



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:23 PM UTC

PDB ID : 3BEZ / pdb_00003bez
Title : Crystal structure of Escherichia coli Signal peptide peptidase (SppA), SeMet crystals
Authors : Paetzel, M.
Deposited on : 2007-11-20
Resolution : 2.76 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
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (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

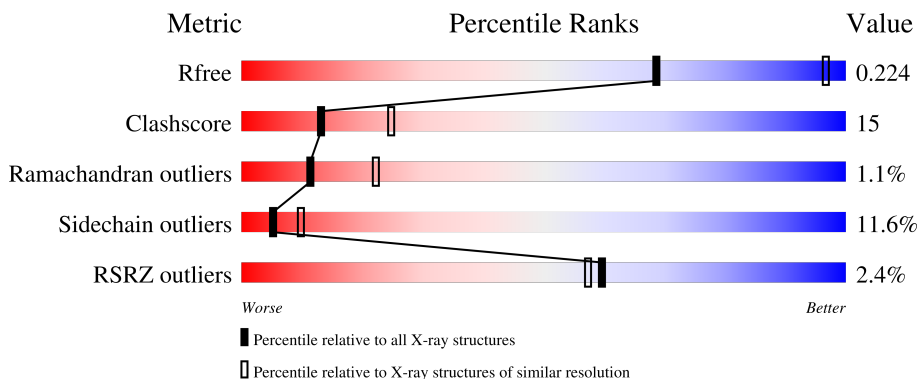
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.76 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	593	2% 54% 21% • • 20%
1	B	593	3% 53% 21% 5% • 20%
1	C	593	2% 53% 23% • • 20%
1	D	593	2% 57% 20% • • 19%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14989 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 Protease 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	473	Total	C	N	O	Se	0	0	0
			3576	2257	610	701	8			
1	B	473	Total	C	N	O	Se	0	0	0
			3572	2254	609	701	8			
1	C	476	Total	C	N	O	Se	0	0	0
			3597	2269	616	704	8			
1	D	479	Total	C	N	O	Se	0	0	0
			3617	2279	619	711	8			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	MSE	-	expression tag	UNP P08395
A	27	GLY	-	expression tag	UNP P08395
A	28	SER	-	expression tag	UNP P08395
A	29	SER	-	expression tag	UNP P08395
A	30	HIS	-	expression tag	UNP P08395
A	31	HIS	-	expression tag	UNP P08395
A	32	HIS	-	expression tag	UNP P08395
A	33	HIS	-	expression tag	UNP P08395
A	34	HIS	-	expression tag	UNP P08395
A	35	HIS	-	expression tag	UNP P08395
A	36	SER	-	expression tag	UNP P08395
A	37	SER	-	expression tag	UNP P08395
A	38	GLY	-	expression tag	UNP P08395
A	39	LEU	-	expression tag	UNP P08395
A	40	VAL	-	expression tag	UNP P08395
A	41	PRO	-	expression tag	UNP P08395
A	42	ARG	-	expression tag	UNP P08395
A	43	GLY	-	expression tag	UNP P08395
A	44	SER	-	expression tag	UNP P08395
A	45	HIS	-	expression tag	UNP P08395
A	46	MSE	-	expression tag	UNP P08395

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Chain	Residue	Modelled	Actual	Comment	Reference
B	26	MSE	-	expression tag	UNP P08395
B	27	GLY	-	expression tag	UNP P08395
B	28	SER	-	expression tag	UNP P08395
B	29	SER	-	expression tag	UNP P08395
B	30	HIS	-	expression tag	UNP P08395
B	31	HIS	-	expression tag	UNP P08395
B	32	HIS	-	expression tag	UNP P08395
B	33	HIS	-	expression tag	UNP P08395
B	34	HIS	-	expression tag	UNP P08395
B	35	HIS	-	expression tag	UNP P08395
B	36	SER	-	expression tag	UNP P08395
B	37	SER	-	expression tag	UNP P08395
B	38	GLY	-	expression tag	UNP P08395
B	39	LEU	-	expression tag	UNP P08395
B	40	VAL	-	expression tag	UNP P08395
B	41	PRO	-	expression tag	UNP P08395
B	42	ARG	-	expression tag	UNP P08395
B	43	GLY	-	expression tag	UNP P08395
B	44	SER	-	expression tag	UNP P08395
B	45	HIS	-	expression tag	UNP P08395
B	46	MSE	-	expression tag	UNP P08395
C	26	MSE	-	expression tag	UNP P08395
C	27	GLY	-	expression tag	UNP P08395
C	28	SER	-	expression tag	UNP P08395
C	29	SER	-	expression tag	UNP P08395
C	30	HIS	-	expression tag	UNP P08395
C	31	HIS	-	expression tag	UNP P08395
C	32	HIS	-	expression tag	UNP P08395
C	33	HIS	-	expression tag	UNP P08395
C	34	HIS	-	expression tag	UNP P08395
C	35	HIS	-	expression tag	UNP P08395
C	36	SER	-	expression tag	UNP P08395
C	37	SER	-	expression tag	UNP P08395
C	38	GLY	-	expression tag	UNP P08395
C	39	LEU	-	expression tag	UNP P08395
C	40	VAL	-	expression tag	UNP P08395
C	41	PRO	-	expression tag	UNP P08395
C	42	ARG	-	expression tag	UNP P08395
C	43	GLY	-	expression tag	UNP P08395
C	44	SER	-	expression tag	UNP P08395
C	45	HIS	-	expression tag	UNP P08395
C	46	MSE	-	expression tag	UNP P08395

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Chain	Residue	Modelled	Actual	Comment	Reference
D	26	MSE	-	expression tag	UNP P08395
D	27	GLY	-	expression tag	UNP P08395
D	28	SER	-	expression tag	UNP P08395
D	29	SER	-	expression tag	UNP P08395
D	30	HIS	-	expression tag	UNP P08395
D	31	HIS	-	expression tag	UNP P08395
D	32	HIS	-	expression tag	UNP P08395
D	33	HIS	-	expression tag	UNP P08395
D	34	HIS	-	expression tag	UNP P08395
D	35	HIS	-	expression tag	UNP P08395
D	36	SER	-	expression tag	UNP P08395
D	37	SER	-	expression tag	UNP P08395
D	38	GLY	-	expression tag	UNP P08395
D	39	LEU	-	expression tag	UNP P08395
D	40	VAL	-	expression tag	UNP P08395
D	41	PRO	-	expression tag	UNP P08395
D	42	ARG	-	expression tag	UNP P08395
D	43	GLY	-	expression tag	UNP P08395
D	44	SER	-	expression tag	UNP P08395
D	45	HIS	-	expression tag	UNP P08395
D	46	MSE	-	expression tag	UNP P08395

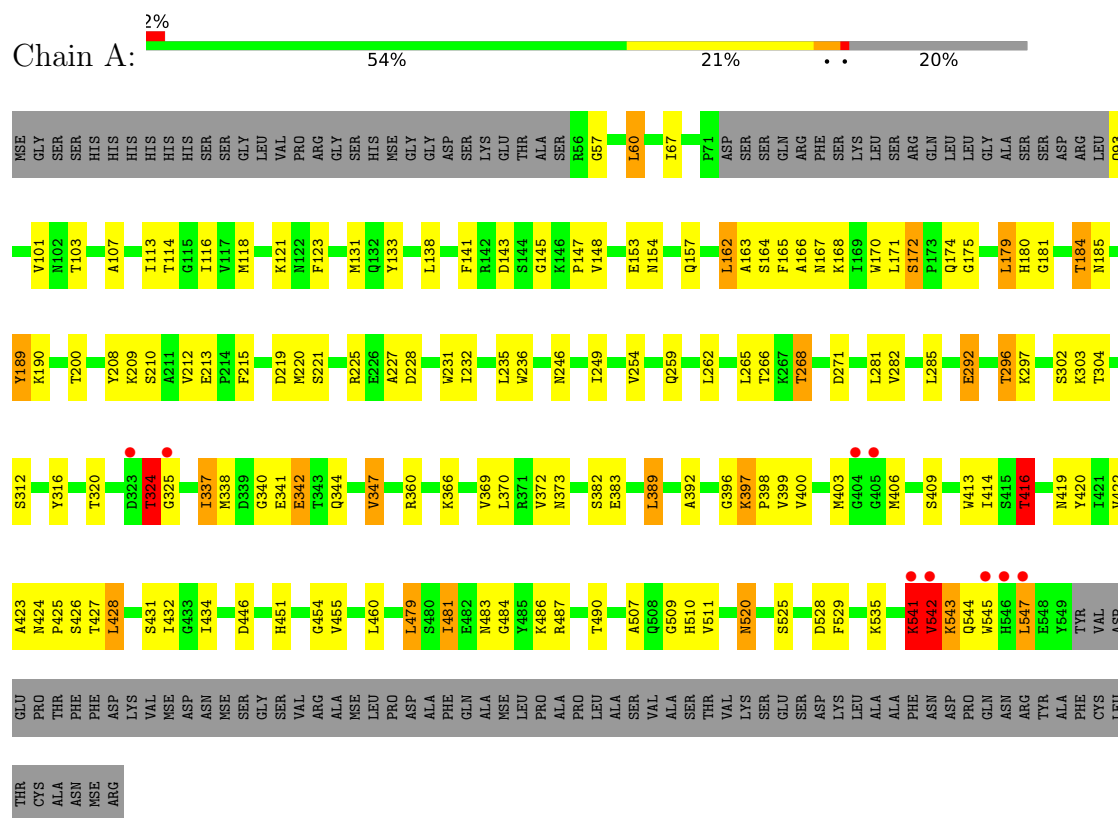
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	149	Total	O	0	0
			149	149		
2	B	166	Total	O	0	0
			166	166		
2	C	139	Total	O	0	0
			139	139		
2	D	173	Total	O	0	0
			173	173		

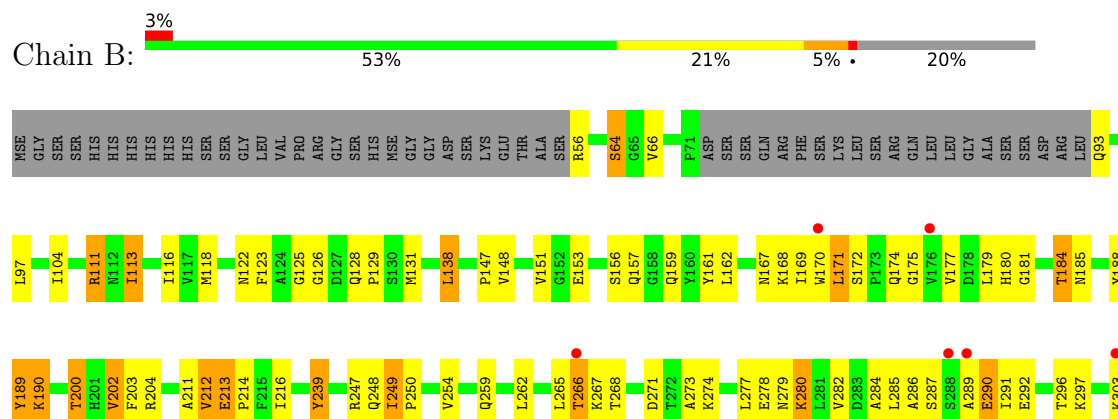
3 Residue-property plots

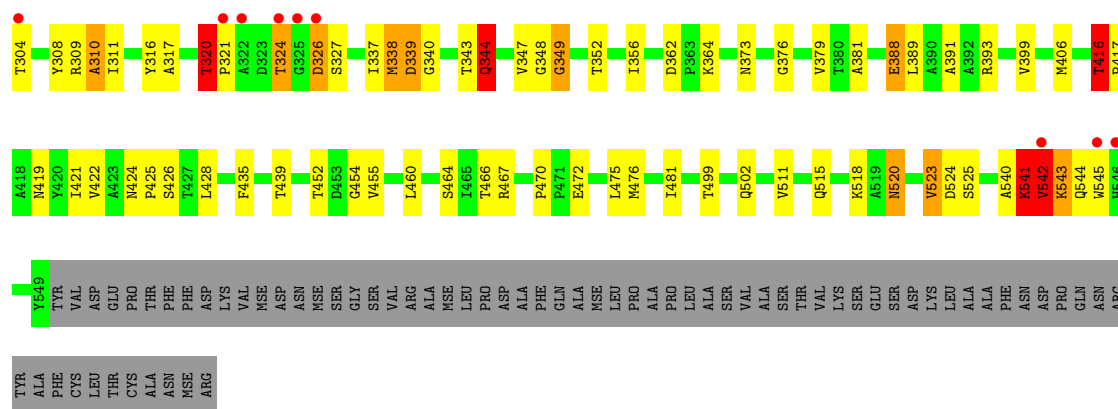
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Protease 4

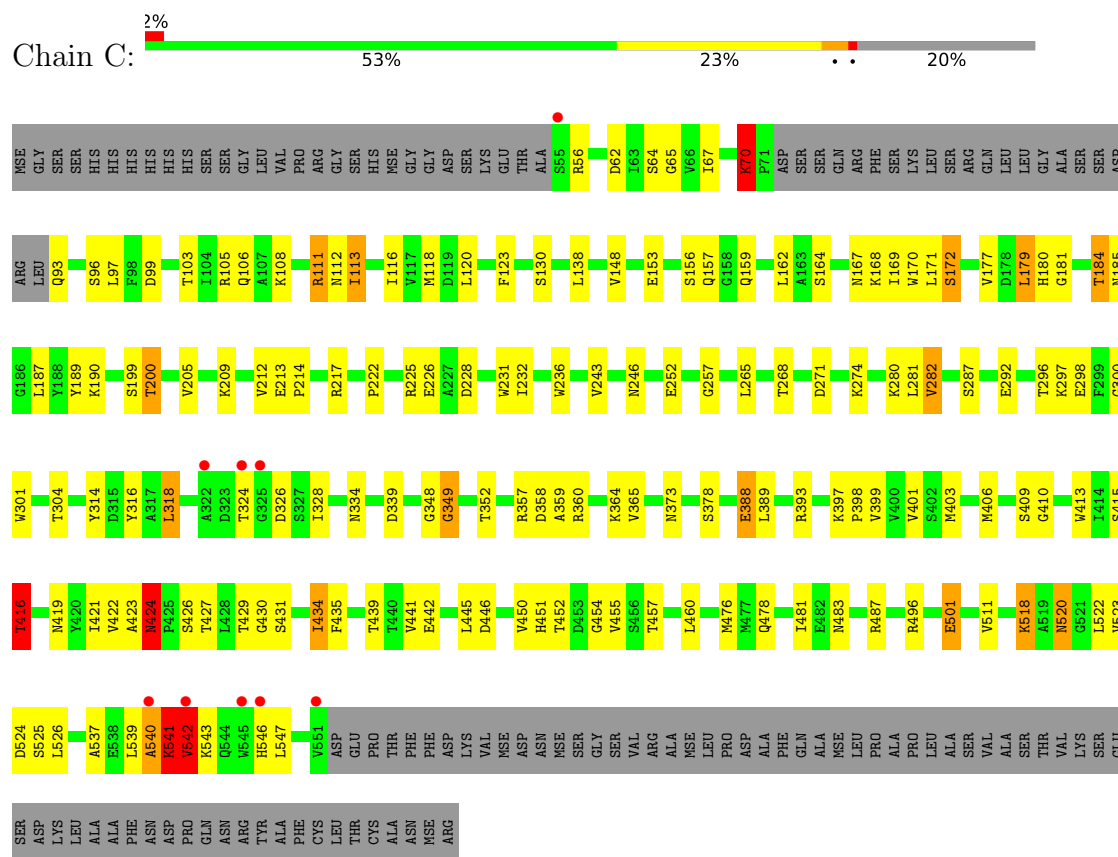


• Molecule 1: Protease 4

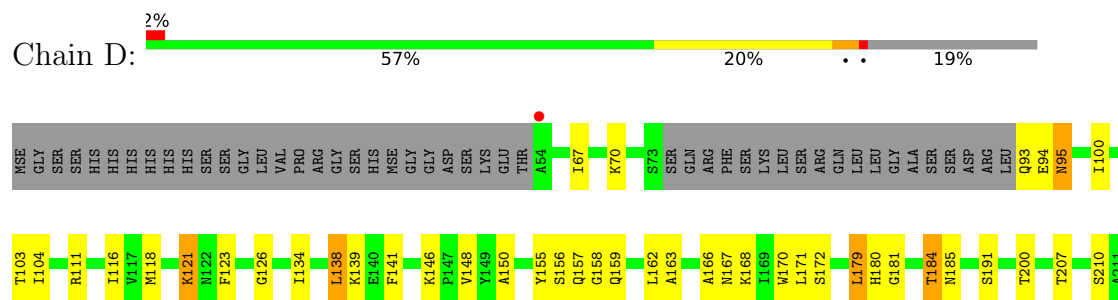


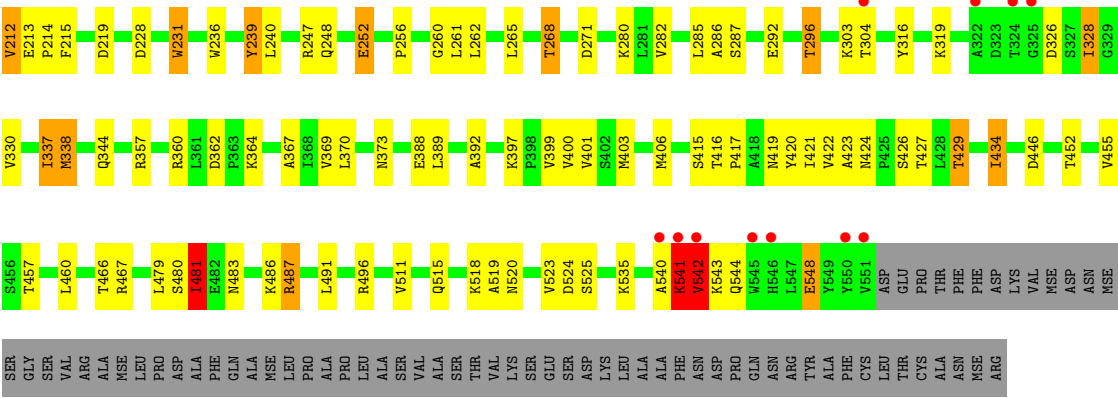


• Molecule 1: Protease 4



• Molecule 1: Protease 4





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	93.02Å 153.04Å 100.21Å 90.00° 104.17° 90.00°	Depositor
Resolution (Å)	46.63 – 2.76 46.63 – 2.76	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.63-2.76) 98.9 (46.63-2.76)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	0.18	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.84 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.200 , 0.259 (Not available) , 0.224	Depositor DCC
R_{free} test set	3501 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	29.4	Xtriage
Anisotropy	0.387	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 40.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	14989	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

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	1.02	1/3635 (0.0%)	1.16	9/4922 (0.2%)
1	B	1.02	3/3631 (0.1%)	1.20	18/4918 (0.4%)
1	C	1.00	1/3656 (0.0%)	1.19	12/4950 (0.2%)
1	D	1.01	3/3676 (0.1%)	1.16	11/4977 (0.2%)
All	All	1.01	8/14598 (0.1%)	1.18	50/19767 (0.3%)

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	2
1	B	0	2
1	C	0	2
1	D	0	1
All	All	0	7

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	542	VAL	CA-CB	10.65	1.69	1.54
1	A	542	VAL	CA-CB	9.63	1.67	1.54
1	D	542	VAL	CA-CB	9.10	1.66	1.54
1	D	429	THR	CA-CB	6.19	1.62	1.54
1	D	481	ILE	CA-CB	5.77	1.61	1.54
1	C	542	VAL	CA-CB	5.66	1.62	1.54
1	B	169	ILE	CA-CB	5.37	1.60	1.53
1	B	470	PRO	CA-C	5.04	1.55	1.52

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	416	THR	CA-C-N	-10.77	109.85	120.83
1	A	416	THR	C-N-CA	-10.77	109.85	120.83
1	A	57	GLY	N-CA-C	8.43	120.59	112.08
1	C	541	LYS	N-CA-C	8.24	121.92	108.99
1	C	542	VAL	N-CA-C	8.22	126.43	109.34
1	D	303	LYS	N-CA-C	-7.30	103.35	111.82
1	D	541	LYS	N-CA-C	7.05	119.24	108.52
1	B	518	LYS	N-CA-C	-6.77	103.14	111.40
1	B	213	GLU	CA-C-N	6.77	127.32	119.47
1	B	213	GLU	C-N-CA	6.77	127.32	119.47
1	B	339	ASP	N-CA-C	6.73	120.71	112.23
1	A	372	VAL	N-CA-C	6.45	117.17	107.75
1	C	416	THR	CA-C-N	-6.38	113.60	121.00
1	C	416	THR	C-N-CA	-6.38	113.60	121.00
1	B	324	THR	N-CA-C	-6.25	104.38	111.07
1	D	483	ASN	N-CA-C	6.23	117.73	111.07
1	B	542	VAL	N-CA-C	6.11	122.06	109.34
1	C	546	HIS	N-CA-C	6.11	119.19	109.24
1	B	416	THR	CA-C-N	-5.99	113.80	120.45
1	B	416	THR	C-N-CA	-5.99	113.80	120.45
1	A	484	GLY	N-CA-C	-5.91	107.34	114.66
1	B	286	ALA	N-CA-C	5.79	117.82	108.32
1	A	542	VAL	N-CA-C	5.76	121.33	109.34
1	D	542	VAL	N-CA-C	5.75	121.29	109.34
1	D	158	GLY	N-CA-C	-5.68	105.75	113.37
1	D	126	GLY	N-CA-C	-5.57	101.11	111.14
1	C	541	LYS	CA-C-N	5.56	131.98	121.97
1	C	541	LYS	C-N-CA	5.56	131.98	121.97
1	D	231	TRP	N-CA-C	5.54	117.40	111.36
1	C	424	ASN	CA-C-N	5.53	125.89	119.47
1	C	424	ASN	C-N-CA	5.53	125.89	119.47
1	D	70	LYS	CA-C-N	5.52	125.44	119.76
1	D	70	LYS	C-N-CA	5.52	125.44	119.76
1	C	281	LEU	N-CA-C	-5.46	106.12	112.89
1	A	396	GLY	N-CA-C	5.41	122.24	115.21
1	C	257	GLY	N-CA-C	-5.40	104.82	112.81
1	B	291	ILE	N-CA-C	-5.32	105.66	110.82
1	B	239	TYR	N-CA-C	-5.31	105.18	110.97
1	C	65	GLY	N-CA-C	5.27	116.88	111.56
1	A	141	PHE	N-CA-C	-5.27	105.44	111.07
1	B	381	ALA	N-CA-C	-5.25	105.63	111.36
1	A	389	LEU	N-CA-C	-5.24	105.64	111.36
1	B	515	GLN	N-CA-C	5.24	116.67	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	97	LEU	N-CA-C	-5.21	105.50	111.07
1	B	320	THR	CA-C-N	5.16	125.15	119.89
1	B	320	THR	C-N-CA	5.16	125.15	119.89
1	B	541	LYS	N-CA-C	5.15	116.35	108.52
1	D	239	TYR	N-CA-C	-5.15	105.36	110.97
1	D	486	LYS	N-CA-C	-5.12	105.39	111.69
1	B	211	ALA	N-CA-C	-5.03	106.98	113.01

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	541	LYS	Peptide
1	A	544	GLN	Peptide
1	B	340	GLY	Peptide
1	B	541	LYS	Peptide
1	C	541	LYS	Peptide
1	C	70	LYS	Peptide
1	D	541	LYS	Peptide

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	3576	0	3537	110	0
1	B	3572	0	3526	114	0
1	C	3597	0	3554	131	0
1	D	3617	0	3571	97	0
2	A	149	0	0	0	0
2	B	166	0	0	6	0
2	C	139	0	0	5	0
2	D	173	0	0	2	0
All	All	14989	0	14188	425	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 15.

All (425) 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:B:123:PHE:CZ	1:B:131:MSE:HE2	1.73	1.21
1:A:403:MSE:HE2	1:A:423:ALA:HB2	1.27	1.17
1:C:403:MSE:HE2	1:C:423:ALA:HB2	1.28	1.15
1:A:541:LYS:HD3	1:A:542:VAL:HG23	1.24	1.14
1:D:121:LYS:HE2	2:D:691:HOH:O	1.55	1.05
1:D:403:MSE:HE2	1:D:423:ALA:HB2	1.42	1.02
1:D:424:ASN:HD22	1:D:426:SER:H	1.03	1.01
1:A:424:ASN:HD22	1:A:426:SER:H	1.01	1.00
1:B:123:PHE:HZ	1:B:131:MSE:HE2	1.10	0.99
1:D:148:VAL:H	1:D:167:ASN:HD22	1.06	0.98
1:B:424:ASN:ND2	1:B:426:SER:H	1.62	0.97
1:C:424:ASN:HD22	1:C:426:SER:H	0.99	0.96
1:A:148:VAL:H	1:A:167:ASN:HD22	1.06	0.96
1:C:148:VAL:H	1:C:167:ASN:HD22	1.08	0.95
1:A:148:VAL:H	1:A:167:ASN:ND2	1.62	0.95
1:B:148:VAL:H	1:B:167:ASN:HD22	0.97	0.94
1:C:148:VAL:H	1:C:167:ASN:ND2	1.65	0.94
1:B:202:VAL:HG21	1:B:216:ILE:HG21	1.48	0.93
1:A:184:THR:HG22	1:C:455:VAL:H	1.32	0.92
1:C:268:THR:CG2	1:C:271:ASP:HB3	1.99	0.92
1:C:424:ASN:ND2	1:C:426:SER:H	1.67	0.91
1:D:157:GLN:HE22	1:D:181:GLY:H	1.19	0.91
1:B:148:VAL:H	1:B:167:ASN:ND2	1.69	0.90
1:D:157:GLN:NE2	1:D:180:HIS:H	1.68	0.90
1:C:541:LYS:HB3	1:C:542:VAL:HG22	1.51	0.90
1:B:157:GLN:HE22	1:B:181:GLY:H	1.13	0.89
1:A:116:ILE:HB	1:A:148:VAL:HG22	1.53	0.88
1:B:540:ALA:O	1:B:541:LYS:HG3	1.75	0.87
1:A:545:TRP:CZ3	1:A:547:LEU:HD23	2.09	0.87
1:B:424:ASN:HD22	1:B:426:SER:H	0.90	0.86
1:B:123:PHE:CZ	1:B:131:MSE:CE	2.59	0.86
1:B:157:GLN:NE2	1:B:180:HIS:H	1.75	0.85
1:A:399:VAL:H	1:A:419:ASN:ND2	1.75	0.83
1:B:118:MSE:HE1	1:B:162:LEU:HD22	1.60	0.83
1:B:424:ASN:HD22	1:B:426:SER:N	1.74	0.83
1:C:424:ASN:HD22	1:C:426:SER:N	1.76	0.82
1:A:171:LEU:O	1:A:285:LEU:HA	1.79	0.82
1:C:184:THR:HG22	1:D:455:VAL:H	1.45	0.82
1:A:403:MSE:HE2	1:A:423:ALA:CB	2.10	0.81
1:B:123:PHE:CE2	1:B:131:MSE:HE2	2.16	0.81
1:D:403:MSE:HE2	1:D:423:ALA:CB	2.10	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:403:MSE:HE1	1:D:429:THR:HG21	1.62	0.80
1:B:104:ILE:HD13	1:B:138:LEU:HD13	1.65	0.79
1:B:202:VAL:HG21	1:B:216:ILE:CG2	2.12	0.79
1:A:184:THR:CG2	1:C:455:VAL:H	1.96	0.78
1:C:168:LYS:HE2	1:C:170:TRP:CZ2	2.18	0.78
1:D:399:VAL:H	1:D:419:ASN:HD22	1.32	0.78
1:B:123:PHE:HZ	1:B:131:MSE:CE	1.95	0.78
1:B:287:SER:HB3	1:B:290:GLU:HG3	1.66	0.77
1:D:424:ASN:ND2	1:D:426:SER:H	1.81	0.77
1:C:403:MSE:HE2	1:C:423:ALA:CB	2.13	0.76
1:A:545:TRP:HZ3	1:A:547:LEU:HD23	1.47	0.76
1:B:287:SER:HB3	1:B:290:GLU:CG	2.16	0.76
1:A:370:LEU:HD22	1:A:389:LEU:HD21	1.67	0.74
1:C:157:GLN:NE2	1:C:180:HIS:H	1.85	0.74
1:A:131:MSE:HE2	1:A:162:LEU:HG	1.69	0.74
1:B:156:SER:H	1:B:159:GLN:NE2	1.86	0.74
1:D:403:MSE:HE3	1:D:427:THR:HG21	1.71	0.72
1:A:157:GLN:HE22	1:A:181:GLY:H	1.37	0.72
1:B:148:VAL:N	1:B:167:ASN:HD22	1.82	0.72
1:C:268:THR:HG22	1:C:271:ASP:HB3	1.70	0.72
1:A:123:PHE:CZ	1:A:131:MSE:HE3	2.25	0.72
1:A:360:ARG:HD3	1:A:392:ALA:HA	1.71	0.72
1:B:502:GLN:HG3	2:B:717:HOH:O	1.89	0.72
1:B:356:ILE:HD13	1:B:389:LEU:HD13	1.72	0.72
1:D:292:GLU:O	1:D:296:THR:HG22	1.90	0.72
1:B:399:VAL:H	1:B:419:ASN:ND2	1.87	0.71
1:C:157:GLN:HE22	1:C:181:GLY:H	1.39	0.70
1:A:403:MSE:CE	1:A:423:ALA:HB2	2.17	0.69
1:B:56:ARG:CB	2:B:655:HOH:O	2.41	0.69
1:C:111:ARG:HB3	1:C:111:ARG:HH11	1.57	0.69
1:C:157:GLN:HE22	1:C:180:HIS:H	1.37	0.69
1:D:148:VAL:H	1:D:167:ASN:ND2	1.86	0.69
1:A:157:GLN:NE2	1:A:180:HIS:H	1.91	0.68
1:C:148:VAL:N	1:C:167:ASN:HD22	1.88	0.68
1:C:156:SER:H	1:C:159:GLN:NE2	1.92	0.68
1:A:542:VAL:HG12	1:A:543:LYS:H	1.58	0.68
1:D:420:TYR:HA	1:D:524:ASP:OD2	1.94	0.68
1:A:370:LEU:HD21	1:A:414:ILE:HD12	1.76	0.67
1:B:277:LEU:HD13	1:B:285:LEU:HG	1.77	0.67
1:A:424:ASN:HD22	1:A:426:SER:N	1.84	0.67
1:B:455:VAL:H	1:D:184:THR:HG22	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:118:MSE:CE	1:B:162:LEU:HD22	2.25	0.66
1:B:213:GLU:N	1:B:214:PRO:HD2	2.11	0.66
1:C:96:SER:HB3	1:C:99:ASP:HB2	1.77	0.66
1:C:268:THR:HG21	1:C:271:ASP:O	1.96	0.66
1:A:446:ASP:OD1	1:A:451:HIS:HE1	1.79	0.65
1:C:540:ALA:O	1:C:541:LYS:HG3	1.95	0.65
1:A:292:GLU:O	1:A:296:THR:HG23	1.97	0.65
1:D:399:VAL:H	1:D:419:ASN:ND2	1.93	0.65
1:A:455:VAL:H	1:B:184:THR:HG22	1.61	0.65
1:C:399:VAL:H	1:C:419:ASN:HD22	1.42	0.65
1:A:148:VAL:N	1:A:167:ASN:HD22	1.89	0.65
1:C:185:ASN:HD21	1:D:452:THR:HB	1.61	0.65
1:C:292:GLU:CD	1:D:357:ARG:HH21	2.04	0.65
1:D:200:THR:HG21	1:D:215:PHE:HB3	1.78	0.65
1:B:499:THR:OG1	1:B:502:GLN:HG2	1.96	0.64
1:B:326:ASP:H	1:B:545:TRP:HE1	1.45	0.64
1:D:157:GLN:HE22	1:D:180:HIS:H	1.42	0.64
1:A:399:VAL:H	1:A:419:ASN:HD22	1.44	0.63
1:B:337:ILE:HG12	1:B:376:GLY:O	1.99	0.63
1:D:373:ASN:HA	1:D:406:MSE:O	1.99	0.63
1:B:520:ASN:C	1:B:520:ASN:HD22	2.07	0.63
1:C:177:VAL:HG12	1:C:179:LEU:HD13	1.81	0.63
1:C:225:ARG:NH2	2:C:673:HOH:O	2.28	0.63
1:A:424:ASN:ND2	1:A:426:SER:H	1.85	0.63
1:C:398:PRO:HG2	1:C:539:LEU:O	1.98	0.63
1:B:200:THR:HG23	1:B:439:THR:HG22	1.81	0.62
1:C:111:ARG:HB3	1:C:111:ARG:NH1	2.15	0.62
1:A:487:ARG:HD3	1:B:174:GLN:HG3	1.82	0.62
1:C:373:ASN:HA	1:C:406:MSE:O	2.00	0.61
1:A:123:PHE:HZ	1:A:131:MSE:HE3	1.63	0.61
1:D:156:SER:H	1:D:159:GLN:NE2	1.98	0.61
1:A:481:ILE:C	1:A:483:ASN:H	2.08	0.61
1:B:126:GLY:N	1:B:131:MSE:HE3	2.15	0.61
1:C:403:MSE:HE3	1:C:427:THR:HG21	1.82	0.61
1:C:483:ASN:O	1:C:487:ARG:HG3	2.01	0.61
1:D:434:ILE:HG23	1:D:481:ILE:HG23	1.81	0.61
1:D:370:LEU:HD22	1:D:389:LEU:HD21	1.83	0.60
1:B:296:THR:HG22	1:B:308:TYR:HB3	1.84	0.60
1:D:424:ASN:HD22	1:D:426:SER:N	1.86	0.60
1:B:213:GLU:N	1:B:214:PRO:CD	2.64	0.60
1:A:157:GLN:HE22	1:A:180:HIS:H	1.48	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:343:THR:O	1:B:344:GLN:HB2	2.01	0.59
1:B:273:ALA:HB1	1:B:285:LEU:HD22	1.83	0.59
1:C:268:THR:HG21	1:C:271:ASP:HB3	1.84	0.59
1:D:452:THR:OG1	1:D:467:ARG:HD2	2.02	0.59
1:A:174:GLN:OE1	1:C:487:ARG:HD2	2.02	0.59
1:B:268:THR:HG22	1:B:271:ASP:HB3	1.85	0.59
1:B:309:ARG:O	1:B:310:ALA:HB2	2.02	0.59
1:B:66:VAL:HG22	1:B:93:GLN:O	2.03	0.59
1:A:520:ASN:C	1:A:520:ASN:HD22	2.12	0.58
1:C:148:VAL:N	1:C:167:ASN:ND2	2.46	0.58
1:A:360:ARG:HD3	1:A:392:ALA:CA	2.33	0.58
1:A:542:VAL:CG1	1:A:543:LYS:N	2.67	0.57
1:B:399:VAL:H	1:B:419:ASN:HD22	1.52	0.57
1:A:184:THR:HG22	1:A:184:THR:O	2.04	0.57
1:C:328:ILE:HD11	1:C:537:ALA:HB2	1.86	0.57
1:B:268:THR:HG22	1:B:268:THR:O	2.04	0.57
1:B:393:ARG:HD2	1:B:419:ASN:ND2	2.20	0.57
1:C:410:GLY:O	1:C:413:TRP:HB3	2.05	0.57
1:D:268:THR:HG23	1:D:271:ASP:H	1.70	0.57
1:D:228:ASP:HA	1:D:231:TRP:NE1	2.19	0.57
1:A:168:LYS:HE2	1:A:170:TRP:CZ2	2.39	0.57
1:A:545:TRP:CH2	1:A:547:LEU:HD23	2.39	0.57
1:A:184:THR:HG22	1:C:455:VAL:N	2.10	0.57
1:C:501:GLU:H	1:C:501:GLU:CD	2.13	0.57
1:A:454:GLY:HA3	1:B:184:THR:O	2.05	0.56
1:B:123:PHE:CE2	1:B:131:MSE:CE	2.88	0.56
1:A:184:THR:CG2	1:A:184:THR:O	2.51	0.56
1:D:103:THR:HG23	1:D:316:TYR:CD2	2.41	0.56
1:C:403:MSE:CE	1:C:423:ALA:HB2	2.19	0.56
1:D:148:VAL:N	1:D:167:ASN:HD22	1.90	0.56
1:A:163:ALA:C	1:A:165:PHE:H	2.13	0.56
1:D:292:GLU:O	1:D:296:THR:CG2	2.54	0.55
1:B:171:LEU:O	1:B:285:LEU:HA	2.07	0.55
1:B:452:THR:HB	1:D:185:ASN:ND2	2.21	0.55
1:A:228:ASP:HA	1:A:231:TRP:NE1	2.22	0.55
1:A:236:TRP:CB	1:C:457:THR:HG21	2.36	0.55
1:B:212:VAL:HG23	1:B:435:PHE:CD2	2.42	0.55
1:A:157:GLN:HE21	1:A:179:LEU:HA	1.71	0.55
1:D:369:VAL:HA	1:D:400:VAL:O	2.07	0.55
1:A:292:GLU:O	1:A:296:THR:CG2	2.56	0.54
1:C:399:VAL:H	1:C:419:ASN:ND2	2.04	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:370:LEU:HB3	1:D:401:VAL:HG22	1.89	0.54
1:B:116:ILE:HB	1:B:148:VAL:HG22	1.88	0.54
1:B:151:VAL:HG12	1:B:170:TRP:HB2	1.89	0.54
1:D:420:TYR:CD2	1:D:535:LYS:HG2	2.43	0.54
1:D:116:ILE:HB	1:D:148:VAL:HG22	1.90	0.54
1:B:148:VAL:N	1:B:167:ASN:ND2	2.48	0.54
1:B:373:ASN:HA	1:B:406:MSE:O	2.08	0.54
1:A:373:ASN:HA	1:A:406:MSE:O	2.08	0.54
1:C:232:ILE:HG21	1:D:455:VAL:HG23	1.89	0.54
1:C:222:PRO:O	1:C:226:GLU:HG2	2.07	0.53
1:C:106:GLN:NE2	1:C:316:TYR:CE2	2.76	0.53
1:C:520:ASN:C	1:C:520:ASN:HD22	2.16	0.53
1:A:370:LEU:HD22	1:A:389:LEU:CD2	2.37	0.53
1:B:184:THR:HG22	1:B:184:THR:O	2.08	0.53
1:C:67:ILE:HG12	1:C:123:PHE:HE1	1.73	0.53
1:A:542:VAL:HG12	1:A:543:LYS:N	2.24	0.52
1:B:454:GLY:HA3	1:D:184:THR:O	2.09	0.52
1:C:153:GLU:O	1:C:172:SER:HB3	2.09	0.52
1:D:252:GLU:HB3	2:D:718:HOH:O	2.09	0.52
1:C:228:ASP:HA	1:C:231:TRP:NE1	2.25	0.52
1:C:268:THR:HG22	1:C:268:THR:O	2.07	0.52
1:B:352:THR:O	1:B:356:ILE:HG13	2.09	0.52
1:C:113:ILE:CD1	1:C:316:TYR:HE1	2.23	0.52
1:B:309:ARG:NH2	2:B:784:HOH:O	2.43	0.52
1:D:93:GLN:HG2	1:D:94:GLU:N	2.25	0.52
1:D:268:THR:HG21	1:D:271:ASP:O	2.09	0.52
1:C:547:LEU:HD23	2:C:688:HOH:O	2.10	0.51
1:C:205:VAL:HG23	1:C:481:ILE:HG22	1.92	0.51
1:A:163:ALA:C	1:A:165:PHE:N	2.68	0.51
1:C:297:LYS:O	1:C:297:LYS:HG2	2.10	0.51
1:C:501:GLU:CD	1:C:501:GLU:N	2.69	0.51
1:A:131:MSE:HE2	1:A:162:LEU:CG	2.40	0.51
1:A:431:SER:HA	1:A:509:GLY:O	2.11	0.51
1:B:452:THR:HB	1:D:185:ASN:HD21	1.75	0.51
1:C:268:THR:HG23	1:C:274:LYS:CD	2.40	0.51
1:C:518:LYS:HG2	1:C:526:LEU:HG	1.92	0.51
1:B:421:ILE:N	1:B:524:ASP:OD2	2.35	0.51
1:C:200:THR:HG23	1:C:439:THR:HG22	1.93	0.51
1:C:217:ARG:HD2	2:C:677:HOH:O	2.09	0.51
1:A:236:TRP:HB2	1:C:457:THR:HG21	1.93	0.50
1:D:213:GLU:N	1:D:214:PRO:CD	2.74	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:THR:HG23	1:A:316:TYR:CD2	2.47	0.50
1:C:268:THR:HG23	1:C:274:LYS:HD3	1.93	0.50
1:C:348:GLY:O	1:C:349:GLY:C	2.54	0.50
1:B:249:ILE:HD11	1:B:254:VAL:HG22	1.91	0.50
1:B:545:TRP:HZ2	2:B:689:HOH:O	1.94	0.50
1:C:184:THR:HG22	1:D:455:VAL:N	2.20	0.50
1:A:324:THR:C	1:A:545:TRP:HE1	2.20	0.50
1:B:416:THR:N	1:B:417:PRO:CD	2.75	0.50
1:D:95:ASN:N	1:D:95:ASN:ND2	2.59	0.50
1:A:446:ASP:OD1	1:A:451:HIS:CE1	2.61	0.50
1:C:105:ARG:O	1:C:108:LYS:HB3	2.11	0.50
1:D:487:ARG:HG2	1:D:487:ARG:HH11	1.76	0.50
1:B:125:GLY:HA2	1:B:131:MSE:CE	2.42	0.50
1:D:100:ILE:HG21	1:D:134:ILE:HD13	1.94	0.50
1:A:189:TYR:O	1:A:190:LYS:C	2.54	0.50
1:A:225:ARG:NH2	1:C:454:GLY:O	2.44	0.50
1:D:67:ILE:HG12	1:D:123:PHE:HE1	1.77	0.50
1:B:327:SER:HB3	1:B:545:TRP:CE2	2.47	0.50
1:C:200:THR:CG2	1:C:439:THR:HG22	2.42	0.50
1:B:177:VAL:HG12	1:B:179:LEU:HD13	1.93	0.49
1:A:341:GLU:O	1:A:342:GLU:C	2.55	0.49
1:C:451:HIS:HA	2:C:654:HOH:O	2.11	0.49
1:D:403:MSE:HE3	1:D:427:THR:CG2	2.42	0.49
1:A:164:SER:O	1:A:246:ASN:HB3	2.12	0.49
1:D:328:ILE:HD13	1:D:367:ALA:HB3	1.93	0.49
1:A:403:MSE:HE3	1:A:427:THR:HG21	1.95	0.49
1:D:415:SER:C	1:D:417:PRO:HD2	2.38	0.49
1:B:161:TYR:HB2	1:B:239:TYR:CE2	2.47	0.49
1:B:324:THR:HB	1:B:545:TRP:CD1	2.48	0.49
1:C:164:SER:HA	1:C:169:ILE:HD11	1.95	0.49
1:C:168:LYS:CE	1:C:170:TRP:CZ2	2.94	0.49
1:C:296:THR:HG21	1:C:301:TRP:HE3	1.77	0.49
1:D:540:ALA:O	1:D:541:LYS:HG3	2.12	0.49
1:A:268:THR:HG21	1:A:271:ASP:O	2.13	0.49
1:A:432:ILE:HB	1:A:507:ALA:HB1	1.94	0.49
1:C:97:LEU:HD22	1:C:130:SER:OG	2.13	0.49
1:C:118:MSE:HE1	1:C:162:LEU:HB3	1.95	0.49
1:C:430:GLY:O	1:C:431:SER:CB	2.61	0.49
1:B:452:THR:HG21	1:B:466:THR:OG1	2.12	0.49
1:A:337:ILE:HA	1:A:347:VAL:HG12	1.95	0.48
1:D:403:MSE:HE1	1:D:429:THR:CG2	2.36	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:359:ALA:HB1	1:C:365:VAL:HG11	1.94	0.48
1:C:430:GLY:O	1:C:431:SER:HB3	2.14	0.48
1:B:424:ASN:CG	1:B:425:PRO:HD2	2.37	0.48
1:D:362:ASP:O	1:D:397:LYS:NZ	2.46	0.48
1:D:421:ILE:HD12	1:D:496:ARG:NH1	2.28	0.48
1:B:153:GLU:O	1:B:175:GLY:HA3	2.13	0.48
1:C:168:LYS:HE2	1:C:170:TRP:CH2	2.49	0.48
1:A:231:TRP:O	1:A:235:LEU:HD13	2.13	0.48
1:A:232:ILE:HG21	1:C:455:VAL:HG23	1.95	0.48
1:A:481:ILE:C	1:A:483:ASN:N	2.72	0.48
1:B:545:TRP:CZ2	2:B:689:HOH:O	2.55	0.48
1:C:164:SER:O	1:C:246:ASN:HB3	2.13	0.48
1:C:189:TYR:O	1:C:190:LYS:C	2.57	0.47
1:A:383:GLU:HA	1:A:383:GLU:OE1	2.14	0.47
1:C:113:ILE:CD1	1:C:316:TYR:CE1	2.96	0.47
1:D:141:PHE:CE1	1:D:146:LYS:HB2	2.49	0.47
1:B:184:THR:O	1:B:184:THR:CG2	2.61	0.47
1:A:153:GLU:HA	1:A:172:SER:HB3	1.96	0.47
1:B:424:ASN:ND2	1:B:426:SER:N	2.46	0.47
1:C:157:GLN:HE21	1:C:179:LEU:HA	1.80	0.47
1:A:103:THR:HG23	1:A:316:TYR:HD2	1.80	0.47
1:A:210:SER:O	1:A:213:GLU:HB2	2.15	0.47
1:C:403:MSE:HE1	1:C:429:THR:HG21	1.97	0.47
1:B:118:MSE:HE3	1:B:138:LEU:HD21	1.97	0.47
1:B:421:ILE:CG2	1:B:523:VAL:HG22	2.45	0.47
1:A:292:GLU:OE1	1:C:357:ARG:NE	2.35	0.47
1:A:424:ASN:ND2	1:A:425:PRO:HD2	2.30	0.47
1:D:168:LYS:HD3	1:D:170:TRP:CH2	2.50	0.46
1:C:113:ILE:HD13	1:C:316:TYR:CE1	2.50	0.46
1:C:213:GLU:N	1:C:214:PRO:HD2	2.30	0.46
1:A:185:ASN:HD21	1:C:452:THR:HB	1.81	0.46
1:C:268:THR:HG22	1:C:271:ASP:CB	2.42	0.46
1:A:545:TRP:HZ3	1:A:547:LEU:CD2	2.22	0.46
1:B:118:MSE:CE	1:B:138:LEU:HD21	2.46	0.46
1:A:131:MSE:HE2	1:A:162:LEU:HB2	1.98	0.46
1:A:154:ASN:OD1	1:A:175:GLY:HA2	2.16	0.46
1:B:472:GLU:O	1:B:476:MSE:HG2	2.16	0.46
1:D:228:ASP:HA	1:D:231:TRP:CD1	2.49	0.46
1:D:360:ARG:HD3	1:D:392:ALA:HA	1.98	0.46
1:B:189:TYR:O	1:B:190:LYS:C	2.57	0.46
1:B:348:GLY:O	1:B:349:GLY:C	2.59	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:518:LYS:O	1:D:519:ALA:C	2.57	0.46
1:A:266:THR:C	1:A:268:THR:H	2.23	0.46
1:C:296:THR:CG2	1:C:301:TRP:HE3	2.29	0.46
1:C:435:PHE:C	1:C:435:PHE:CD2	2.93	0.46
1:C:116:ILE:HB	1:C:148:VAL:HG22	1.97	0.45
1:D:256:PRO:HB2	1:D:260:GLY:HA3	1.98	0.45
1:B:268:THR:CG2	1:B:274:LYS:HB3	2.46	0.45
1:B:266:THR:C	1:B:268:THR:H	2.24	0.45
1:B:467:ARG:NH2	2:B:720:HOH:O	2.46	0.45
1:D:210:SER:C	1:D:212:VAL:H	2.24	0.45
1:A:360:ARG:O	1:A:360:ARG:HG3	2.17	0.45
1:B:204:ARG:HD3	1:B:212:VAL:HG11	1.98	0.45
1:C:62:ASP:O	1:C:314:TYR:HE2	2.00	0.45
1:B:327:SER:HB3	1:B:545:TRP:NE1	2.31	0.45
1:D:157:GLN:HE21	1:D:179:LEU:HA	1.81	0.45
1:A:101:VAL:HG21	1:A:133:TYR:CE2	2.52	0.45
1:C:213:GLU:N	1:C:214:PRO:CD	2.80	0.45
1:C:56:ARG:H	1:C:112:ASN:HB3	1.82	0.45
1:C:296:THR:HG21	1:C:301:TRP:CE3	2.51	0.45
1:C:393:ARG:HA	1:C:397:LYS:O	2.17	0.45
1:C:413:TRP:O	1:C:416:THR:HB	2.17	0.45
1:A:479:LEU:O	1:A:483:ASN:HB2	2.17	0.44
1:B:113:ILE:HD13	1:B:316:TYR:CE1	2.52	0.44
1:A:227:ALA:O	1:A:231:TRP:HD1	2.00	0.44
1:B:203:PHE:O	1:B:435:PHE:HA	2.17	0.44
1:C:187:LEU:HD23	1:D:452:THR:HG22	1.99	0.44
1:D:330:VAL:HB	1:D:548:GLU:HG3	2.00	0.44
1:A:118:MSE:CE	1:A:166:ALA:HB2	2.47	0.44
1:A:324:THR:HG22	1:A:325:GLY:H	1.82	0.44
1:C:520:ASN:ND2	1:C:522:LEU:HG	2.33	0.44
1:A:228:ASP:HA	1:A:231:TRP:CD1	2.53	0.44
1:B:388:GLU:O	1:B:391:ALA:HB3	2.17	0.44
1:C:539:LEU:O	1:C:540:ALA:HB2	2.18	0.44
1:D:219:ASP:C	1:D:219:ASP:OD1	2.60	0.44
1:D:268:THR:HG23	1:D:271:ASP:N	2.31	0.44
1:B:249:ILE:HB	1:B:250:PRO:HD2	1.98	0.44
1:D:168:LYS:HE2	1:D:170:TRP:CZ2	2.53	0.44
1:D:236:TRP:O	1:D:239:TYR:HB3	2.18	0.44
1:D:240:LEU:HD23	1:D:240:LEU:HA	1.67	0.44
1:A:60:LEU:HB3	1:A:312:SER:HA	1.99	0.44
1:D:157:GLN:HE21	1:D:180:HIS:H	1.57	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:369:VAL:HA	1:A:400:VAL:O	2.18	0.44
1:B:170:TRP:CD1	1:B:284:ALA:HB3	2.52	0.44
1:B:279:ASN:O	1:B:280:LYS:HB2	2.16	0.43
1:C:103:THR:OG1	1:C:318:LEU:HD23	2.17	0.43
1:C:106:GLN:HE21	1:C:316:TYR:HE2	1.66	0.43
1:D:541:LYS:C	1:D:542:VAL:HG22	2.44	0.43
1:A:420:TYR:CD2	1:A:535:LYS:HG2	2.53	0.43
1:C:434:ILE:HG23	1:C:481:ILE:HG23	2.01	0.43
1:B:64:SER:HA	1:B:122:ASN:O	2.19	0.43
1:B:104:ILE:HD13	1:B:138:LEU:CD1	2.42	0.43
1:C:120:LEU:HA	1:C:120:LEU:HD23	1.75	0.43
1:A:107:ALA:HB1	1:A:113:ILE:HG21	2.01	0.43
1:A:208:TYR:HB2	1:A:510:HIS:N	2.34	0.43
1:C:70:LYS:HB2	1:C:334:ASN:HD22	1.84	0.43
1:D:162:LEU:O	1:D:163:ALA:C	2.62	0.43
1:A:200:THR:HG21	1:A:215:PHE:HB3	2.01	0.43
1:A:337:ILE:HD13	1:A:347:VAL:HG11	2.01	0.43
1:B:111:ARG:HH11	1:B:111:ARG:HB3	1.84	0.43
1:B:320:THR:HA	1:B:321:PRO:HD3	1.86	0.43
1:D:210:SER:C	1:D:212:VAL:N	2.76	0.43
1:A:67:ILE:HG12	1:A:123:PHE:HE1	1.84	0.42
1:A:528:ASP:O	1:A:529:PHE:C	2.62	0.42
1:C:70:LYS:HB2	1:C:334:ASN:ND2	2.34	0.42
1:C:185:ASN:ND2	1:D:452:THR:HB	2.33	0.42
1:C:389:LEU:HD12	1:C:389:LEU:HA	1.88	0.42
1:C:445:LEU:HB3	1:C:450:VAL:HB	2.00	0.42
1:D:104:ILE:HD13	1:D:138:LEU:HD13	2.00	0.42
1:D:338:MSE:HE2	1:D:338:MSE:HA	2.01	0.42
1:A:302:SER:C	1:A:304:THR:N	2.73	0.42
1:B:289:ALA:HA	1:B:292:GLU:HB2	1.99	0.42
1:B:337:ILE:HA	1:B:347:VAL:HB	2.01	0.42
1:C:118:MSE:HE2	1:C:120:LEU:HD11	2.01	0.42
1:A:428:LEU:HD12	1:A:428:LEU:HA	1.81	0.42
1:D:491:LEU:HA	1:D:491:LEU:HD23	1.76	0.42
1:A:143:ASP:C	1:A:145:GLY:H	2.26	0.42
1:B:540:ALA:C	1:B:541:LYS:HG3	2.44	0.42
1:D:118:MSE:CE	1:D:166:ALA:HB2	2.50	0.42
1:A:366:LYS:O	1:A:398:PRO:HD2	2.19	0.42
1:C:421:ILE:HG22	1:C:523:VAL:HG22	2.02	0.42
1:B:157:GLN:HE21	1:B:179:LEU:HA	1.85	0.42
1:C:357:ARG:O	1:C:360:ARG:HB3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:403:MSE:HE1	1:C:429:THR:CB	2.50	0.42
1:D:403:MSE:CE	1:D:423:ALA:HB2	2.30	0.42
1:D:67:ILE:HG12	1:D:123:PHE:CE1	2.54	0.42
1:D:155:TYR:HA	1:D:159:GLN:HE21	1.84	0.42
1:B:247:ARG:O	1:B:248:GLN:C	2.63	0.42
1:C:171:LEU:HB2	1:C:282:VAL:HG21	2.02	0.42
1:C:236:TRP:CB	1:D:457:THR:HG21	2.49	0.42
1:D:467:ARG:HE	1:D:467:ARG:HB3	1.66	0.42
1:A:213:GLU:OE2	1:A:221:SER:OG	2.36	0.42
1:B:188:TYR:CD1	1:B:188:TYR:N	2.87	0.42
1:C:476:MSE:HE3	1:C:476:MSE:HB3	1.93	0.42
1:C:401:VAL:HG21	1:C:415:SER:HA	2.01	0.41
1:A:231:TRP:CE3	1:A:428:LEU:HD21	2.55	0.41
1:A:382:SER:HB3	1:A:413:TRP:CD1	2.55	0.41
1:B:338:MSE:HE2	1:B:338:MSE:HA	2.03	0.41
1:D:542:VAL:HB	1:D:543:LYS:H	1.65	0.41
1:A:114:THR:O	1:A:147:PRO:HD2	2.19	0.41
1:B:455:VAL:N	1:D:184:THR:HG22	2.31	0.41
1:C:199:SER:O	1:C:439:THR:HA	2.21	0.41
1:C:496:ARG:HH22	1:C:524:ASP:CG	2.28	0.41
1:D:148:VAL:O	1:D:167:ASN:N	2.46	0.41
1:A:209:LYS:NZ	1:A:409:SER:HB2	2.35	0.41
1:B:128:GLN:HB3	1:B:129:PRO:HD3	2.03	0.41
1:B:520:ASN:C	1:B:520:ASN:ND2	2.75	0.41
1:C:70:LYS:HE3	1:C:334:ASN:HB3	2.02	0.41
1:C:409:SER:O	1:C:410:GLY:C	2.63	0.41
1:D:452:THR:HG21	1:D:466:THR:OG1	2.20	0.41
1:B:147:PRO:HA	1:B:167:ASN:ND2	2.36	0.41
1:D:172:SER:HA	1:D:286:ALA:O	2.20	0.41
1:B:249:ILE:HB	1:B:250:PRO:CD	2.51	0.41
1:B:362:ASP:OD1	1:B:362:ASP:C	2.63	0.41
1:C:388:GLU:O	1:C:389:LEU:C	2.64	0.41
1:C:441:VAL:O	1:C:442:GLU:C	2.63	0.41
1:D:181:GLY:HA2	1:D:236:TRP:CZ2	2.56	0.41
1:B:157:GLN:HE22	1:B:180:HIS:H	1.60	0.41
1:B:157:GLN:HE21	1:B:180:HIS:H	1.62	0.41
1:B:540:ALA:O	1:B:541:LYS:CG	2.59	0.41
1:C:111:ARG:HG2	1:C:112:ASN:N	2.36	0.41
1:C:209:LYS:NZ	1:C:409:SER:OG	2.53	0.41
1:C:292:GLU:OE1	1:D:357:ARG:NH2	2.44	0.41
1:D:337:ILE:O	1:D:338:MSE:HE3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:434:ILE:HD13	1:D:481:ILE:HD12	2.03	0.41
1:A:259:GLN:HE21	1:A:259:GLN:HA	1.86	0.41
1:A:397:LYS:HA	1:A:398:PRO:HD3	1.87	0.41
1:C:113:ILE:HD11	1:C:316:TYR:CE1	2.56	0.41
1:D:118:MSE:HE3	1:D:150:ALA:HB2	2.03	0.40
1:C:268:THR:CG2	1:C:268:THR:O	2.69	0.40
1:D:171:LEU:O	1:D:285:LEU:HA	2.21	0.40
1:A:249:ILE:HD11	1:A:254:VAL:HG22	2.03	0.40
1:A:302:SER:O	1:A:303:LYS:C	2.65	0.40
1:A:413:TRP:O	1:A:416:THR:HB	2.21	0.40
1:B:542:VAL:CG1	1:B:543:LYS:N	2.84	0.40
1:C:153:GLU:HG2	2:C:635:HOH:O	2.21	0.40
1:D:247:ARG:O	1:D:248:GLN:HB2	2.22	0.40
1:C:113:ILE:HD11	1:C:316:TYR:HE1	1.86	0.40
1:C:296:THR:O	1:C:300:GLY:HA2	2.22	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	469/593 (79%)	426 (91%)	39 (8%)	4 (1%)	14	28
1	B	469/593 (79%)	428 (91%)	32 (7%)	9 (2%)	6	12
1	C	472/593 (80%)	433 (92%)	33 (7%)	6 (1%)	9	18
1	D	475/593 (80%)	437 (92%)	37 (8%)	1 (0%)	43	64
All	All	1885/2372 (80%)	1724 (92%)	141 (8%)	20 (1%)	11	22

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	542	VAL
1	B	326	ASP
1	B	344	GLN
1	B	542	VAL
1	C	326	ASP
1	C	540	ALA
1	D	542	VAL
1	A	324	THR
1	A	340	GLY
1	C	349	GLY
1	C	542	VAL
1	B	267	LYS
1	B	297	LYS
1	C	70	LYS
1	A	342	GLU
1	B	317	ALA
1	C	324	THR
1	B	310	ALA
1	B	172	SER
1	B	349	GLY

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	373/460 (81%)	330 (88%)	43 (12%)	5	10
1	B	372/460 (81%)	325 (87%)	47 (13%)	4	8
1	C	374/460 (81%)	335 (90%)	39 (10%)	7	13
1	D	377/460 (82%)	333 (88%)	44 (12%)	5	9
All	All	1496/1840 (81%)	1323 (88%)	173 (12%)	5	10

All (173) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	LEU
1	A	93	GLN

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Mol	Chain	Res	Type
1	A	121	LYS
1	A	138	LEU
1	A	162	LEU
1	A	172	SER
1	A	179	LEU
1	A	184	THR
1	A	189	TYR
1	A	212	VAL
1	A	219	ASP
1	A	220	MSE
1	A	262	LEU
1	A	265	LEU
1	A	268	THR
1	A	281	LEU
1	A	282	VAL
1	A	292	GLU
1	A	296	THR
1	A	297	LYS
1	A	320	THR
1	A	324	THR
1	A	337	ILE
1	A	338	MSE
1	A	344	GLN
1	A	347	VAL
1	A	397	LYS
1	A	416	THR
1	A	422	VAL
1	A	428	LEU
1	A	434	ILE
1	A	460	LEU
1	A	479	LEU
1	A	481	ILE
1	A	486	LYS
1	A	490	THR
1	A	511	VAL
1	A	520	ASN
1	A	525	SER
1	A	541	LYS
1	A	542	VAL
1	A	543	LYS
1	A	547	LEU
1	B	64	SER

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Mol	Chain	Res	Type
1	B	111	ARG
1	B	113	ILE
1	B	138	LEU
1	B	168	LYS
1	B	171	LEU
1	B	184	THR
1	B	185	ASN
1	B	189	TYR
1	B	190	LYS
1	B	200	THR
1	B	202	VAL
1	B	212	VAL
1	B	249	ILE
1	B	259	GLN
1	B	262	LEU
1	B	265	LEU
1	B	266	THR
1	B	278	GLU
1	B	280	LYS
1	B	282	VAL
1	B	290	GLU
1	B	303	LYS
1	B	304	THR
1	B	311	ILE
1	B	320	THR
1	B	338	MSE
1	B	339	ASP
1	B	344	GLN
1	B	364	LYS
1	B	379	VAL
1	B	388	GLU
1	B	416	THR
1	B	422	VAL
1	B	428	LEU
1	B	460	LEU
1	B	464	SER
1	B	475	LEU
1	B	481	ILE
1	B	511	VAL
1	B	520	ASN
1	B	523	VAL
1	B	525	SER

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Mol	Chain	Res	Type
1	B	541	LYS
1	B	542	VAL
1	B	543	LYS
1	B	544	GLN
1	C	64	SER
1	C	93	GLN
1	C	111	ARG
1	C	113	ILE
1	C	138	LEU
1	C	172	SER
1	C	179	LEU
1	C	184	THR
1	C	200	THR
1	C	212	VAL
1	C	243	VAL
1	C	252	GLU
1	C	265	LEU
1	C	280	LYS
1	C	282	VAL
1	C	287	SER
1	C	298	GLU
1	C	304	THR
1	C	318	LEU
1	C	339	ASP
1	C	352	THR
1	C	358	ASP
1	C	364	LYS
1	C	378	SER
1	C	388	GLU
1	C	416	THR
1	C	422	VAL
1	C	424	ASN
1	C	434	ILE
1	C	446	ASP
1	C	460	LEU
1	C	478	GLN
1	C	501	GLU
1	C	511	VAL
1	C	518	LYS
1	C	520	ASN
1	C	525	SER
1	C	542	VAL

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Mol	Chain	Res	Type
1	C	543	LYS
1	D	95	ASN
1	D	111	ARG
1	D	121	LYS
1	D	138	LEU
1	D	139	LYS
1	D	179	LEU
1	D	184	THR
1	D	191	SER
1	D	207	THR
1	D	212	VAL
1	D	252	GLU
1	D	261	LEU
1	D	262	LEU
1	D	265	LEU
1	D	268	THR
1	D	280	LYS
1	D	282	VAL
1	D	287	SER
1	D	296	THR
1	D	304	THR
1	D	319	LYS
1	D	326	ASP
1	D	328	ILE
1	D	337	ILE
1	D	338	MSE
1	D	344	GLN
1	D	364	LYS
1	D	388	GLU
1	D	416	THR
1	D	422	VAL
1	D	434	ILE
1	D	446	ASP
1	D	460	LEU
1	D	479	LEU
1	D	480	SER
1	D	481	ILE
1	D	487	ARG
1	D	511	VAL
1	D	515	GLN
1	D	520	ASN
1	D	523	VAL

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Mol	Chain	Res	Type
1	D	525	SER
1	D	544	GLN
1	D	548	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (57) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	157	GLN
1	A	159	GLN
1	A	167	ASN
1	A	185	ASN
1	A	238	ASN
1	A	241	ASN
1	A	259	GLN
1	A	419	ASN
1	A	424	ASN
1	A	451	HIS
1	A	478	GLN
1	A	520	ASN
1	B	93	GLN
1	B	122	ASN
1	B	128	GLN
1	B	157	GLN
1	B	159	GLN
1	B	167	ASN
1	B	185	ASN
1	B	201	HIS
1	B	238	ASN
1	B	241	ASN
1	B	248	GLN
1	B	259	GLN
1	B	344	GLN
1	B	346	ASN
1	B	419	ASN
1	B	424	ASN
1	B	520	ASN
1	C	122	ASN
1	C	157	GLN
1	C	159	GLN
1	C	167	ASN
1	C	174	GLN
1	C	180	HIS

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Mol	Chain	Res	Type
1	C	185	ASN
1	C	238	ASN
1	C	334	ASN
1	C	344	GLN
1	C	346	ASN
1	C	419	ASN
1	C	424	ASN
1	C	474	GLN
1	C	478	GLN
1	C	520	ASN
1	D	157	GLN
1	D	159	GLN
1	D	167	ASN
1	D	180	HIS
1	D	185	ASN
1	D	238	ASN
1	D	241	ASN
1	D	346	ASN
1	D	419	ASN
1	D	424	ASN
1	D	478	GLN
1	D	502	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 ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	465/593 (78%)	-0.26	9 (1%) 66 64	9, 25, 46, 60	0
1	B	465/593 (78%)	-0.21	15 (3%) 50 48	11, 23, 50, 62	0
1	C	468/593 (78%)	-0.33	9 (1%) 66 64	9, 21, 43, 60	0
1	D	471/593 (79%)	-0.34	12 (2%) 58 56	10, 23, 44, 55	0
All	All	1869/2372 (78%)	-0.29	45 (2%) 59 57	9, 23, 47, 62	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	289	ALA	5.0
1	D	545	TRP	4.1
1	A	542	VAL	4.0
1	B	326	ASP	3.5
1	B	542	VAL	3.3
1	A	546	HIS	3.2
1	C	551	VAL	3.2
1	A	545	TRP	3.2
1	D	324	THR	3.1
1	D	325	GLY	3.1
1	C	324	THR	3.0
1	D	322	ALA	3.0
1	D	542	VAL	3.0
1	D	54	ALA	2.9
1	C	540	ALA	2.8
1	D	551	VAL	2.8
1	C	545	TRP	2.8
1	B	288	SER	2.8
1	C	55	SER	2.7
1	D	304	THR	2.7
1	B	266	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	546	HIS	2.6
1	C	322	ALA	2.6
1	C	325	GLY	2.6
1	B	322	ALA	2.6
1	A	323	ASP	2.5
1	B	303	LYS	2.5
1	C	542	VAL	2.5
1	A	325	GLY	2.4
1	B	324	THR	2.4
1	D	541	LYS	2.3
1	A	547	LEU	2.3
1	B	176	VAL	2.2
1	B	545	TRP	2.2
1	D	540	ALA	2.2
1	A	541	LYS	2.2
1	B	325	GLY	2.2
1	C	546	HIS	2.2
1	B	170	TRP	2.1
1	B	304	THR	2.1
1	A	405	GLY	2.1
1	B	546	HIS	2.1
1	A	404	GLY	2.1
1	D	550	TYR	2.1
1	B	321	PRO	2.0

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.