



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:57 PM UTC

PDB ID : 3H8F / pdb_00003h8f
Title : High pH native structure of leucine aminopeptidase from *Pseudomonas putida*
Authors : Kale, A.; Dijkstra, B.W.; Sonke, T.; Thunnissen, A.M.W.H.
Deposited on : 2009-04-29
Resolution : 2.20 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

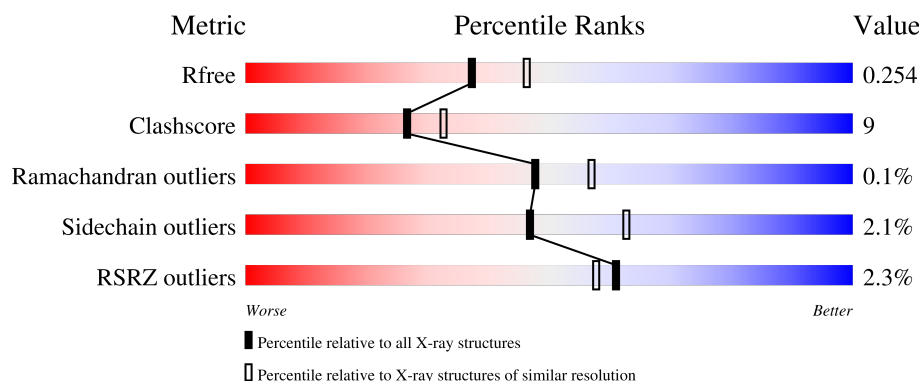
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	497	3% 82% 16% .
1	B	497	2% 83% 16% .
1	C	497	% 84% 15% .
1	D	497	2% 82% 16% .
1	E	497	3% 85% 12% .

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Mol	Chain	Length	Quality of chain
1	F	497	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	BCT	E	504	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 22836 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 Cytosol aminopeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	497	Total	C	N	O	S	0	0	0
			3682	2324	644	696	18			
1	B	497	Total	C	N	O	S	0	0	0
			3682	2324	644	696	18			
1	C	497	Total	C	N	O	S	0	0	0
			3682	2324	644	696	18			
1	D	497	Total	C	N	O	S	0	0	0
			3682	2324	644	696	18			
1	E	497	Total	C	N	O	S	0	0	0
			3682	2324	644	696	18			
1	F	497	Total	C	N	O	S	0	0	0
			3682	2324	644	696	18			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		
2	E	1	Total	Zn	0	0
			1	1		
2	F	1	Total	Zn	0	0
			1	1		

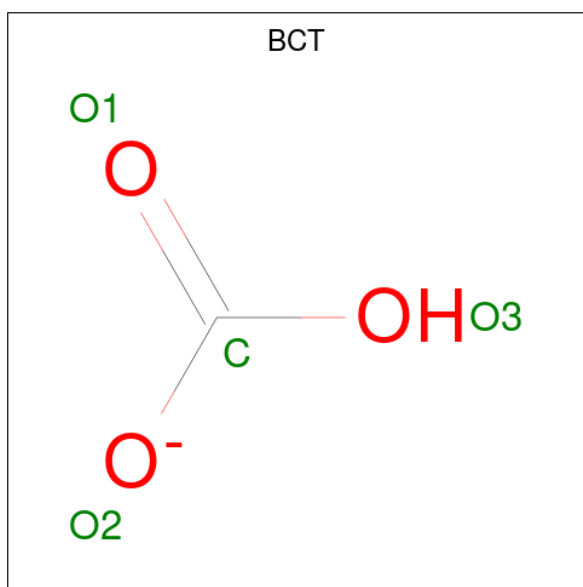
- Molecule 3 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mn 1 1	0	0
3	B	1	Total Mn 1 1	0	0
3	C	1	Total Mn 1 1	0	0
3	D	1	Total Mn 1 1	0	0
3	E	1	Total Mn 1 1	0	0
3	F	1	Total Mn 1 1	0	0

- Molecule 4 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total K 1 1	0	0
4	B	1	Total K 1 1	0	0
4	C	1	Total K 1 1	0	0
4	D	1	Total K 1 1	0	0
4	E	1	Total K 1 1	0	0
4	F	1	Total K 1 1	0	0

- Molecule 5 is BICARBONATE ION (CCD ID: BCT) (formula: CHO_3).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	1	3		
5	B	1	Total	C	O	0	0
			4	1	3		
5	C	1	Total	C	O	0	0
			4	1	3		
5	D	1	Total	C	O	0	0
			4	1	3		
5	E	1	Total	C	O	0	0
			4	1	3		
5	F	1	Total	C	O	0	0
			4	1	3		

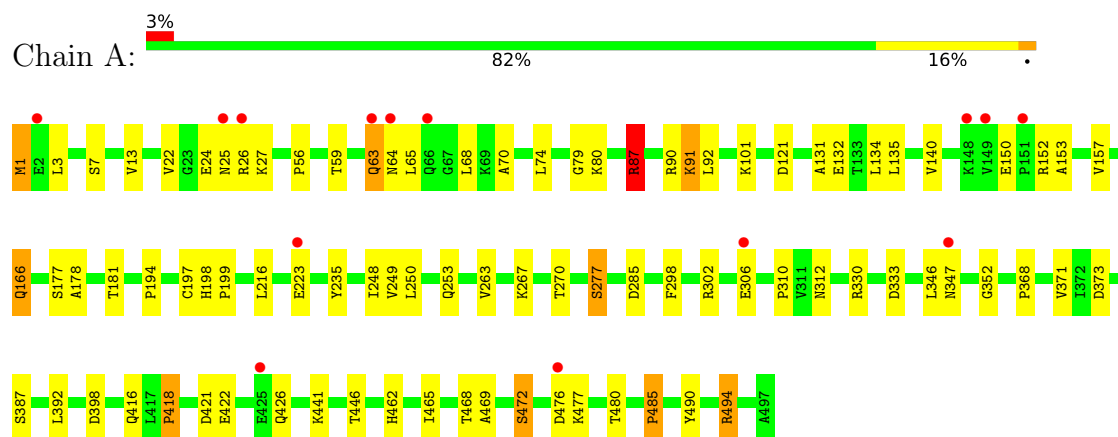
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	114	Total	O	0	0
			114	114		
6	B	99	Total	O	0	0
			99	99		
6	C	162	Total	O	0	0
			162	162		
6	D	111	Total	O	0	0
			111	111		
6	E	123	Total	O	0	0
			123	123		
6	F	93	Total	O	0	0
			93	93		

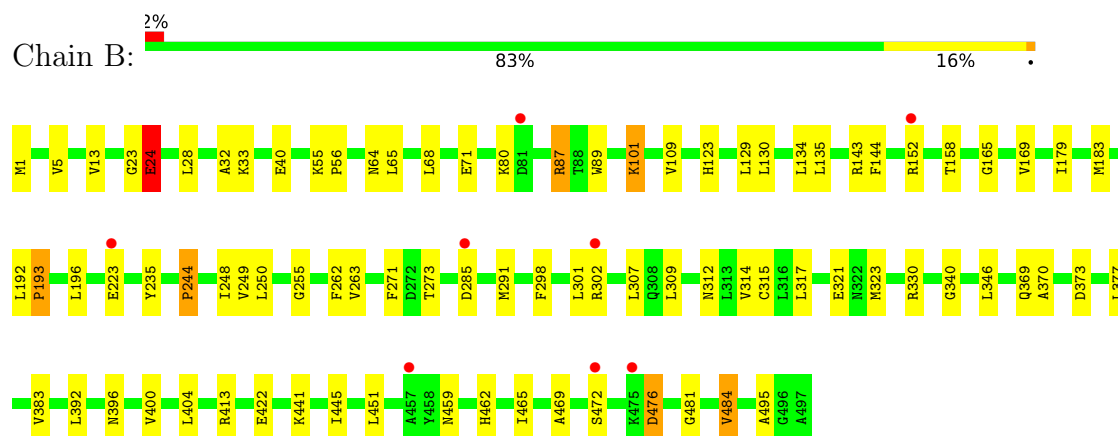
3 Residue-property plots [i](#)

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.

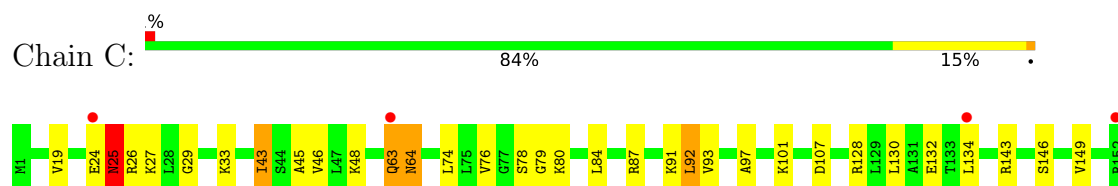
• Molecule 1: Cytosol aminopeptidase

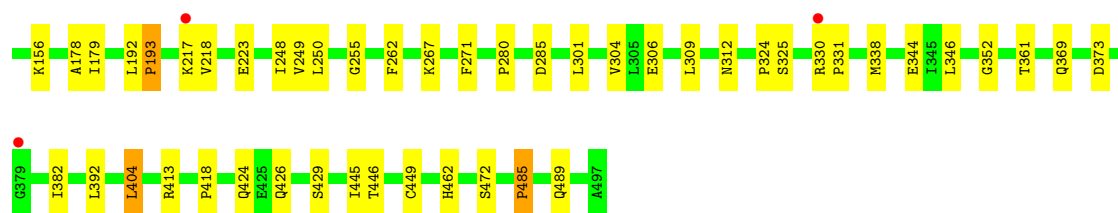


• Molecule 1: Cytosol aminopeptidase

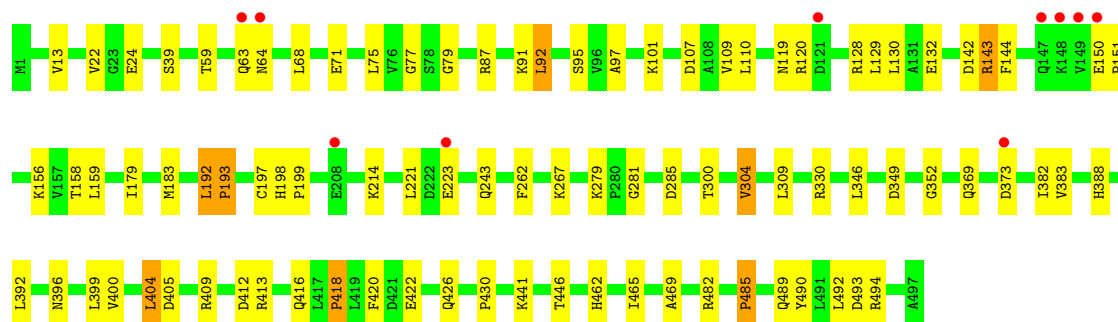
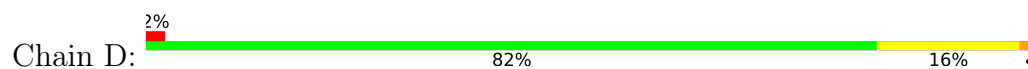


• Molecule 1: Cytosol aminopeptidase

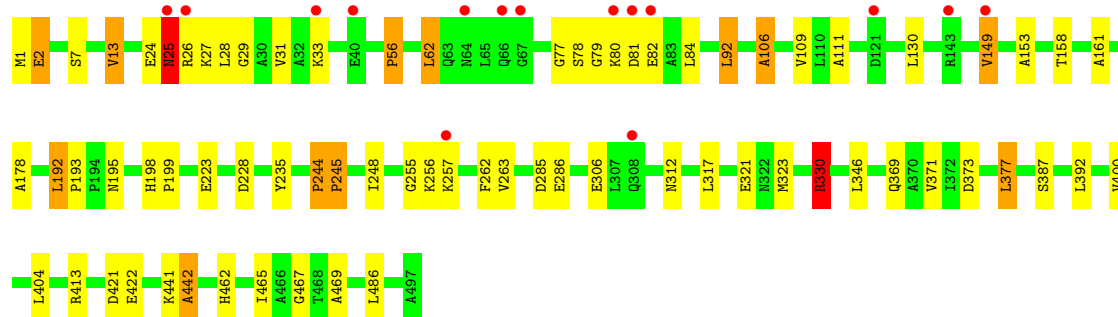
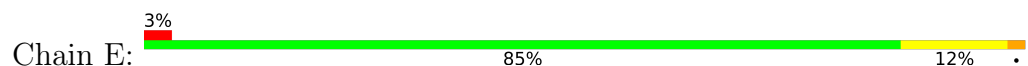




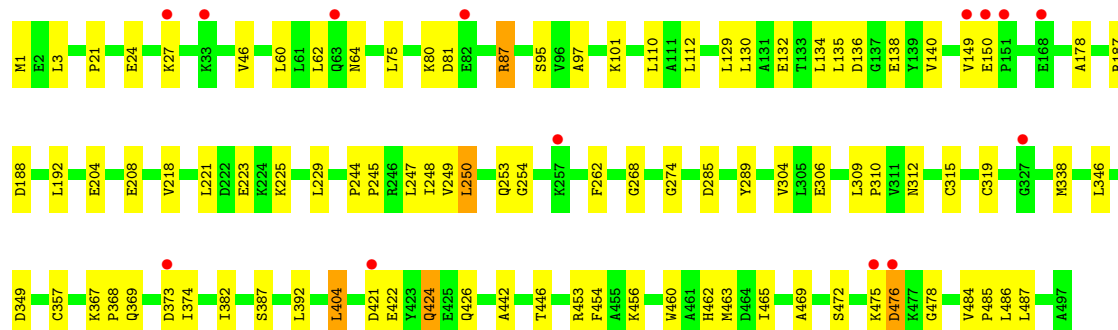
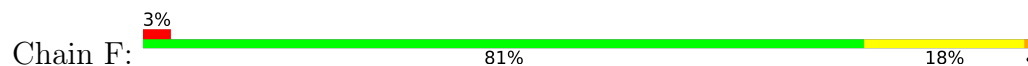
• Molecule 1: Cytosol aminopeptidase



• Molecule 1: Cytosol aminopeptidase



• Molecule 1: Cytosol aminopeptidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	95.79Å 95.93Å 96.31Å 68.43° 76.31° 94.86°	Depositor
Resolution (Å)	47.45 – 2.20 47.45 – 2.20	Depositor EDS
% Data completeness (in resolution range)	93.5 (47.45-2.20) 93.5 (47.45-2.20)	Depositor EDS
R_{merge}	0.37	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.45 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.3.0022	Depositor
R, R_{free}	0.192 , 0.251 0.203 , 0.254	Depositor DCC
R_{free} test set	16590 reflections (9.95%)	wwPDB-VP
Wilson B-factor (Å ²)	21.6	Xtriage
Anisotropy	0.310	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 32.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	22836	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.75% 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

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, BCT, K, MN

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	1.05	11/3737 (0.3%)	1.12	15/5054 (0.3%)
1	B	0.97	7/3737 (0.2%)	1.07	12/5054 (0.2%)
1	C	1.02	4/3737 (0.1%)	1.11	11/5054 (0.2%)
1	D	0.94	5/3737 (0.1%)	1.08	14/5054 (0.3%)
1	E	0.97	6/3737 (0.2%)	1.09	12/5054 (0.2%)
1	F	0.85	1/3737 (0.0%)	1.01	4/5054 (0.1%)
All	All	0.97	34/22422 (0.2%)	1.08	68/30324 (0.2%)

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	C	0	2
1	D	0	1
1	E	0	2
All	All	0	6

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	301	LEU	C-N	12.32	1.48	1.33
1	A	91	LYS	C-N	-8.58	1.23	1.33
1	D	144	PHE	C-O	-7.95	1.13	1.24
1	B	40	GLU	C-N	-7.48	1.23	1.33
1	A	418	PRO	N-CD	-7.26	1.37	1.47
1	D	430	PRO	N-CD	-6.91	1.38	1.47
1	A	121	ASP	CB-CG	-6.44	1.35	1.52
1	C	301	LEU	C-O	-6.42	1.16	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	149	VAL	CA-CB	6.42	1.60	1.54
1	A	87	ARG	C-O	-6.41	1.16	1.24
1	C	485	PRO	N-CD	-6.29	1.39	1.47
1	A	421	ASP	C-O	-6.29	1.16	1.24
1	D	214	LYS	C-O	-6.19	1.16	1.24
1	B	302	ARG	C-N	-6.12	1.26	1.33
1	A	494	ARG	C-O	-6.11	1.17	1.24
1	F	424	GLN	C-O	-6.04	1.16	1.24
1	B	193	PRO	N-CD	-6.02	1.39	1.47
1	B	33	LYS	C-O	-5.98	1.17	1.24
1	A	90	ARG	C-N	-5.98	1.25	1.33
1	A	70	ALA	C-O	-5.93	1.16	1.23
1	B	71	GLU	C-O	-5.91	1.17	1.24
1	C	193	PRO	N-CD	-5.75	1.39	1.47
1	A	131	ALA	C-N	-5.59	1.26	1.33
1	E	330	ARG	C-O	-5.55	1.16	1.23
1	D	214	LYS	CB-CG	-5.52	1.35	1.52
1	E	245	PRO	N-CD	-5.41	1.40	1.47
1	D	418	PRO	N-CD	-5.38	1.40	1.47
1	E	2	GLU	C-O	-5.35	1.17	1.23
1	E	149	VAL	CA-CB	5.34	1.60	1.54
1	A	485	PRO	N-CD	-5.31	1.40	1.47
1	A	310	PRO	N-CD	-5.18	1.40	1.47
1	E	192	LEU	C-O	-5.12	1.18	1.24
1	B	23	GLY	C-O	-5.07	1.17	1.23
1	E	62	LEU	C-O	-5.05	1.17	1.23

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1	MET	CA-C-N	12.22	139.59	122.72
1	A	1	MET	C-N-CA	12.22	139.59	122.72
1	B	24	GLU	N-CA-C	10.13	123.36	111.71
1	E	245	PRO	CA-N-CD	-10.08	97.89	112.00
1	C	193	PRO	CA-N-CD	-9.91	98.12	112.00
1	B	193	PRO	CA-N-CD	-9.45	98.77	112.00
1	E	244	PRO	CA-N-CD	-9.39	98.85	112.00
1	D	418	PRO	CA-N-CD	-9.35	98.92	112.00
1	D	485	PRO	CA-N-CD	-9.35	98.91	112.00
1	A	310	PRO	CA-N-CD	-9.34	98.92	112.00
1	A	418	PRO	CA-N-CD	-9.13	99.22	112.00
1	B	244	PRO	CA-N-CD	-8.97	99.44	112.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	56	PRO	CA-N-CD	-8.88	99.57	112.00
1	C	485	PRO	CA-N-CD	-8.77	99.73	112.00
1	A	485	PRO	CA-N-CD	-8.75	99.75	112.00
1	C	280	PRO	CA-C-N	-8.52	104.71	121.41
1	C	280	PRO	C-N-CA	-8.52	104.71	121.41
1	D	430	PRO	CA-N-CD	-8.38	100.26	112.00
1	B	301	LEU	O-C-N	-8.32	113.50	122.07
1	D	119	ASN	CA-C-N	7.81	133.50	122.72
1	D	119	ASN	C-N-CA	7.81	133.50	122.72
1	B	472	SER	N-CA-C	7.78	119.55	111.14
1	C	192	LEU	CA-C-N	7.68	125.19	119.66
1	C	192	LEU	C-N-CA	7.68	125.19	119.66
1	B	192	LEU	CA-C-N	7.44	125.02	119.66
1	B	192	LEU	C-N-CA	7.44	125.02	119.66
1	E	330	ARG	N-CA-C	6.88	118.47	109.93
1	D	243	GLN	CA-C-N	-6.60	113.58	120.38
1	D	243	GLN	C-N-CA	-6.60	113.58	120.38
1	D	120	ARG	O-C-N	-6.46	115.97	123.27
1	A	472	SER	N-CA-C	6.28	118.13	111.28
1	A	480	THR	N-CA-C	-6.25	101.21	110.46
1	F	484	VAL	N-CA-CB	6.24	114.43	110.50
1	F	472	SER	N-CA-C	5.88	117.76	111.36
1	B	445	ILE	N-CA-C	5.79	115.97	110.42
1	A	480	THR	CA-C-N	-5.76	110.13	121.41
1	A	480	THR	C-N-CA	-5.76	110.13	121.41
1	B	193	PRO	O-C-N	5.73	123.84	121.15
1	D	142	ASP	CA-C-O	5.70	126.32	119.59
1	D	214	LYS	N-CA-C	5.70	118.43	111.82
1	E	13	VAL	N-CA-C	5.66	116.89	109.58
1	C	64	ASN	CB-CA-C	-5.62	100.73	109.61
1	B	64	ASN	CB-CA-C	-5.50	102.01	110.26
1	D	400	VAL	N-CA-CB	5.47	118.78	110.58
1	B	413	ARG	N-CA-C	5.46	118.19	110.50
1	A	421	ASP	N-CA-C	5.40	117.17	111.28
1	F	244	PRO	CA-C-N	5.39	125.37	120.03
1	F	244	PRO	C-N-CA	5.39	125.37	120.03
1	D	193	PRO	O-C-N	5.39	123.68	121.15
1	E	228	ASP	N-CA-C	5.33	117.84	111.71
1	C	429	SER	CA-C-N	-5.33	114.18	119.56
1	C	429	SER	C-N-CA	-5.33	114.18	119.56
1	C	472	SER	N-CA-C	5.27	117.11	111.36
1	B	484	VAL	N-CA-CB	5.24	113.80	110.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	22	VAL	CA-C-N	5.22	126.43	120.53
1	A	22	VAL	C-N-CA	5.22	126.43	120.53
1	E	442	ALA	CA-C-N	5.21	125.88	119.99
1	E	442	ALA	C-N-CA	5.21	125.88	119.99
1	D	281	GLY	N-CA-C	5.21	118.98	112.73
1	C	280	PRO	O-C-N	-5.14	115.35	122.24
1	E	421	ASP	N-CA-C	5.14	117.62	111.71
1	A	166	GLN	N-CA-C	5.13	117.27	111.11
1	E	413	ARG	N-CA-C	5.10	117.69	110.50
1	A	13	VAL	N-CA-C	5.09	115.91	108.53
1	A	480	THR	O-C-N	-5.08	116.95	122.84
1	E	25	ASN	CA-C-N	5.05	129.72	122.40
1	E	25	ASN	C-N-CA	5.05	129.72	122.40
1	D	413	ARG	N-CA-C	5.05	117.95	110.48

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	63	GLN	Peptide
1	C	25	ASN	Peptide
1	C	63	GLN	Peptide
1	D	63	GLN	Peptide
1	E	106	ALA	Peptide
1	E	25	ASN	Peptide

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	3682	0	3777	67	1
1	B	3682	0	3777	84	2
1	C	3682	0	3777	80	1
1	D	3682	0	3777	69	0
1	E	3682	0	3777	74	4
1	F	3682	0	3777	75	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
5	A	4	0	0	0	0
5	B	4	0	0	0	0
5	C	4	0	0	0	0
5	D	4	0	0	0	0
5	E	4	0	0	6	0
5	F	4	0	0	0	0
6	A	114	0	0	7	0
6	B	99	0	0	7	0
6	C	162	0	0	3	0
6	D	111	0	0	6	0
6	E	123	0	0	3	0
6	F	93	0	0	4	0
All	All	22836	0	22662	418	4

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 9.

All (418) 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:F:101:LYS:HE3	1:F:138:GLU:OE2	1.28	1.33
1:F:27:LYS:HE2	1:F:81:ASP:OD1	1.30	1.31
1:C:223:GLU:HB3	6:C:3723:HOH:O	1.48	1.13
1:B:248:ILE:HG22	1:B:250:LEU:HD11	1.31	1.12

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:THR:HG21	1:A:306:GLU:OE2	1.51	1.10
1:D:39:SER:HB3	1:D:68:LEU:HD21	1.34	1.03
1:E:377:LEU:CD1	5:E:504:BCT:C	2.37	1.03
1:B:262:PHE:CZ	1:B:369:GLN:NE2	2.26	1.01
1:F:262:PHE:CZ	1:F:369:GLN:NE2	2.27	1.01
1:D:223:GLU:HB3	6:D:3693:HOH:O	1.61	1.00
1:C:262:PHE:CZ	1:C:369:GLN:NE2	2.29	1.00
1:D:39:SER:CB	1:D:68:LEU:HD21	1.92	0.99
1:F:101:LYS:CE	1:F:138:GLU:OE2	2.10	0.99
1:B:193:PRO:HD2	1:B:193:PRO:O	1.58	0.98
1:F:1:MET:HE3	1:F:135:LEU:HD23	1.44	0.98
1:E:262:PHE:CZ	1:E:369:GLN:NE2	2.34	0.95
1:A:87:ARG:HE	1:A:87:ARG:HA	1.31	0.95
1:B:307:LEU:HB2	1:B:309:LEU:CD2	1.97	0.94
1:B:248:ILE:HG22	1:B:250:LEU:CD1	1.96	0.94
1:E:27:LYS:NZ	1:E:81:ASP:OD1	2.01	0.93
1:F:178:ALA:HA	1:F:306:GLU:OE1	1.68	0.93
1:F:188:ASP:O	1:F:192:LEU:HD23	1.66	0.93
1:B:373:ASP:OD1	1:B:462:HIS:HA	1.70	0.92
1:F:188:ASP:O	1:F:192:LEU:CD2	2.17	0.92
1:E:377:LEU:HD12	5:E:504:BCT:O2	1.71	0.91
1:C:130:LEU:HD21	1:C:134:LEU:HD11	1.54	0.89
1:D:192:LEU:HG	1:D:197:CYS:HB2	1.53	0.89
1:A:181:THR:CG2	1:A:306:GLU:OE2	2.22	0.88
1:E:223:GLU:HB3	6:E:3720:HOH:O	1.75	0.87
1:F:248:ILE:HG22	1:F:250:LEU:HD11	1.57	0.87
1:B:248:ILE:CG2	1:B:250:LEU:HD11	2.05	0.86
1:F:1:MET:CE	1:F:135:LEU:HD23	2.07	0.85
1:B:307:LEU:CB	1:B:309:LEU:CD2	2.53	0.85
1:F:248:ILE:HG22	1:F:250:LEU:CD1	2.05	0.85
1:D:392:LEU:HD23	1:D:404:LEU:HD13	1.57	0.85
1:B:1:MET:HE1	1:B:134:LEU:HG	1.59	0.83
1:A:418:PRO:HD2	1:A:418:PRO:O	1.76	0.83
1:C:25:ASN:OD1	1:C:27:LYS:HD2	1.78	0.83
1:B:249:VAL:C	1:B:250:LEU:HD12	2.04	0.83
1:C:330:ARG:NH2	1:E:195:ASN:HA	1.94	0.82
1:A:178:ALA:HA	1:A:306:GLU:OE1	1.80	0.81
1:A:223:GLU:HB3	6:A:3688:HOH:O	1.80	0.81
1:C:193:PRO:HD2	1:C:193:PRO:O	1.81	0.80
1:D:373:ASP:OD1	1:D:462:HIS:HA	1.81	0.80
1:C:128:ARG:HH12	1:C:489:GLN:HE22	1.29	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:373:ASP:OD1	1:E:462:HIS:HA	1.80	0.80
1:E:377:LEU:CD1	5:E:504:BCT:O2	2.29	0.80
1:E:377:LEU:HD11	5:E:504:BCT:C	2.11	0.80
1:D:418:PRO:HD2	1:D:418:PRO:O	1.80	0.80
1:B:285:ASP:OD2	1:E:346:LEU:HD22	1.83	0.79
1:D:392:LEU:HD23	1:D:404:LEU:CD1	2.12	0.78
1:C:248:ILE:HG22	1:C:250:LEU:CD1	2.13	0.77
1:C:424:GLN:NE2	1:C:449:CYS:SG	2.58	0.77
1:D:300:THR:O	1:D:304:VAL:HG12	1.84	0.77
1:D:39:SER:CB	1:D:68:LEU:CD2	2.61	0.77
1:C:304:VAL:HG12	1:C:309:LEU:HD22	1.67	0.77
1:A:248:ILE:HG22	1:A:250:LEU:HD11	1.66	0.76
1:A:285:ASP:OD2	1:D:346:LEU:HD22	1.84	0.76
1:F:27:LYS:HE2	1:F:81:ASP:CG	2.10	0.76
1:B:250:LEU:HD12	1:B:250:LEU:N	2.00	0.76
1:F:1:MET:HE1	1:F:134:LEU:HG	1.68	0.76
1:E:392:LEU:CD2	1:E:404:LEU:HD22	2.15	0.76
1:B:451:LEU:CD1	1:B:462:HIS:CD2	2.68	0.75
1:E:245:PRO:HD2	1:E:245:PRO:O	1.87	0.75
1:B:223:GLU:HB3	6:B:3698:HOH:O	1.86	0.75
1:E:56:PRO:HD3	1:E:79:GLY:CA	2.17	0.75
1:B:244:PRO:HD2	1:B:244:PRO:O	1.86	0.74
1:A:373:ASP:OD1	1:A:462:HIS:HA	1.87	0.74
1:C:25:ASN:HB2	1:C:27:LYS:HB2	1.70	0.74
1:D:262:PHE:CZ	1:D:369:GLN:NE2	2.55	0.74
1:A:249:VAL:C	1:A:250:LEU:HD12	2.13	0.74
1:C:130:LEU:CD2	1:C:134:LEU:HD11	2.18	0.73
1:A:248:ILE:HG22	1:A:250:LEU:CD1	2.18	0.73
1:F:392:LEU:HD23	1:F:404:LEU:CD1	2.18	0.73
1:B:373:ASP:OD2	1:B:462:HIS:HD2	1.72	0.72
1:B:451:LEU:CD1	1:B:462:HIS:NE2	2.53	0.72
1:B:24:GLU:OE1	1:B:80:LYS:HG3	1.88	0.72
1:C:373:ASP:OD1	1:C:462:HIS:HA	1.89	0.72
1:E:377:LEU:HD12	1:E:377:LEU:O	1.89	0.72
1:C:267:LYS:HE3	1:C:352:GLY:HA3	1.71	0.72
1:D:267:LYS:HE3	1:D:352:GLY:HA3	1.72	0.71
1:D:405:ASP:OD2	1:D:409:ARG:NH1	2.24	0.71
1:F:249:VAL:C	1:F:250:LEU:HD12	2.15	0.71
1:B:1:MET:HE3	1:B:135:LEU:HD23	1.72	0.71
1:B:400:VAL:O	1:B:404:LEU:HD23	1.89	0.71
1:B:392:LEU:HD23	1:B:404:LEU:CD2	2.21	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:392:LEU:CD2	1:B:404:LEU:HD22	2.21	0.70
1:D:426:GLN:HE21	1:D:441:LYS:HE3	1.56	0.69
1:B:87:ARG:HD2	1:F:422:GLU:OE2	1.92	0.69
1:B:373:ASP:OD2	1:B:462:HIS:CD2	2.46	0.69
1:D:39:SER:HB3	1:D:68:LEU:CD2	2.19	0.69
1:E:109:VAL:HG22	1:E:158:THR:HB	1.73	0.69
1:C:217:LYS:HD2	1:C:218:VAL:N	2.07	0.69
1:B:451:LEU:HD12	1:B:462:HIS:CD2	2.27	0.69
1:F:373:ASP:OD1	1:F:462:HIS:HA	1.92	0.69
1:D:39:SER:HA	1:D:68:LEU:CD2	2.24	0.68
1:E:244:PRO:HD2	1:E:244:PRO:O	1.92	0.68
1:E:377:LEU:HD12	5:E:504:BCT:C	2.18	0.68
1:D:39:SER:CA	1:D:68:LEU:HD21	2.22	0.68
1:F:223:GLU:HB2	6:F:3716:HOH:O	1.92	0.68
1:D:267:LYS:HE3	1:D:352:GLY:CA	2.24	0.68
1:A:422:GLU:OE1	1:C:413:ARG:NH2	2.27	0.67
1:B:451:LEU:HD12	1:B:462:HIS:NE2	2.09	0.67
1:E:392:LEU:HD23	1:E:404:LEU:CD2	2.24	0.67
1:A:426:GLN:OE1	1:A:446:THR:HG21	1.94	0.67
1:B:307:LEU:CB	1:B:309:LEU:HD22	2.22	0.67
1:B:309:LEU:HD12	1:B:495:ALA:CB	2.24	0.67
1:D:285:ASP:OD2	1:F:346:LEU:HD22	1.95	0.67
1:C:25:ASN:O	1:C:26:ARG:HB2	1.94	0.67
1:F:188:ASP:O	1:F:192:LEU:HD22	1.95	0.67
1:A:87:ARG:HD3	1:A:91:LYS:HE3	1.77	0.66
1:C:267:LYS:HE3	1:C:352:GLY:CA	2.25	0.66
1:C:248:ILE:HG22	1:C:250:LEU:HD11	1.77	0.66
1:B:307:LEU:HB3	1:B:309:LEU:CD2	2.25	0.65
1:F:1:MET:HE3	1:F:135:LEU:CD2	2.24	0.65
1:B:383:VAL:O	1:F:442:ALA:HA	1.97	0.65
1:B:193:PRO:O	1:B:193:PRO:CD	2.38	0.65
1:B:451:LEU:HD13	1:B:462:HIS:CD2	2.31	0.65
1:B:392:LEU:HD23	1:B:404:LEU:HD21	1.78	0.65
1:F:456:LYS:NZ	6:F:3247:HOH:O	2.21	0.65
1:D:128:ARG:HH12	1:D:489:GLN:HE22	1.45	0.65
1:B:250:LEU:CD1	1:B:250:LEU:N	2.59	0.65
1:D:192:LEU:HD12	1:D:193:PRO:HD2	1.78	0.65
1:C:330:ARG:HD3	1:E:321:GLU:OE2	1.97	0.64
1:C:338:MET:HE3	1:C:361:THR:HG22	1.79	0.64
1:A:472:SER:O	1:A:477:LYS:HD2	1.97	0.64
1:B:346:LEU:CD2	1:C:285:ASP:OD2	2.45	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:307:LEU:HB3	1:B:309:LEU:HD22	1.79	0.64
1:A:346:LEU:HD22	1:F:285:ASP:OD2	1.98	0.64
1:B:1:MET:CE	1:B:135:LEU:HD23	2.28	0.64
1:F:248:ILE:CG2	1:F:250:LEU:HD11	2.27	0.64
1:E:400:VAL:O	1:E:404:LEU:HD23	1.98	0.63
1:C:255:GLY:H	1:C:312:ASN:ND2	1.97	0.63
1:C:392:LEU:HD23	1:C:404:LEU:CD1	2.28	0.63
1:E:24:GLU:C	1:E:26:ARG:H	2.05	0.63
1:B:244:PRO:HD2	6:B:3694:HOH:O	1.99	0.62
1:C:178:ALA:HA	1:C:306:GLU:OE1	1.98	0.62
1:E:244:PRO:HD2	6:E:516:HOH:O	1.99	0.62
1:F:3:LEU:HD13	1:F:135:LEU:HD21	1.82	0.62
1:C:330:ARG:HH21	1:E:195:ASN:HA	1.65	0.62
1:A:426:GLN:OE1	1:A:446:THR:CG2	2.48	0.61
1:B:271:PHE:CD2	1:E:330:ARG:HG2	2.34	0.61
1:A:267:LYS:HE3	1:A:352:GLY:HA3	1.81	0.61
1:A:285:ASP:OD2	1:D:346:LEU:CD2	2.46	0.61
1:B:285:ASP:OD2	1:E:346:LEU:CD2	2.48	0.61
1:E:29:GLY:O	1:E:33:LYS:HG3	2.01	0.61
1:A:101:LYS:HD2	1:A:152:ARG:NE	2.15	0.61
1:C:97:ALA:O	1:C:101:LYS:HG3	2.00	0.61
1:C:130:LEU:CD2	1:C:134:LEU:CD1	2.77	0.61
1:D:39:SER:CA	1:D:68:LEU:CD2	2.78	0.61
1:B:307:LEU:HB2	1:B:309:LEU:HD23	1.81	0.60
1:B:307:LEU:CB	1:B:309:LEU:HD21	2.31	0.60
1:D:426:GLN:NE2	1:D:441:LYS:HE3	2.17	0.60
1:C:43:ILE:O	1:C:46:VAL:HG22	2.01	0.60
1:E:56:PRO:HD3	1:E:79:GLY:N	2.16	0.60
1:E:25:ASN:HB3	1:E:27:LYS:HB2	1.84	0.60
1:C:76:VAL:HG22	1:C:92:LEU:CD2	2.32	0.59
1:D:409:ARG:NH2	1:D:493:ASP:OD2	2.36	0.59
1:E:387:SER:HA	1:E:467:GLY:O	2.02	0.59
1:C:130:LEU:O	1:C:134:LEU:HG	2.02	0.59
1:A:418:PRO:O	1:A:418:PRO:CD	2.51	0.59
1:B:392:LEU:CD2	1:B:404:LEU:CD2	2.79	0.59
1:C:223:GLU:CB	6:C:3723:HOH:O	2.26	0.59
1:B:370:ALA:HB2	6:B:3777:HOH:O	2.02	0.59
1:D:426:GLN:OE1	1:D:446:THR:HG21	2.03	0.58
1:C:130:LEU:HD21	1:C:134:LEU:CD1	2.32	0.58
1:C:128:ARG:HH12	1:C:489:GLN:NE2	1.99	0.58
1:E:392:LEU:HD23	1:E:404:LEU:HD22	1.82	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:441:LYS:HG2	6:D:3713:HOH:O	2.03	0.58
1:F:253:GLN:HA	1:F:312:ASN:HD22	1.69	0.58
1:D:223:GLU:CB	6:D:3693:HOH:O	2.36	0.56
1:D:22:VAL:O	1:D:77:GLY:HA2	2.05	0.56
1:E:56:PRO:HD3	1:E:79:GLY:HA3	1.88	0.56
1:F:424:GLN:HE21	1:F:453:ARG:HH21	1.53	0.56
1:A:64:ASN:OD1	1:A:64:ASN:O	2.24	0.56
1:F:426:GLN:OE1	1:F:446:THR:HG21	2.06	0.56
1:A:330:ARG:NH1	1:A:333:ASP:OD1	2.39	0.55
1:B:101:LYS:HD2	1:B:152:ARG:HH12	1.69	0.55
1:C:217:LYS:HD2	1:C:218:VAL:H	1.70	0.55
1:E:255:GLY:H	1:E:312:ASN:ND2	2.05	0.55
1:A:3:LEU:CD2	1:A:157:VAL:HB	2.37	0.55
1:A:277:SER:CB	1:A:347:ASN:OD1	2.55	0.55
1:C:424:GLN:NE2	1:C:449:CYS:CB	2.70	0.55
1:A:346:LEU:CD2	1:F:285:ASP:OD2	2.54	0.55
1:C:424:GLN:HE22	1:C:449:CYS:CB	2.20	0.55
1:B:250:LEU:HB2	1:B:315:CYS:HB2	1.89	0.55
1:E:462:HIS:HE1	6:E:550:HOH:O	1.89	0.55
1:A:223:GLU:HG2	1:A:235:TYR:HE1	1.72	0.55
1:E:56:PRO:CD	1:E:79:GLY:CA	2.83	0.54
1:B:101:LYS:HG2	1:B:152:ARG:NH1	2.22	0.54
1:F:465:ILE:HD12	1:F:469:ALA:HB2	1.88	0.54
1:A:25:ASN:O	1:A:26:ARG:HB2	2.06	0.54
1:B:263:VAL:HG22	1:B:314:VAL:HB	1.88	0.54
1:C:248:ILE:CG2	1:C:250:LEU:HD11	2.36	0.54
1:E:285:ASP:OD1	1:E:286:GLU:N	2.41	0.54
1:C:76:VAL:CG2	1:C:92:LEU:CD2	2.85	0.54
1:D:87:ARG:HG3	1:D:91:LYS:HE2	1.90	0.54
1:B:144:PHE:HZ	1:B:196:LEU:HD12	1.73	0.54
1:D:279:LYS:HE2	1:D:349:ASP:OD1	2.07	0.54
1:C:249:VAL:C	1:C:250:LEU:HD12	2.33	0.53
1:E:223:GLU:OE1	1:E:244:PRO:HB3	2.08	0.53
1:A:248:ILE:CG2	1:A:250:LEU:HD11	2.34	0.53
1:C:330:ARG:HH21	1:E:195:ASN:CA	2.20	0.53
1:D:418:PRO:O	1:D:418:PRO:CD	2.55	0.53
1:A:1:MET:HE2	1:A:135:LEU:HD23	1.91	0.53
1:B:89:TRP:CZ2	1:B:130:LEU:HD12	2.43	0.53
1:F:387:SER:HB3	6:F:865:HOH:O	2.08	0.53
1:A:267:LYS:HE3	1:A:352:GLY:CA	2.39	0.53
1:E:377:LEU:HD11	5:E:504:BCT:O1	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:132:GLU:OE2	1:F:485:PRO:HG3	2.09	0.52
1:E:28:LEU:O	1:E:33:LYS:HD2	2.08	0.52
1:D:426:GLN:OE1	1:D:446:THR:CG2	2.57	0.52
1:D:465:ILE:HD12	1:D:469:ALA:HB2	1.92	0.52
1:E:223:GLU:HB2	1:E:245:PRO:HD2	1.92	0.52
1:C:193:PRO:O	1:C:193:PRO:CD	2.56	0.52
1:C:330:ARG:O	1:C:331:PRO:C	2.52	0.52
1:C:130:LEU:HD23	1:C:134:LEU:CD1	2.40	0.52
1:F:475:LYS:HG3	1:F:476:ASP:OD1	2.09	0.52
1:C:426:GLN:OE1	1:C:446:THR:HG21	2.10	0.51
1:E:77:GLY:O	1:E:92:LEU:HD22	2.09	0.51
1:E:130:LEU:C	1:E:130:LEU:HD23	2.35	0.51
1:A:63:GLN:OE1	1:E:62:LEU:CD2	2.58	0.51
1:C:392:LEU:HD23	1:C:404:LEU:HD13	1.90	0.51
1:F:229:LEU:HB3	1:F:338:MET:HE3	1.92	0.51
1:F:486:LEU:HD23	1:F:486:LEU:C	2.36	0.51
1:E:7:SER:HA	1:E:161:ALA:O	2.11	0.51
1:B:346:LEU:HD23	1:C:285:ASP:OD2	2.10	0.51
1:B:400:VAL:HG12	1:B:404:LEU:HD23	1.93	0.51
1:E:392:LEU:CD2	1:E:404:LEU:CD2	2.83	0.51
1:B:101:LYS:HD2	1:B:152:ARG:NH1	2.26	0.51
1:F:223:GLU:HG3	1:F:245:PRO:HD2	1.92	0.51
1:C:218:VAL:HG22	1:C:250:LEU:HG	1.91	0.51
1:B:244:PRO:CD	6:B:3694:HOH:O	2.56	0.50
1:A:59:THR:HB	1:A:74:LEU:HD11	1.93	0.50
1:D:198:HIS:HB2	1:D:199:PRO:CD	2.41	0.50
1:F:188:ASP:C	1:F:192:LEU:HD23	2.35	0.50
1:D:109:VAL:HG22	1:D:158:THR:HB	1.93	0.50
1:E:223:GLU:HB2	1:E:245:PRO:CD	2.42	0.49
1:F:250:LEU:HB2	1:F:315:CYS:HB2	1.93	0.49
1:A:462:HIS:HE1	6:A:538:HOH:O	1.95	0.49
1:C:24:GLU:HG3	1:C:79:GLY:C	2.36	0.49
1:D:382:ILE:HG22	1:E:441:LYS:HG2	1.92	0.49
1:A:87:ARG:HA	1:A:87:ARG:NE	2.13	0.49
1:D:396:ASN:HD22	1:D:399:LEU:H	1.60	0.49
1:F:130:LEU:C	1:F:130:LEU:HD23	2.37	0.49
1:A:7:SER:HB2	1:A:166:GLN:OE1	2.13	0.49
1:C:76:VAL:HG21	1:C:92:LEU:HD22	1.95	0.49
1:E:245:PRO:O	1:E:245:PRO:CD	2.59	0.49
1:F:424:GLN:NE2	1:F:453:ARG:HE	2.10	0.49
1:D:132:GLU:OE2	1:D:485:PRO:HG3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:27:LYS:CE	1:F:81:ASP:OD1	2.26	0.49
1:C:130:LEU:HD23	1:C:134:LEU:HD12	1.95	0.49
1:E:178:ALA:HA	1:E:306:GLU:OE1	2.12	0.49
1:A:223:GLU:CB	6:A:3688:HOH:O	2.52	0.48
1:B:255:GLY:H	1:B:312:ASN:ND2	2.11	0.48
1:B:255:GLY:HA2	6:B:3748:HOH:O	2.13	0.48
1:C:179:ILE:HG23	1:C:485:PRO:HD3	1.95	0.48
1:F:97:ALA:HB1	1:F:101:LYS:NZ	2.28	0.48
1:F:426:GLN:OE1	1:F:446:THR:CG2	2.60	0.48
1:D:373:ASP:OD1	1:D:462:HIS:CA	2.58	0.48
1:F:21:PRO:HD3	1:F:112:LEU:HD23	1.94	0.48
1:D:143:ARG:HB3	6:D:3730:HOH:O	2.13	0.48
1:E:244:PRO:O	1:E:244:PRO:CD	2.62	0.48
1:A:101:LYS:HD2	1:A:152:ARG:HE	1.78	0.48
1:A:177:SER:HB3	6:A:760:HOH:O	2.12	0.48
1:F:149:VAL:HG12	1:F:150:GLU:O	2.13	0.48
1:C:424:GLN:NE2	1:C:449:CYS:HB3	2.28	0.47
1:D:392:LEU:CD2	1:D:404:LEU:HD13	2.36	0.47
1:A:132:GLU:OE2	1:A:485:PRO:HG3	2.14	0.47
1:A:250:LEU:HD12	1:A:250:LEU:N	2.29	0.47
1:C:330:ARG:HH22	1:E:195:ASN:HA	1.74	0.47
1:E:373:ASP:OD1	1:E:462:HIS:CA	2.58	0.47
1:E:392:LEU:HD23	1:E:404:LEU:HD21	1.93	0.47
1:E:56:PRO:CD	1:E:79:GLY:HA3	2.44	0.47
1:B:321:GLU:HG2	1:B:323:MET:HG2	1.95	0.47
1:A:263:VAL:HG23	1:A:368:PRO:HB3	1.96	0.47
1:B:1:MET:CE	1:B:134:LEU:HG	2.36	0.47
1:E:321:GLU:HG2	1:E:323:MET:HG2	1.96	0.47
1:A:277:SER:HB3	1:A:347:ASN:OD1	2.15	0.47
1:F:250:LEU:HD12	1:F:250:LEU:N	2.28	0.47
1:A:24:GLU:O	1:A:25:ASN:CB	2.62	0.47
1:F:218:VAL:HG22	1:F:250:LEU:HG	1.97	0.47
1:F:374:ILE:HD12	1:F:374:ILE:N	2.30	0.46
1:C:324:PRO:O	1:C:325:SER:HB3	2.15	0.46
1:A:253:GLN:HA	1:A:312:ASN:HD22	1.80	0.46
1:D:396:ASN:ND2	1:D:399:LEU:H	2.14	0.46
1:F:262:PHE:HZ	1:F:369:GLN:NE2	2.06	0.46
1:C:45:ALA:HA	1:C:48:LYS:HD2	1.97	0.46
1:E:486:LEU:C	1:E:486:LEU:HD23	2.39	0.46
1:D:97:ALA:O	1:D:101:LYS:HG3	2.16	0.46
1:B:55:LYS:O	1:B:56:PRO:C	2.59	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:392:LEU:HD13	1:F:463:MET:HG2	1.97	0.46
1:A:24:GLU:O	1:A:25:ASN:HB2	2.16	0.46
1:D:179:ILE:CG2	1:D:485:PRO:HD3	2.45	0.46
1:F:136:ASP:OD1	1:F:187:ARG:NH1	2.48	0.46
1:B:262:PHE:CE2	1:B:369:GLN:NE2	2.81	0.46
1:D:392:LEU:O	1:D:416:GLN:HA	2.16	0.46
1:D:462:HIS:HE1	6:D:528:HOH:O	1.98	0.45
1:E:106:ALA:HA	1:E:153:ALA:O	2.17	0.45
1:E:198:HIS:HB2	1:E:199:PRO:CD	2.47	0.45
1:E:465:ILE:HD12	1:E:469:ALA:HB2	1.98	0.45
1:F:374:ILE:HG12	1:F:487:LEU:HD11	1.98	0.45
1:C:78:SER:O	1:C:84:LEU:HD21	2.16	0.45
1:D:130:LEU:C	1:D:130:LEU:HD23	2.41	0.45
1:F:262:PHE:CE2	1:F:369:GLN:NE2	2.82	0.45
1:D:412:ASP:OD1	1:D:482:ARG:NE	2.47	0.45
1:C:130:LEU:HD23	1:C:130:LEU:C	2.42	0.45
1:F:460:TRP:CZ2	1:F:462:HIS:HB2	2.52	0.45
1:E:248:ILE:HB	1:E:317:LEU:HB2	1.99	0.44
1:D:77:GLY:O	1:D:92:LEU:HD22	2.18	0.44
1:D:128:ARG:HH22	1:D:489:GLN:NE2	2.15	0.44
1:A:490:TYR:O	1:A:494:ARG:HG2	2.16	0.44
1:B:5:VAL:HG21	1:B:169:VAL:HG12	1.98	0.44
1:B:441:LYS:HG2	1:F:382:ILE:HG22	1.99	0.44
1:E:78:SER:O	1:E:84:LEU:HD21	2.17	0.44
1:A:465:ILE:HD12	1:A:469:ALA:HB2	1.99	0.44
1:A:302:ARG:HG3	6:A:3785:HOH:O	2.16	0.44
1:C:250:LEU:HD12	1:C:250:LEU:N	2.32	0.44
1:C:338:MET:HE3	1:C:361:THR:CG2	2.44	0.44
1:D:24:GLU:HG3	1:D:79:GLY:C	2.43	0.44
1:F:250:LEU:CD1	1:F:250:LEU:N	2.81	0.44
1:A:65:LEU:HB2	1:A:68:LEU:HD12	1.99	0.44
1:B:330:ARG:HG2	1:C:271:PHE:CD2	2.53	0.44
1:F:221:LEU:HD22	1:F:247:LEU:HD23	1.99	0.44
1:B:271:PHE:HE2	1:B:273:THR:HG22	1.83	0.44
1:B:404:LEU:HD12	6:B:535:HOH:O	2.17	0.44
1:E:223:GLU:HG2	1:E:235:TYR:HE1	1.83	0.44
1:F:24:GLU:OE1	1:F:80:LYS:HG3	2.17	0.44
1:E:80:LYS:O	1:E:82:GLU:N	2.51	0.44
1:B:465:ILE:HD12	1:B:469:ALA:HB2	1.98	0.43
1:D:383:VAL:O	1:E:442:ALA:HA	2.19	0.43
1:A:3:LEU:HD23	1:A:157:VAL:HB	1.98	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:GLN:O	1:C:64:ASN:O	2.36	0.43
1:C:91:LYS:HG3	6:C:3783:HOH:O	2.17	0.43
1:E:263:VAL:HB	1:E:371:VAL:HG22	2.00	0.43
1:B:476:ASP:N	1:B:476:ASP:OD1	2.51	0.43
1:E:257:LYS:HD2	1:E:257:LYS:HA	1.66	0.43
1:A:1:MET:HE1	1:A:134:LEU:HG	1.99	0.43
1:A:216:LEU:HD11	1:A:250:LEU:HB3	1.99	0.43
1:B:143:ARG:HB2	6:B:513:HOH:O	2.19	0.43
1:C:25:ASN:OD1	1:C:27:LYS:HB2	2.17	0.43
1:B:109:VAL:HG22	1:B:158:THR:HB	2.00	0.43
1:B:441:LYS:HG2	1:F:382:ILE:CG2	2.49	0.43
1:A:198:HIS:HB2	1:A:199:PRO:CD	2.48	0.43
1:F:268:GLY:O	1:F:319:CYS:HA	2.19	0.43
1:F:367:LYS:N	1:F:368:PRO:HD3	2.34	0.43
1:A:153:ALA:HA	6:A:3779:HOH:O	2.19	0.43
1:A:392:LEU:O	1:A:416:GLN:HA	2.18	0.43
1:C:63:GLN:O	1:C:64:ASN:C	2.62	0.43
1:E:192:LEU:HD23	1:E:193:PRO:HD2	2.00	0.43
1:C:29:GLY:O	1:C:33:LYS:HG3	2.18	0.43
1:C:330:ARG:HH21	1:E:195:ASN:N	2.16	0.43
1:D:64:ASN:HB3	1:D:71:GLU:O	2.18	0.43
1:A:1:MET:HE2	1:A:1:MET:HB3	1.78	0.42
1:A:387:SER:O	1:A:468:THR:HG22	2.19	0.42
1:B:223:GLU:HG2	1:B:235:TYR:HE1	1.84	0.42
1:D:128:ARG:HH12	1:D:489:GLN:NE2	2.14	0.42
1:F:204:GLU:O	1:F:208:GLU:HG3	2.19	0.42
1:A:253:GLN:HA	1:A:312:ASN:ND2	2.35	0.42
1:B:346:LEU:HD22	1:C:285:ASP:OD2	2.17	0.42
1:F:392:LEU:HD23	1:F:404:LEU:HD12	2.00	0.42
1:C:392:LEU:CD2	1:C:404:LEU:HD13	2.48	0.42
1:B:144:PHE:CZ	1:B:196:LEU:HD12	2.52	0.42
1:B:396:ASN:HD22	1:B:459:ASN:HB3	1.84	0.42
1:C:426:GLN:OE1	1:C:446:THR:CG2	2.67	0.42
1:C:132:GLU:OE2	1:C:485:PRO:HG3	2.19	0.42
1:D:490:TYR:O	1:D:494:ARG:HG2	2.20	0.42
1:F:274:GLY:HA2	1:F:349:ASP:OD1	2.20	0.42
1:F:424:GLN:HE22	1:F:453:ARG:HE	1.67	0.42
1:A:250:LEU:CD1	1:A:250:LEU:N	2.83	0.42
1:A:87:ARG:HE	1:A:87:ARG:CA	1.96	0.42
1:B:179:ILE:O	1:B:183:MET:HG3	2.20	0.42
1:D:183:MET:HE1	6:D:1219:HOH:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:330:ARG:NH2	1:E:195:ASN:CA	2.71	0.42
1:C:392:LEU:CD2	1:C:404:LEU:CD1	2.96	0.42
1:D:110:LEU:HB2	1:D:159:LEU:HD23	2.02	0.42
1:E:29:GLY:O	1:E:33:LYS:CG	2.68	0.42
1:A:197:CYS:O	1:A:270:THR:HB	2.19	0.41
1:F:225:LYS:NZ	1:F:229:LEU:HD21	2.35	0.41
1:F:254:GLY:O	1:F:310:PRO:HA	2.19	0.41
1:A:441:LYS:HG2	1:C:382:ILE:HG23	2.02	0.41
1:D:409:ARG:NH2	1:D:493:ASP:CG	2.78	0.41
1:F:60:LEU:HD22	1:F:75:LEU:HD12	2.02	0.41
1:D:150:GLU:HA	1:D:151:PRO:HD3	1.87	0.41
1:F:46:VAL:HG11	1:F:62:LEU:HD21	2.03	0.41
1:B:481:GLY:O	1:B:484:VAL:HG23	2.20	0.41
1:D:420:PHE:HB3	1:D:422:GLU:OE1	2.20	0.41
1:A:263:VAL:HB	1:A:371:VAL:HG22	2.02	0.41
1:E:262:PHE:HZ	1:E:369:GLN:NE2	2.10	0.41
1:B:248:ILE:HB	1:B:317:LEU:HB2	2.03	0.41
1:B:330:ARG:HH11	1:B:330:ARG:HD2	1.63	0.41
1:B:377:LEU:HD22	1:B:462:HIS:CE1	2.56	0.41
1:B:422:GLU:OE2	1:F:87:ARG:NH1	2.49	0.41
1:F:304:VAL:HG13	1:F:309:LEU:HB2	2.01	0.41
1:B:28:LEU:HD22	1:B:32:ALA:HB1	2.02	0.41
1:C:19:VAL:HG22	1:C:74:LEU:HB3	2.02	0.41
1:F:46:VAL:HG23	6:F:1725:HOH:O	2.21	0.41
1:E:31:VAL:HG11	1:E:111:ALA:HB1	2.03	0.41
1:A:56:PRO:HD3	1:A:79:GLY:HA3	2.03	0.41
1:C:93:VAL:HG13	1:C:134:LEU:HD21	2.03	0.41
1:D:128:ARG:NH1	1:D:489:GLN:HE22	2.16	0.41
1:D:388:HIS:CE1	1:E:422:GLU:HB3	2.56	0.41
1:C:25:ASN:CB	1:C:27:LYS:HB2	2.47	0.41
1:C:107:ASP:HA	1:C:156:LYS:O	2.21	0.41
1:D:198:HIS:HB2	1:D:199:PRO:HD2	2.01	0.41
1:F:289:TYR:OH	1:F:478:GLY:HA2	2.20	0.41
1:A:476:ASP:HB2	6:A:3746:HOH:O	2.20	0.40
1:D:107:ASP:HA	1:D:156:LYS:O	2.22	0.40
1:D:309:LEU:HD21	1:D:492:LEU:HD23	2.03	0.40
1:A:194:PRO:HB2	1:D:330:ARG:HD2	2.03	0.40
1:D:59:THR:HA	1:D:75:LEU:O	2.21	0.40
1:E:28:LEU:HB2	1:E:33:LYS:HG2	2.02	0.40
1:F:357:CYS:HB2	1:F:454:PHE:CD1	2.57	0.40
1:B:65:LEU:HB2	1:B:68:LEU:HD12	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:400:VAL:CG1	1:B:404:LEU:HD23	2.51	0.40
1:C:418:PRO:HD2	1:C:445:ILE:HD11	2.04	0.40
1:B:123:HIS:CE1	1:B:165:GLY:HA3	2.56	0.40
1:C:344:GLU:HG2	1:C:346:LEU:HD12	2.03	0.40

All (4) 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:A:398:ASP:OD2	1:E:256:LYS:CE[1_455]	1.66	0.54
1:C:146:SER:O	1:E:26:ARG:NH2[1_554]	1.77	0.43
1:B:340:GLY:CA	1:E:25:ASN:ND2[1_554]	2.12	0.08
1:B:340:GLY:C	1:E:25:ASN:ND2[1_554]	2.18	0.02

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	495/497 (100%)	478 (97%)	17 (3%)	0	100	100
1	B	495/497 (100%)	475 (96%)	20 (4%)	0	100	100
1	C	495/497 (100%)	478 (97%)	16 (3%)	1 (0%)	43	51
1	D	495/497 (100%)	478 (97%)	17 (3%)	0	100	100
1	E	495/497 (100%)	475 (96%)	20 (4%)	0	100	100
1	F	495/497 (100%)	479 (97%)	15 (3%)	1 (0%)	43	51
All	All	2970/2982 (100%)	2863 (96%)	105 (4%)	2 (0%)	48	57

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	64	ASN

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Mol	Chain	Res	Type
1	C	25	ASN

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	378/378 (100%)	370 (98%)	8 (2%)	47	63
1	B	378/378 (100%)	370 (98%)	8 (2%)	47	63
1	C	378/378 (100%)	372 (98%)	6 (2%)	55	71
1	D	378/378 (100%)	369 (98%)	9 (2%)	43	58
1	E	378/378 (100%)	371 (98%)	7 (2%)	50	66
1	F	378/378 (100%)	369 (98%)	9 (2%)	43	58
All	All	2268/2268 (100%)	2221 (98%)	47 (2%)	47	63

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LYS
1	A	80	LYS
1	A	87	ARG
1	A	92	LEU
1	A	140	VAL
1	A	150	GLU
1	A	277	SER
1	A	298	PHE
1	B	13	VAL
1	B	24	GLU
1	B	87	ARG
1	B	101	LYS
1	B	129	LEU
1	B	291	MET
1	B	298	PHE
1	B	476	ASP
1	C	43	ILE

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Mol	Chain	Res	Type
1	C	80	LYS
1	C	87	ARG
1	C	92	LEU
1	C	143	ARG
1	C	404	LEU
1	D	13	VAL
1	D	92	LEU
1	D	95	SER
1	D	129	LEU
1	D	143	ARG
1	D	192	LEU
1	D	221	LEU
1	D	304	VAL
1	D	404	LEU
1	E	1	MET
1	E	2	GLU
1	E	13	VAL
1	E	92	LEU
1	E	149	VAL
1	E	330	ARG
1	E	377	LEU
1	F	87	ARG
1	F	95	SER
1	F	110	LEU
1	F	129	LEU
1	F	140	VAL
1	F	250	LEU
1	F	404	LEU
1	F	421	ASP
1	F	476	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	205	GLN
1	A	213	HIS
1	A	253	GLN
1	A	312	ASN
1	A	459	ASN
1	A	462	HIS
1	B	312	ASN
1	B	462	HIS

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Mol	Chain	Res	Type
1	C	119	ASN
1	C	123	HIS
1	C	312	ASN
1	C	424	GLN
1	C	459	ASN
1	C	489	GLN
1	D	205	GLN
1	D	312	ASN
1	D	341	GLN
1	D	388	HIS
1	D	396	ASN
1	D	426	GLN
1	D	462	HIS
1	D	489	GLN
1	E	118	ASN
1	E	119	ASN
1	E	123	HIS
1	E	308	GLN
1	E	312	ASN
1	E	462	HIS
1	F	104	ASN
1	F	312	ASN
1	F	424	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 24 ligands modelled in this entry, 18 are monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	BCT	B	504	-	3,3,3	0.64	0	2,3,3	0.47	0
5	BCT	D	504	-	3,3,3	0.49	0	2,3,3	0.74	0
5	BCT	C	504	-	3,3,3	0.58	0	2,3,3	0.69	0
5	BCT	F	504	-	3,3,3	0.78	0	2,3,3	1.01	0
5	BCT	A	504	-	3,3,3	0.66	0	2,3,3	0.69	0
5	BCT	E	504	-	3,3,3	0.65	0	2,3,3	0.51	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	504	BCT	6	0

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	497/497 (100%)	-0.05	14 (2%) 55 52	2, 11, 27, 41	0
1	B	497/497 (100%)	0.05	8 (1%) 70 67	3, 15, 28, 41	0
1	C	497/497 (100%)	-0.26	7 (1%) 73 71	2, 9, 23, 35	0
1	D	497/497 (100%)	-0.01	10 (2%) 65 62	4, 13, 28, 51	0
1	E	497/497 (100%)	-0.02	15 (3%) 52 49	2, 12, 30, 43	0
1	F	497/497 (100%)	0.10	14 (2%) 55 52	3, 13, 26, 38	0
All	All	2982/2982 (100%)	-0.03	68 (2%) 61 58	2, 12, 27, 51	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	25	ASN	7.4
1	D	63	GLN	6.0
1	A	306	GLU	5.8
1	A	63	GLN	5.3
1	F	63	GLN	5.2
1	C	63	GLN	4.8
1	D	149	VAL	4.8
1	F	149	VAL	3.9
1	D	148	LYS	3.8
1	E	149	VAL	3.8
1	E	81	ASP	3.7
1	A	149	VAL	3.7
1	B	475	LYS	3.6
1	A	26	ARG	3.5
1	D	147	GLN	3.5
1	E	121	ASP	3.4
1	A	425	GLU	3.4
1	D	64	ASN	3.2
1	F	82	GLU	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	81	ASP	3.0
1	E	33	LYS	3.0
1	A	64	ASN	3.0
1	E	67	GLY	3.0
1	E	143	ARG	2.8
1	B	223	GLU	2.8
1	A	148	LYS	2.7
1	E	80	LYS	2.7
1	A	66	GLN	2.6
1	E	26	ARG	2.5
1	E	66	GLN	2.5
1	A	151	PRO	2.4
1	C	152	ARG	2.4
1	E	308	GLN	2.4
1	F	150	GLU	2.4
1	B	285	ASP	2.3
1	E	257	LYS	2.3
1	F	168	GLU	2.3
1	C	330	ARG	2.3
1	F	27	LYS	2.3
1	F	476	ASP	2.2
1	D	373	ASP	2.2
1	B	472	SER	2.2
1	A	476	ASP	2.2
1	D	121	ASP	2.2
1	E	40	GLU	2.2
1	A	347	ASN	2.2
1	F	475	LYS	2.2
1	B	457	ALA	2.2
1	B	302	ARG	2.1
1	D	208	GLU	2.1
1	A	25	ASN	2.1
1	A	223	GLU	2.1
1	D	223	GLU	2.1
1	E	64	ASN	2.1
1	F	151	PRO	2.1
1	A	2	GLU	2.1
1	F	373	ASP	2.1
1	C	379	GLY	2.1
1	D	150	GLU	2.0
1	E	82	GLU	2.0
1	F	327	GLY	2.0

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Mol	Chain	Res	Type	RSRZ
1	F	33	LYS	2.0
1	F	257	LYS	2.0
1	C	24	GLU	2.0
1	C	134	LEU	2.0
1	F	421	ASP	2.0
1	B	152	ARG	2.0
1	C	217	LYS	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	BCT	F	504	4/4	0.94	0.10	11,11,12,13	0
3	MN	C	502	1/1	0.95	0.13	34,34,34,34	0
4	K	F	503	1/1	0.96	0.08	21,21,21,21	0
3	MN	E	502	1/1	0.96	0.14	33,33,33,33	0
5	BCT	A	504	4/4	0.97	0.07	10,11,12,14	0
5	BCT	E	504	4/4	0.97	0.09	7,7,8,9	0
3	MN	B	502	1/1	0.97	0.14	34,34,34,34	0
5	BCT	B	504	4/4	0.98	0.03	12,12,13,13	0
5	BCT	C	504	4/4	0.98	0.05	7,9,9,9	0
5	BCT	D	504	4/4	0.98	0.07	11,12,12,13	0
3	MN	D	502	1/1	0.98	0.10	35,35,35,35	0
3	MN	A	502	1/1	0.98	0.14	37,37,37,37	0
4	K	A	503	1/1	0.99	0.12	9,9,9,9	0
4	K	B	503	1/1	0.99	0.06	13,13,13,13	0
4	K	D	503	1/1	0.99	0.09	10,10,10,10	0
4	K	E	503	1/1	0.99	0.05	12,12,12,12	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	F	501	1/1	0.99	0.08	20,20,20,20	0
2	ZN	A	501	1/1	0.99	0.09	20,20,20,20	0
2	ZN	B	501	1/1	0.99	0.07	21,21,21,21	0
2	ZN	C	501	1/1	0.99	0.06	17,17,17,17	0
2	ZN	D	501	1/1	0.99	0.08	21,21,21,21	0
2	ZN	E	501	1/1	0.99	0.06	18,18,18,18	0
3	MN	F	502	1/1	0.99	0.12	37,37,37,37	0
4	K	C	503	1/1	1.00	0.03	12,12,12,12	0

6.5 Other polymers [i](#)

There are no such residues in this entry.