	CA-C-N	6.39	126.15	119.76
1	B	61	ARG	C-N-CA	6.39	126.15	119.76
1	A	88	THR	CB-CA-C	-6.37	100.47	111.30
1	A	90	THR	CA-C-N	6.37	135.07	121.45
1	A	90	THR	C-N-CA	6.37	135.07	121.45
1	B	58	VAL	CA-C-O	-6.29	114.44	121.36
1	C	9	GLU	N-CA-CB	6.29	118.78	110.29
1	C	92	GLU	CA-CB-CG	6.27	126.64	114.10
1	B	76	GLU	CB-CG-CD	6.26	123.24	112.60
1	A	54	ASN	OD1-CG-ND2	6.25	128.85	122.60
1	C	71	GLN	N-CA-C	-6.24	100.13	110.17
1	C	21	GLU	CB-CG-CD	6.22	123.17	112.60
1	A	53	CYS	N-CA-C	6.21	116.47	108.34
1	B	42	PRO	CA-C-N	6.17	133.50	121.66
1	B	42	PRO	C-N-CA	6.17	133.50	121.66
1	A	64	GLN	OE1-CD-NE2	6.17	128.77	122.60
1	C	68	ARG	O-C-N	6.16	128.67	121.10
1	B	62	PRO	N-CA-C	6.07	120.72	111.19
1	B	21	GLU	CA-C-O	-6.05	114.03	121.11
1	B	95	LEU	N-CA-C	5.99	118.76	111.82
1	A	48	ARG	CD-NE-CZ	-5.97	116.05	124.40
1	C	80	LYS	CA-C-N	5.95	136.32	121.80
1	C	80	LYS	C-N-CA	5.95	136.32	121.80
1	B	43	CYS	CA-C-N	-5.92	116.05	121.97
1	B	43	CYS	C-N-CA	-5.92	116.05	121.97
1	B	72	VAL	CA-C-O	-5.88	114.08	120.67
1	B	40	TRP	N-CA-CB	5.85	120.78	110.37
1	A	96	ALA	N-CA-CB	-5.80	102.22	111.62
1	B	71	GLN	N-CA-CB	5.80	119.63	110.23
1	A	59	GLN	CB-CG-CD	-5.80	102.75	112.60
1	B	67	LEU	CA-C-O	5.79	126.75	120.32
1	B	15	GLU	CA-C-O	-5.79	114.82	121.07
1	C	13	ILE	CB-CA-C	5.78	119.47	111.38
1	A	58	VAL	O-C-N	5.73	129.14	123.18
1	A	24	GLU	CB-CG-CD	5.71	122.31	112.60
1	B	44	VAL	N-CA-CB	-5.66	102.37	109.84
1	A	57	ASN	CB-CA-C	5.66	121.67	110.42
1	A	58	VAL	N-CA-CB	5.62	119.46	111.82
1	A	90	THR	N-CA-CB	5.57	119.30	110.84
1	A	100	GLU	CA-C-O	5.54	127.42	121.16

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	50	SER	CA-C-O	-5.50	114.00	120.60
1	A	48	ARG	CA-C-N	5.50	130.19	121.99
1	A	48	ARG	C-N-CA	5.50	130.19	121.99
1	A	63	THR	CB-CA-C	5.50	118.29	109.27
1	B	16	CYS	N-CA-C	5.50	117.19	108.67
1	A	40	TRP	CA-C-O	-5.49	115.16	119.66
1	A	67	LEU	CA-CB-CG	5.49	135.50	116.30
1	A	59	GLN	CB-CA-C	5.41	119.99	109.33
1	B	45	GLU	CG-CD-OE1	5.41	130.85	118.40
1	B	58	VAL	O-C-N	5.40	128.35	122.63
1	A	103	ALA	N-CA-C	-5.39	102.22	110.30
1	A	39	VAL	CA-C-O	-5.36	112.94	121.36
1	B	11	ALA	O-C-N	5.33	128.99	122.96
1	A	60	CYS	CA-C-O	-5.33	115.38	120.92
1	C	73	ARG	N-CA-CB	5.31	117.96	110.36
1	A	66	GLN	OE1-CD-NE2	5.31	127.91	122.60
1	B	86	LYS	N-CA-CB	5.31	118.91	110.84
1	B	24	GLU	CB-CG-CD	5.30	121.62	112.60
1	C	64	GLN	CA-C-O	-5.30	115.17	121.16
1	A	25	ILE	CA-C-N	5.28	130.58	122.77
1	A	25	ILE	C-N-CA	5.28	130.58	122.77
1	A	56	ARG	CD-NE-CZ	5.28	131.78	124.40
1	A	99	CYS	O-C-N	5.26	129.42	122.89
1	B	63	THR	CA-CB-OG1	-5.26	101.72	109.60
1	C	13	ILE	CA-C-N	5.25	131.57	121.54
1	C	13	ILE	C-N-CA	5.25	131.57	121.54
1	B	93	ASP	CA-C-O	-5.25	115.47	121.82
1	B	69	PRO	CA-C-O	-5.23	115.20	121.48
1	B	23	PHE	N-CA-C	5.21	117.44	107.75
1	C	48	ARG	N-CA-C	5.20	115.22	107.88
1	A	48	ARG	N-CA-C	5.18	118.19	109.95
1	B	81	LYS	O-C-N	5.18	125.41	121.27
1	A	52	CYS	N-CA-C	5.17	116.53	109.18
1	A	69	PRO	CB-CA-C	-5.17	103.24	110.63
1	B	57	ASN	N-CA-C	-5.16	105.74	111.36
1	A	63	THR	CA-CB-CG2	5.15	119.25	110.50
1	B	98	LYS	CA-C-N	5.13	128.63	121.24
1	B	98	LYS	C-N-CA	5.13	128.63	121.24
1	C	43	CYS	N-CA-C	5.10	117.92	107.37
1	A	76	GLU	CA-C-O	-5.10	114.91	120.36
1	A	58	VAL	CA-C-N	-5.07	114.85	122.81
1	A	58	VAL	C-N-CA	-5.07	114.85	122.81

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	52	CYS	CB-CA-C	-5.05	100.50	111.11
1	C	9	GLU	CB-CG-CD	5.02	121.13	112.60
1	B	82	PRO	N-CA-C	-5.01	102.14	112.47
1	B	39	VAL	CB-CA-C	5.01	119.51	111.29

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	48	ARG	Sidechain
1	B	19	ARG	Sidechain
1	B	68	ARG	Sidechain

5.2 Too-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 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	692	0	713	108	0
1	B	659	0	681	104	0
1	C	684	0	708	104	2
All	All	2035	0	2102	303	2

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 73.

All (303) 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:78:VAL:HG12	1:C:81:LYS:NZ	1.52	1.21
1:C:68:ARG:NH1	1:C:91:LEU:HD12	1.62	1.15
1:A:19:ARG:HG2	1:A:19:ARG:HH11	1.09	1.13
1:C:78:VAL:HG12	1:C:81:LYS:HZ1	1.05	1.08
1:C:68:ARG:HH12	1:C:91:LEU:CD1	1.66	1.07
1:A:78:VAL:HB	1:A:83:ILE:HD11	1.33	1.07
1:A:38:LEU:HD13	1:A:38:LEU:H	0.94	1.06

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:PRO:HB2	1:B:43:CYS:SG	1.96	1.06
1:A:38:LEU:H	1:A:38:LEU:CD1	1.72	1.02
1:A:38:LEU:HD13	1:A:38:LEU:N	1.76	1.01
1:B:48:ARG:HD3	1:B:94:HIS:CD2	1.96	1.01
1:A:86:LYS:HZ3	1:A:86:LYS:HB2	1.24	1.00
1:A:38:LEU:HD23	1:A:75:ILE:HD12	1.46	0.98
1:C:59:GLN:HE21	1:C:102:VAL:HG11	1.25	0.97
1:B:59:GLN:HE22	1:B:61:ARG:NH1	1.62	0.96
1:B:76:GLU:HG2	1:B:85:LYS:HZ3	1.30	0.96
1:A:19:ARG:HH11	1:A:19:ARG:CG	1.76	0.95
1:C:68:ARG:HH12	1:C:91:LEU:HD12	0.79	0.95
1:A:86:LYS:HB2	1:A:86:LYS:NZ	1.81	0.94
1:B:55:ASN:H	1:B:55:ASN:ND2	1.60	0.94
1:B:55:ASN:HD22	1:B:55:ASN:N	1.63	0.94
1:B:23:PHE:HE2	1:B:91:LEU:CD2	1.81	0.93
1:C:98:LYS:HG3	1:C:99:CYS:H	1.30	0.93
1:C:66:GLN:HB2	1:C:95:LEU:CD1	1.99	0.93
1:A:38:LEU:CD2	1:A:75:ILE:HB	2.00	0.92
1:C:98:LYS:HG3	1:C:99:CYS:N	1.80	0.92
1:B:38:LEU:O	1:B:38:LEU:HD23	1.70	0.90
1:C:78:VAL:O	1:C:81:LYS:HD2	1.72	0.89
1:B:76:GLU:HG2	1:B:85:LYS:NZ	1.89	0.88
1:B:61:ARG:HB3	1:B:62:PRO:HD2	1.56	0.86
1:C:59:GLN:HB2	1:C:102:VAL:HB	1.54	0.86
1:C:7:ILE:CG1	1:C:8:ALA:H	1.87	0.86
1:B:23:PHE:CE2	1:B:91:LEU:CD2	2.59	0.86
1:B:25:ILE:HG13	1:B:42:PRO:O	1.75	0.86
1:C:9:GLU:HG3	1:C:10:PRO:HD2	1.59	0.85
1:C:79:ARG:O	1:C:81:LYS:HG3	1.76	0.85
1:A:55:ASN:HB2	1:A:58:VAL:HG22	1.59	0.85
1:C:7:ILE:HG13	1:C:8:ALA:H	1.41	0.84
1:C:9:GLU:HG3	1:C:10:PRO:CD	2.07	0.84
1:B:48:ARG:HA	1:B:94:HIS:CD2	2.13	0.83
1:A:78:VAL:HB	1:A:83:ILE:CD1	2.09	0.83
1:C:78:VAL:HG12	1:C:81:LYS:HZ2	1.44	0.80
1:B:48:ARG:HD3	1:B:94:HIS:NE2	1.97	0.79
1:A:14:ALA:HB1	1:A:19:ARG:HH21	1.49	0.78
1:A:21:GLU:HG3	1:A:48:ARG:HB2	1.65	0.77
1:A:77:ILE:HG12	1:A:82:PRO:HB3	1.66	0.77
1:A:56:ARG:CB	1:A:56:ARG:HH11	1.98	0.77
1:B:76:GLU:CG	1:B:85:LYS:NZ	2.49	0.76

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:46:VAL:HG21	1:C:92:GLU:O	1.86	0.75
1:C:66:GLN:NE2	1:C:95:LEU:HD11	2.02	0.75
1:A:19:ARG:HG2	1:A:19:ARG:NH1	1.95	0.75
1:B:24:GLU:HA	1:B:43:CYS:HA	1.68	0.74
1:A:38:LEU:HD23	1:A:75:ILE:CD1	2.15	0.74
1:B:38:LEU:HD23	1:B:38:LEU:C	2.13	0.73
1:C:98:LYS:CG	1:C:99:CYS:H	2.02	0.72
1:B:23:PHE:HE2	1:B:91:LEU:HD21	1.52	0.72
1:B:69:PRO:HA	1:B:90:THR:HA	1.71	0.72
1:B:23:PHE:HE2	1:B:91:LEU:HD23	1.52	0.71
1:A:38:LEU:HD21	1:A:75:ILE:HB	1.72	0.71
1:B:97:CYS:O	1:B:98:LYS:HD2	1.91	0.71
1:A:11:ALA:HB2	1:B:70:VAL:CG2	2.21	0.70
1:C:78:VAL:CG1	1:C:81:LYS:HZ1	1.95	0.70
1:B:59:GLN:NE2	1:B:61:ARG:NH1	2.38	0.70
1:C:78:VAL:O	1:C:81:LYS:CD	2.39	0.70
1:C:66:GLN:HB2	1:C:95:LEU:HD11	1.72	0.69
1:C:18:THR:HA	1:C:48:ARG:O	1.93	0.69
1:A:46:VAL:HG11	1:A:91:LEU:HD22	1.76	0.68
1:C:55:ASN:O	1:C:57:ASN:N	2.25	0.68
1:C:78:VAL:HB	1:C:83:ILE:HG13	1.76	0.67
1:B:23:PHE:CE2	1:B:91:LEU:HD23	2.28	0.67
1:A:42:PRO:HD2	1:B:13:ILE:HD12	1.77	0.67
1:A:55:ASN:HB2	1:A:58:VAL:CG2	2.25	0.67
1:B:55:ASN:H	1:B:55:ASN:HD22	0.80	0.66
1:B:79:ARG:O	1:B:81:LYS:HG2	1.94	0.66
1:A:49:CYS:HB2	1:A:94:HIS:CE1	2.31	0.66
1:A:49:CYS:HB2	1:A:94:HIS:ND1	2.10	0.66
1:A:62:PRO:HG3	1:A:94:HIS:CE1	2.31	0.66
1:C:66:GLN:HB3	1:C:93:ASP:HB2	1.76	0.66
1:A:76:GLU:O	1:A:83:ILE:HD12	1.96	0.66
1:B:48:ARG:HG3	1:B:48:ARG:HH11	1.60	0.65
1:C:44:VAL:HG21	1:C:91:LEU:HD11	1.77	0.65
1:A:16:CYS:O	1:A:16:CYS:SG	2.53	0.65
1:B:55:ASN:HB2	1:B:58:VAL:HG23	1.79	0.65
1:C:74:LYS:O	1:C:74:LYS:HG3	1.96	0.65
1:C:59:GLN:NE2	1:C:102:VAL:HG11	2.07	0.65
1:B:47:GLN:O	1:B:48:ARG:HG2	1.96	0.65
1:B:76:GLU:HG3	1:B:85:LYS:HZ2	1.62	0.64
1:C:44:VAL:CG2	1:C:91:LEU:HD11	2.27	0.64
1:A:55:ASN:H	1:A:58:VAL:CG2	2.11	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:79:ARG:HB3	1:C:81:LYS:HE3	1.80	0.64
1:A:24:GLU:HA	1:A:43:CYS:HB3	1.79	0.63
1:C:66:GLN:HB2	1:C:95:LEU:HD12	1.77	0.63
1:C:95:LEU:O	1:C:96:ALA:HB2	1.97	0.63
1:A:56:ARG:HH11	1:A:56:ARG:HB3	1.62	0.63
1:C:71:GLN:NE2	1:C:86:LYS:HD3	2.13	0.63
1:A:46:VAL:CG1	1:A:91:LEU:HD22	2.28	0.63
1:B:76:GLU:OE1	1:B:78:VAL:HG13	1.97	0.62
1:C:58:VAL:HG13	1:C:100:GLU:O	1.99	0.62
1:B:76:GLU:CG	1:B:85:LYS:HZ3	2.07	0.62
1:B:23:PHE:CE2	1:B:91:LEU:HD21	2.30	0.62
1:B:24:GLU:HA	1:B:43:CYS:CA	2.29	0.62
1:C:68:ARG:HG3	1:C:68:ARG:HH11	1.65	0.61
1:A:14:ALA:HB1	1:A:19:ARG:NH2	2.14	0.61
1:A:47:GLN:O	1:A:94:HIS:HB2	1.99	0.61
1:B:39:VAL:HG12	1:B:72:VAL:HG13	1.81	0.61
1:C:55:ASN:N	1:C:55:ASN:ND2	2.48	0.61
1:C:55:ASN:N	1:C:55:ASN:HD22	1.98	0.61
1:C:76:GLU:HB3	1:C:83:ILE:O	2.00	0.61
1:C:46:VAL:CG2	1:C:92:GLU:O	2.48	0.61
1:A:83:ILE:HD12	1:A:83:ILE:H	1.64	0.61
1:B:38:LEU:HD21	1:B:75:ILE:HD12	1.81	0.61
1:A:48:ARG:NH2	1:A:92:GLU:O	2.34	0.60
1:A:102:VAL:HG22	1:A:104:ALA:HB2	1.81	0.60
1:C:69:PRO:HA	1:C:89:VAL:O	2.01	0.60
1:C:66:GLN:HG3	1:C:67:LEU:H	1.65	0.60
1:B:70:VAL:HG11	1:B:91:LEU:HD12	1.82	0.60
1:A:42:PRO:HD2	1:B:13:ILE:CD1	2.31	0.60
1:A:74:LYS:HE3	1:A:76:GLU:HG2	1.84	0.60
1:C:85:LYS:HG2	1:C:86:LYS:H	1.67	0.60
1:B:59:GLN:NE2	1:B:61:ARG:HH11	2.01	0.59
1:B:56:ARG:HH11	1:B:56:ARG:HG3	1.67	0.59
1:B:48:ARG:HD3	1:B:94:HIS:HE2	1.67	0.59
1:B:72:VAL:HG21	1:B:89:VAL:CG1	2.33	0.59
1:C:71:GLN:CB	1:C:88:THR:HG23	2.32	0.58
1:A:19:ARG:CG	1:A:19:ARG:NH1	2.48	0.58
1:C:17:LYS:HG2	1:C:18:THR:H	1.68	0.58
1:A:77:ILE:CG1	1:A:82:PRO:HB3	2.33	0.58
1:A:62:PRO:HG3	1:A:94:HIS:NE2	2.19	0.58
1:A:77:ILE:HG12	1:A:82:PRO:CB	2.33	0.58
1:C:18:THR:HG23	1:C:18:THR:O	2.03	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:GLU:CG	1:A:48:ARG:HB2	2.34	0.57
1:B:61:ARG:CB	1:B:62:PRO:HD2	2.24	0.57
1:B:21:GLU:HG3	1:B:48:ARG:HB2	1.86	0.57
1:A:83:ILE:HD12	1:A:83:ILE:N	2.19	0.57
1:A:10:PRO:HB3	1:B:71:GLN:HG3	1.86	0.57
1:A:19:ARG:HG3	1:A:50:SER:OG	2.05	0.57
1:C:63:THR:HG22	1:C:64:GLN:CG	2.35	0.57
1:C:15:GLU:HG2	1:C:16:CYS:N	2.19	0.56
1:A:13:ILE:HD12	1:B:40:TRP:CZ2	2.40	0.56
1:A:20:THR:HG22	1:A:47:GLN:CG	2.34	0.56
1:A:37:PHE:O	1:A:39:VAL:HG22	2.06	0.56
1:B:76:GLU:CG	1:B:85:LYS:HZ2	2.17	0.56
1:C:79:ARG:CB	1:C:81:LYS:HE3	2.36	0.55
1:C:14:ALA:O	1:C:15:GLU:O	2.24	0.55
1:A:20:THR:HG22	1:A:47:GLN:HG2	1.88	0.55
1:A:38:LEU:HD23	1:A:75:ILE:CG1	2.35	0.55
1:B:43:CYS:O	1:B:44:VAL:HB	2.05	0.55
1:C:72:VAL:N	1:C:87:ALA:O	2.32	0.55
1:A:41:PRO:HG2	1:B:11:ALA:HB1	1.88	0.55
1:A:71:GLN:HB2	1:B:10:PRO:HA	1.87	0.55
1:A:20:THR:HG22	1:A:47:GLN:HE21	1.72	0.55
1:B:76:GLU:HB3	1:B:83:ILE:O	2.07	0.55
1:C:82:PRO:HB2	1:C:84:PHE:CE1	2.42	0.55
1:C:7:ILE:HG13	1:C:8:ALA:N	2.14	0.54
1:B:40:TRP:HA	1:B:42:PRO:HD3	1.89	0.54
1:B:70:VAL:CG1	1:B:91:LEU:HD12	2.37	0.54
1:A:55:ASN:H	1:A:58:VAL:HG21	1.73	0.54
1:B:38:LEU:HG	1:B:39:VAL:O	2.08	0.54
1:C:82:PRO:HB2	1:C:84:PHE:HE1	1.73	0.53
1:A:74:LYS:HE3	1:A:76:GLU:OE2	2.08	0.53
1:C:57:ASN:N	1:C:57:ASN:HD22	2.06	0.53
1:B:72:VAL:CG2	1:B:89:VAL:HG12	2.39	0.53
1:C:41:PRO:HD2	1:C:72:VAL:CG2	2.38	0.53
1:A:78:VAL:CB	1:A:83:ILE:HD11	2.23	0.53
1:A:76:GLU:O	1:A:83:ILE:CD1	2.57	0.53
1:C:71:GLN:HE22	1:C:86:LYS:HG2	1.74	0.52
1:A:11:ALA:HB2	1:B:70:VAL:HG21	1.90	0.52
1:B:70:VAL:HG11	1:B:91:LEU:CD1	2.38	0.52
1:A:15:GLU:CG	1:A:17:LYS:HD3	2.39	0.52
1:B:38:LEU:HD21	1:B:75:ILE:CD1	2.39	0.52
1:C:68:ARG:NH1	1:C:68:ARG:HG3	2.24	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:ALA:O	1:A:15:GLU:C	2.53	0.51
1:B:70:VAL:HG22	1:B:89:VAL:HG13	1.92	0.51
1:C:63:THR:HG22	1:C:64:GLN:HG2	1.92	0.51
1:A:74:LYS:HE3	1:A:76:GLU:CG	2.40	0.51
1:C:66:GLN:CB	1:C:93:ASP:HB2	2.39	0.51
1:C:62:PRO:HG3	1:C:94:HIS:CD2	2.46	0.51
1:B:68:ARG:NE	1:B:93:ASP:OD2	2.34	0.51
1:C:41:PRO:HD2	1:C:72:VAL:HG22	1.93	0.51
1:A:14:ALA:C	1:A:15:GLU:O	2.52	0.51
1:A:37:PHE:CD1	1:A:37:PHE:N	2.79	0.51
1:A:14:ALA:O	1:A:15:GLU:O	2.29	0.50
1:A:56:ARG:HH11	1:A:56:ARG:HB2	1.75	0.50
1:B:72:VAL:HG21	1:B:89:VAL:HG11	1.93	0.50
1:B:68:ARG:NH1	1:B:91:LEU:CD1	2.74	0.50
1:C:47:GLN:O	1:C:48:ARG:HG2	2.11	0.50
1:A:66:GLN:HB2	1:A:95:LEU:HD22	1.92	0.50
1:B:56:ARG:HH11	1:B:56:ARG:CG	2.23	0.50
1:C:58:VAL:HG12	1:C:59:GLN:N	2.26	0.50
1:B:68:ARG:NH1	1:B:91:LEU:HD13	2.27	0.50
1:C:93:ASP:O	1:C:95:LEU:HD12	2.11	0.50
1:B:66:GLN:HG2	1:B:67:LEU:N	2.26	0.50
1:B:23:PHE:O	1:B:43:CYS:HA	2.12	0.49
1:C:79:ARG:CA	1:C:81:LYS:HE3	2.41	0.49
1:A:38:LEU:CD2	1:A:75:ILE:CB	2.84	0.49
1:C:59:GLN:HB2	1:C:102:VAL:CB	2.36	0.49
1:C:18:THR:O	1:C:18:THR:CG2	2.60	0.49
1:B:39:VAL:CG1	1:B:72:VAL:HG13	2.42	0.49
1:B:74:LYS:O	1:B:85:LYS:N	2.38	0.49
1:C:9:GLU:HG3	1:C:10:PRO:CG	2.42	0.49
1:C:79:ARG:O	1:C:80:LYS:C	2.55	0.49
1:A:41:PRO:HG2	1:B:11:ALA:CB	2.43	0.49
1:A:57:ASN:HB3	1:A:104:ALA:O	2.13	0.49
1:A:58:VAL:HA	1:A:100:GLU:O	2.12	0.49
1:C:85:LYS:HD3	1:C:86:LYS:O	2.12	0.49
1:C:14:ALA:O	1:C:15:GLU:C	2.55	0.49
1:C:59:GLN:O	1:C:99:CYS:HA	2.13	0.49
1:C:78:VAL:CG1	1:C:81:LYS:NZ	2.47	0.49
1:C:14:ALA:HB3	1:C:52:CYS:SG	2.52	0.48
1:A:25:ILE:HD13	1:A:89:VAL:HG11	1.93	0.48
1:A:15:GLU:HG2	1:A:17:LYS:HD3	1.95	0.48
1:A:76:GLU:OE1	1:A:85:LYS:HD3	2.14	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:80:LYS:H	1:C:80:LYS:HD3	1.78	0.48
1:C:63:THR:HG22	1:C:64:GLN:HG3	1.95	0.48
1:A:27:ARG:NH1	1:A:37:PHE:CD1	2.81	0.48
1:A:74:LYS:HE3	1:A:76:GLU:CD	2.38	0.48
1:C:23:PHE:HD2	1:C:25:ILE:HG23	1.79	0.48
1:B:25:ILE:HG21	1:B:89:VAL:HG21	1.94	0.47
1:B:63:THR:O	1:B:63:THR:HG23	2.14	0.47
1:C:68:ARG:O	1:C:91:LEU:HB2	2.13	0.47
1:A:48:ARG:HA	1:A:48:ARG:HD3	1.70	0.47
1:A:102:VAL:HG22	1:A:102:VAL:O	2.13	0.47
1:A:48:ARG:HH11	1:A:48:ARG:HD2	1.57	0.47
1:B:65:VAL:HG23	1:B:92:GLU:HG2	1.96	0.47
1:C:71:GLN:HB3	1:C:88:THR:HG23	1.97	0.47
1:A:11:ALA:HB2	1:B:70:VAL:HB	1.97	0.47
1:B:39:VAL:CG1	1:B:72:VAL:CG1	2.93	0.46
1:A:38:LEU:HD22	1:A:75:ILE:HB	1.95	0.46
1:B:48:ARG:HD3	1:B:94:HIS:HD2	1.65	0.46
1:C:7:ILE:CG1	1:C:8:ALA:N	2.61	0.46
1:A:15:GLU:OE1	1:A:15:GLU:N	2.44	0.46
1:C:59:GLN:HG3	1:C:102:VAL:HG11	1.98	0.46
1:C:71:GLN:HA	1:C:88:THR:HA	1.98	0.46
1:A:9:GLU:HA	1:A:10:PRO:HD3	1.67	0.45
1:A:69:PRO:HB2	1:B:8:ALA:HB2	1.98	0.45
1:B:55:ASN:O	1:B:58:VAL:N	2.49	0.45
1:C:13:ILE:HG13	1:C:14:ALA:N	2.31	0.45
1:B:80:LYS:HD3	1:B:80:LYS:HA	1.80	0.45
1:C:37:PHE:CE1	1:C:74:LYS:NZ	2.79	0.45
1:B:72:VAL:HG21	1:B:89:VAL:HG12	1.99	0.45
1:C:66:GLN:HE21	1:C:95:LEU:HD11	1.79	0.45
1:A:16:CYS:HA	1:A:51:GLY:HA3	1.99	0.45
1:A:69:PRO:HB2	1:B:8:ALA:CB	2.47	0.45
1:B:38:LEU:C	1:B:38:LEU:CD2	2.82	0.45
1:B:77:ILE:HG22	1:B:77:ILE:O	2.16	0.45
1:A:49:CYS:CB	1:A:94:HIS:ND1	2.78	0.44
1:B:54:ASN:O	1:B:55:ASN:C	2.60	0.44
1:B:60:CYS:HA	1:B:99:CYS:HA	1.99	0.44
1:B:81:LYS:HB3	1:B:81:LYS:HE3	1.68	0.44
1:A:46:VAL:HG12	1:A:91:LEU:HD13	1.98	0.44
1:C:57:ASN:N	1:C:57:ASN:ND2	2.65	0.44
1:A:55:ASN:N	1:A:58:VAL:HG23	2.32	0.44
1:C:69:PRO:HA	1:C:90:THR:HA	2.00	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:GLU:H	1:B:84:PHE:HA	1.82	0.44
1:A:70:VAL:O	1:A:88:THR:HA	2.18	0.44
1:B:72:VAL:CG2	1:B:89:VAL:CG1	2.96	0.43
1:B:72:VAL:HG23	1:B:89:VAL:HG12	2.00	0.43
1:A:18:THR:O	1:A:18:THR:OG1	2.33	0.43
1:A:11:ALA:HB2	1:B:70:VAL:CB	2.48	0.43
1:A:38:LEU:CD2	1:A:75:ILE:HD12	2.32	0.43
1:B:65:VAL:CG2	1:B:92:GLU:CG	2.96	0.43
1:C:14:ALA:O	1:C:51:GLY:CA	2.67	0.43
1:B:76:GLU:CB	1:B:83:ILE:O	2.65	0.43
1:A:64:GLN:HG2	1:A:96:ALA:HB3	2.00	0.43
1:C:14:ALA:O	1:C:51:GLY:HA2	2.18	0.43
1:C:84:PHE:N	1:C:84:PHE:CD1	2.85	0.43
1:B:46:VAL:HG23	1:B:93:ASP:OD1	2.17	0.43
1:C:66:GLN:HG3	1:C:67:LEU:N	2.33	0.43
1:A:41:PRO:HA	1:A:42:PRO:HD3	1.89	0.42
1:A:92:GLU:O	1:A:92:GLU:HG2	2.19	0.42
1:B:65:VAL:CG2	1:B:92:GLU:HG2	2.49	0.42
1:C:18:THR:OG1	1:C:47:GLN:HG2	2.18	0.42
1:B:25:ILE:CG1	1:B:42:PRO:O	2.55	0.42
1:A:61:ARG:HB2	1:A:100:GLU:OE1	2.19	0.42
1:B:76:GLU:OE1	1:B:78:VAL:HG22	2.19	0.42
1:C:74:LYS:HG2	1:C:85:LYS:HB3	2.01	0.42
1:B:65:VAL:CG2	1:B:66:GLN:H	2.32	0.42
1:C:23:PHE:CD2	1:C:25:ILE:CG2	3.03	0.42
1:A:55:ASN:N	1:A:58:VAL:CG2	2.80	0.42
1:C:23:PHE:CD2	1:C:25:ILE:HG23	2.55	0.42
1:B:66:GLN:O	1:B:67:LEU:HD23	2.20	0.42
1:A:77:ILE:HG12	1:A:82:PRO:CA	2.50	0.41
1:B:83:ILE:O	1:B:83:ILE:HG13	2.20	0.41
1:A:64:GLN:CG	1:A:96:ALA:HB3	2.49	0.41
1:A:70:VAL:HG23	1:A:89:VAL:HG13	2.01	0.41
1:A:102:VAL:O	1:A:103:ALA:C	2.64	0.41
1:A:68:ARG:HA	1:A:69:PRO:HD3	1.82	0.41
1:C:67:LEU:HD12	1:C:67:LEU:HA	1.72	0.41
1:C:74:LYS:CG	1:C:85:LYS:HB3	2.50	0.41
1:A:74:LYS:CE	1:A:76:GLU:OE2	2.68	0.41
1:B:23:PHE:CD2	1:B:46:VAL:HG11	2.55	0.41
1:B:46:VAL:HB	1:B:68:ARG:HH22	1.86	0.41
1:B:68:ARG:HH21	1:B:93:ASP:CG	2.27	0.41
1:C:23:PHE:HD2	1:C:25:ILE:CG2	2.33	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:GLN:HB2	1:A:95:LEU:CD2	2.49	0.41
1:B:48:ARG:HA	1:B:94:HIS:CG	2.53	0.41
1:A:15:GLU:HG3	1:A:17:LYS:HD3	2.02	0.41
1:B:48:ARG:CD	1:B:94:HIS:HE2	2.31	0.41
1:C:58:VAL:CG1	1:C:100:GLU:O	2.68	0.41
1:C:12:MET:HE2	1:C:13:ILE:H	1.86	0.40
1:C:85:LYS:HE3	1:C:86:LYS:O	2.22	0.40
1:C:25:ILE:HD13	1:C:25:ILE:HG21	1.87	0.40
1:A:37:PHE:O	1:A:38:LEU:C	2.63	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:43:CYS:SG	1:C:52:CYS:SG[2_555]	1.72	0.48
1:C:19:ARG:NH2	1:C:45:GLU:OE1[2_555]	2.06	0.14

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	83/109 (76%)	68 (82%)	14 (17%)	1 (1%)	10	40
1	B	79/109 (72%)	61 (77%)	13 (16%)	5 (6%)	1	6
1	C	82/109 (75%)	59 (72%)	16 (20%)	7 (8%)	0	3
All	All	244/327 (75%)	188 (77%)	43 (18%)	13 (5%)	1	9

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	42	PRO
1	B	54	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	79	ARG
1	C	15	GLU
1	C	56	ARG
1	A	15	GLU
1	C	8	ALA
1	C	99	CYS
1	C	79	ARG
1	C	80	LYS
1	C	98	LYS
1	B	40	TRP
1	B	83	ILE

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	80/98 (82%)	47 (59%)	33 (41%)	0	0
1	B	77/98 (79%)	42 (54%)	35 (46%)	0	0
1	C	80/98 (82%)	54 (68%)	26 (32%)	0	1
All	All	237/294 (81%)	143 (60%)	94 (40%)	0	1

All (94) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	MET
1	A	15	GLU
1	A	19	ARG
1	A	20	THR
1	A	21	GLU
1	A	22	VAL
1	A	23	PHE
1	A	26	SER
1	A	38	LEU
1	A	39	VAL
1	A	40	TRP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	43	CYS
1	A	48	ARG
1	A	56	ARG
1	A	57	ASN
1	A	58	VAL
1	A	59	GLN
1	A	61	ARG
1	A	64	GLN
1	A	65	VAL
1	A	66	GLN
1	A	75	ILE
1	A	77	ILE
1	A	78	VAL
1	A	83	ILE
1	A	85	LYS
1	A	86	LYS
1	A	89	VAL
1	A	90	THR
1	A	94	HIS
1	A	95	LEU
1	A	98	LYS
1	A	102	VAL
1	B	7	ILE
1	B	9	GLU
1	B	15	GLU
1	B	17	LYS
1	B	18	THR
1	B	19	ARG
1	B	24	GLU
1	B	25	ILE
1	B	38	LEU
1	B	39	VAL
1	B	40	TRP
1	B	42	PRO
1	B	43	CYS
1	B	46	VAL
1	B	52	CYS
1	B	55	ASN
1	B	56	ARG
1	B	60	CYS
1	B	61	ARG
1	B	63	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	65	VAL
1	B	70	VAL
1	B	71	GLN
1	B	72	VAL
1	B	73	ARG
1	B	74	LYS
1	B	76	GLU
1	B	78	VAL
1	B	80	LYS
1	B	85	LYS
1	B	86	LYS
1	B	90	THR
1	B	98	LYS
1	B	99	CYS
1	B	100	GLU
1	C	7	ILE
1	C	12	MET
1	C	13	ILE
1	C	20	THR
1	C	24	GLU
1	C	50	SER
1	C	55	ASN
1	C	56	ARG
1	C	57	ASN
1	C	60	CYS
1	C	65	VAL
1	C	67	LEU
1	C	71	GLN
1	C	72	VAL
1	C	73	ARG
1	C	76	GLU
1	C	81	LYS
1	C	83	ILE
1	C	85	LYS
1	C	86	LYS
1	C	88	THR
1	C	92	GLU
1	C	95	LEU
1	C	97	CYS
1	C	101	THR
1	C	102	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	57	ASN
1	B	55	ASN
1	B	59	GLN
1	B	66	GLN
1	C	55	ASN
1	C	59	GLN
1	C	64	GLN
1	C	66	GLN
1	C	71	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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.