



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:56 AM UTC

PDB ID : 2J5I / pdb_00002j5i
Title : Crystal Structure of Hydroxycinnamoyl-CoA Hydratase-Lyase
Authors : Leonard, P.M.; Brzozowski, A.M.; Lebedev, A.; Marshall, C.M.; Smith, D.J.;
Verma, C.S.; Walton, N.J.; Grogan, G.
Deposited on : 2006-09-18
Resolution : 1.80 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:

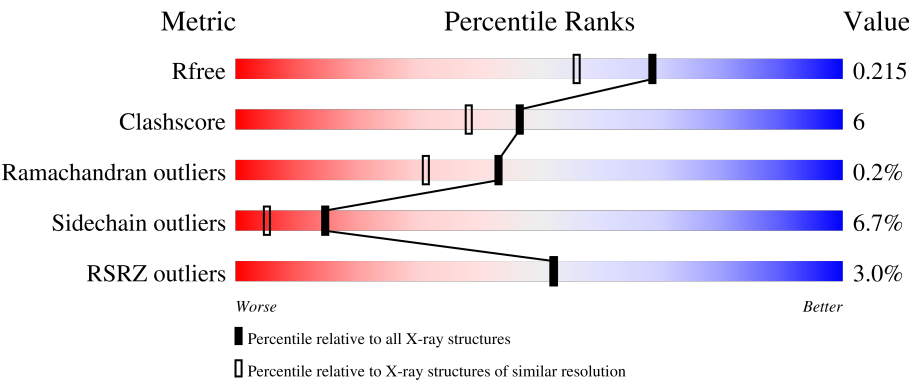
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	276	2% 79%9%10%
2	B	276	3% 75%12%11%
2	C	276	% 79%9%11%
2	D	276	4% 75%13%10%
2	E	276	4% 78%12%10%

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Mol	Chain	Length	Quality of chain
2	F	276	2% 75% 13% 11%
2	G	276	% 78% 9% • 11%
2	H	276	2% 78% 9% • 11%
2	J	276	3% 71% 17% • 11%
2	K	276	4% 76% 12% •• 11%
2	L	276	4% 73% 13% •• 11%
3	I	276	% 72% 15% • 11%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 25264 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 P-HYDROXYCINNAMOYL COA HYDRATASE/LYASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	248	Total	C	N	O	S	25	2	1
			1958	1242	339	363	14			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	169	MET	TYR	conflict	UNP O69762

- Molecule 2 is a protein called P-HYDROXYCINNAMOYL COA HYDRATASE/LYASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	245	Total	C	N	O	S	64	4	1
			1934	1231	334	356	13			
2	C	246	Total	C	N	O	S	46	2	0
			1959	1245	340	362	12			
2	D	248	Total	C	N	O	S	46	3	1
			1968	1252	340	364	12			
2	E	248	Total	C	N	O	S	79	2	1
			1959	1245	338	363	13			
2	F	246	Total	C	N	O	S	45	2	1
			1946	1239	338	357	12			
2	G	246	Total	C	N	O	S	19	3	0
			1960	1247	341	359	13			
2	H	247	Total	C	N	O	S	42	1	1
			1946	1237	336	360	13			
2	J	247	Total	C	N	O	S	82	0	1
			1941	1233	336	360	12			
2	K	247	Total	C	N	O	S	78	2	1
			1952	1242	339	359	12			
2	L	246	Total	C	N	O	S	87	1	1
			1936	1231	335	358	12			

- Molecule 3 is a protein called P-HYDROXYCINNAMOYL COA HYDRATASE/LYASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	246	Total 1944	C 1235	N 340	O 357	S 12	34	1	1

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	232	GLN	GLU	conflict	UNP O69762

- Molecule 4 is water.

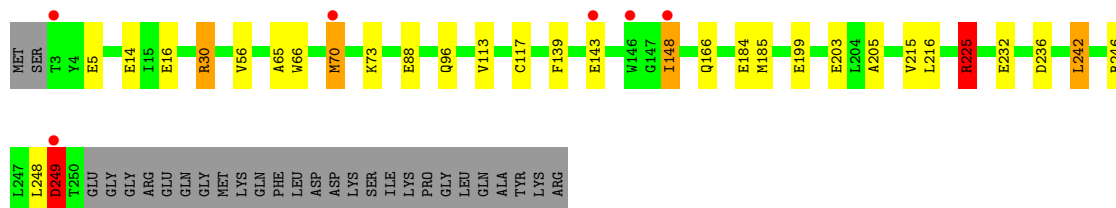
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	163	Total 163	O 163	0	0
4	B	115	Total 115	O 115	0	0
4	C	149	Total 149	O 149	0	0
4	D	149	Total 149	O 149	0	0
4	E	169	Total 169	O 169	0	0
4	F	175	Total 175	O 175	0	0
4	G	198	Total 198	O 198	0	0
4	H	152	Total 152	O 152	0	0
4	I	148	Total 148	O 148	0	0
4	J	146	Total 146	O 146	0	0
4	K	151	Total 151	O 151	0	0
4	L	146	Total 146	O 146	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

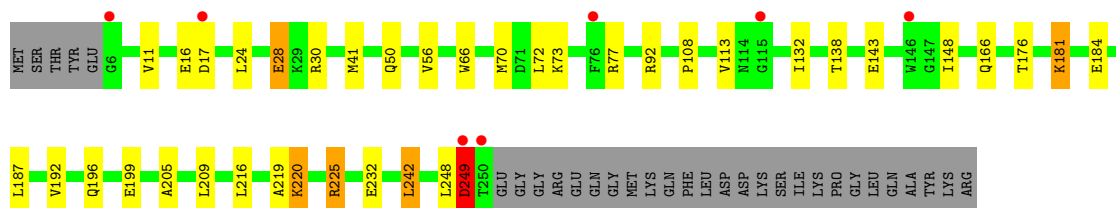
• Molecule 1: P-HYDROXYCINNAMOYL COA HYDRATASE/LYASE

Chain A: 



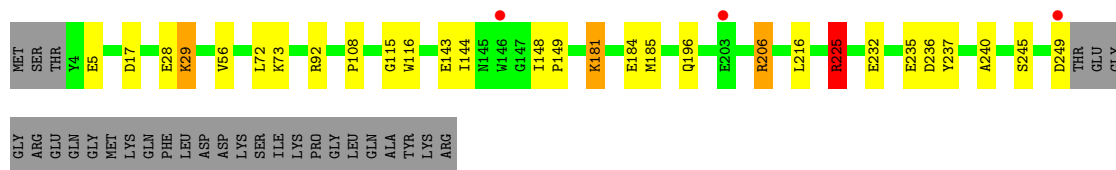
• Molecule 2: P-HYDROXYCINNAMOYL COA HYDRATASE/LYASE

Chain B: 



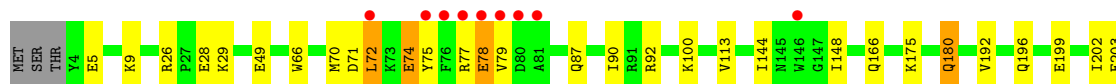
• Molecule 2: P-HYDROXYCINNAMOYL COA HYDRATASE/LYASE

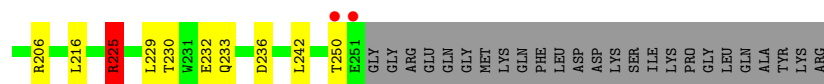
Chain C: 



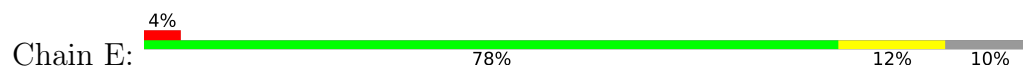
• Molecule 2: P-HYDROXYCINNAMOYL COA HYDRATASE/LYASE

Chain D: 

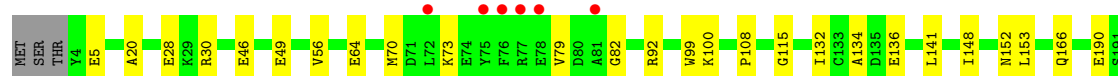
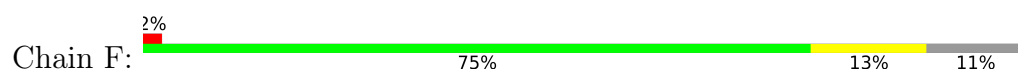




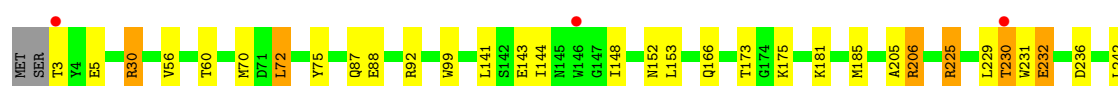
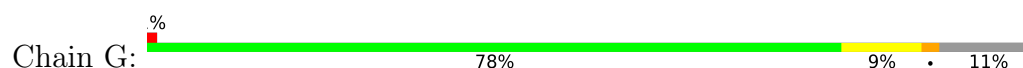
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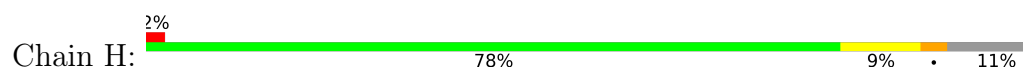
• Molecule 2: P-HYDROXYCINNAMOYL COA HYDRATASE/LYASE



• Molecule 2: P-HYDROXYCINNAMOYL COA HYDRATASE/LYASE

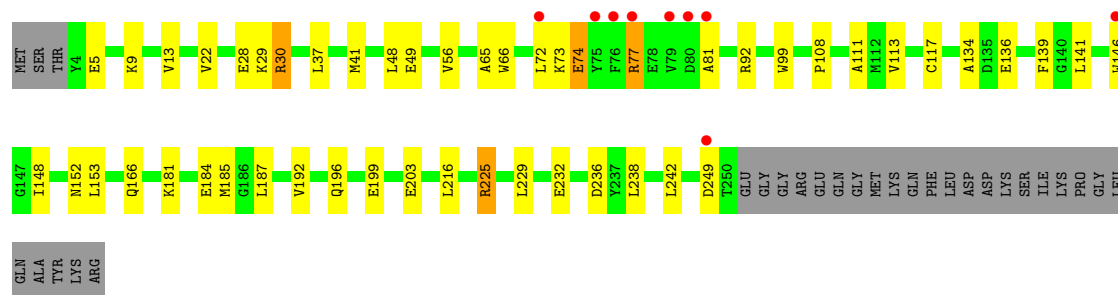


• Molecule 2: P-HYDROXYCINNAMOYL COA HYDRATASE/LYASE

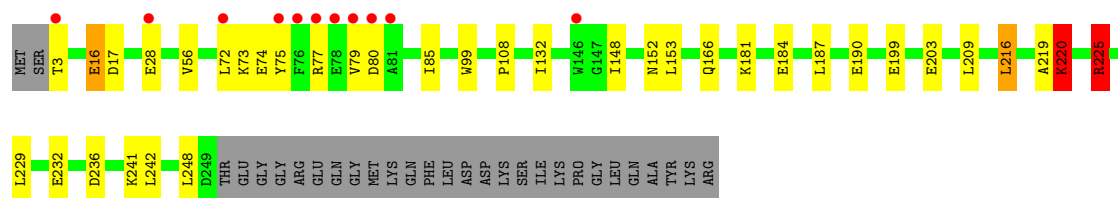
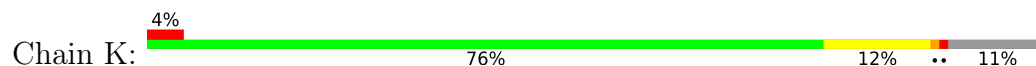


• Molecule 2: P-HYDROXYCINNAMOYL COA HYDRATASE/LYASE

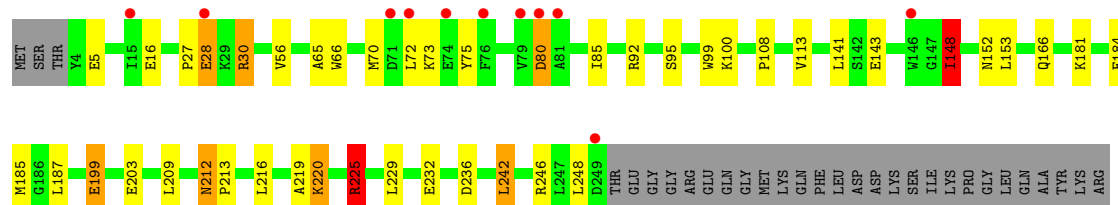




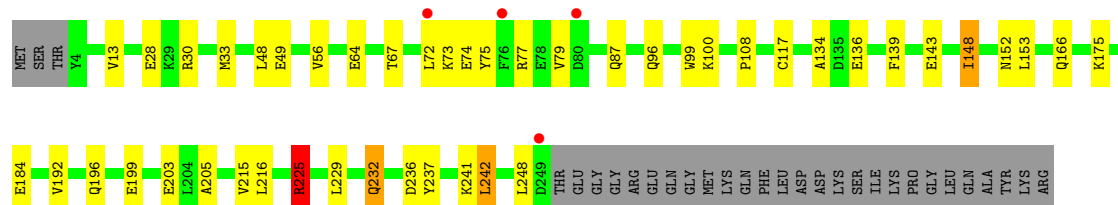
• Molecule 2: P-HYDROXYCINNAMOYL COA HYDRATASE/LYASE



• Molecule 2: P-HYDROXYCINNAMOYL COA HYDRATASE/LYASE



• Molecule 3: P-HYDROXYCINNAMOYL COA HYDRATASE/LYASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	154.24Å 167.49Å 130.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	111.80 – 1.80 111.80 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.5 (111.80-1.80) 99.5 (111.80-1.80)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.09 (at 1.80Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.179 , 0.215 0.180 , 0.215	Depositor DCC
R_{free} test set	15557 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	20.1	Xtrriage
Anisotropy	0.621	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	25264	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 24.10 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.0906e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.99	1/1999 (0.1%)	0.98	3/2704 (0.1%)
2	B	0.98	1/1984 (0.1%)	1.03	8/2685 (0.3%)
2	C	0.94	1/1998 (0.1%)	0.96	2/2703 (0.1%)
2	D	1.01	2/2017 (0.1%)	1.00	2/2730 (0.1%)
2	E	1.03	2/2004 (0.1%)	0.97	0/2712
2	F	1.07	1/1991 (0.1%)	0.99	4/2694 (0.1%)
2	G	1.04	1/2008 (0.0%)	1.00	5/2715 (0.2%)
2	H	1.01	1/1988 (0.1%)	1.02	5/2690 (0.2%)
2	J	0.95	2/1980 (0.1%)	0.97	1/2680 (0.0%)
2	K	1.05	3/1997 (0.2%)	1.07	8/2703 (0.3%)
2	L	0.99	3/1978 (0.2%)	1.04	9/2677 (0.3%)
3	I	1.07	4/1983 (0.2%)	1.00	1/2683 (0.0%)
All	All	1.01	22/23927 (0.1%)	1.00	48/32376 (0.1%)

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
2	E	0	1
2	G	1	0
All	All	1	1

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	K	80	ASP	CA-CB	-13.26	1.32	1.53
2	H	249	ASP	C-N	-7.83	1.22	1.33
2	L	75	TYR	CA-CB	-7.16	1.41	1.53
2	J	249	ASP	C-N	-7.08	1.23	1.33
2	D	250	THR	C-N	-6.98	1.23	1.33
2	K	79	VAL	CB-CG1	-6.97	1.29	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	249	ASP	C-N	-6.91	1.23	1.33
2	B	249	ASP	C-N	-6.88	1.23	1.33
2	G	206	ARG	CD-NE	-6.80	1.36	1.46
2	L	80	ASP	CA-CB	-6.46	1.43	1.53
2	F	20	ALA	CA-CB	5.86	1.60	1.53
3	I	73	LYS	C-O	-5.80	1.17	1.24
2	J	111	ALA	CA-CB	5.53	1.61	1.53
3	I	184	GLU	C-O	-5.53	1.17	1.24
2	L	248	LEU	C-N	-5.48	1.25	1.33
2	E	123	SER	N-CA	5.44	1.50	1.46
3	I	248	LEU	C-N	-5.37	1.25	1.33
2	D	192	VAL	C-O	-5.29	1.20	1.24
2	K	248	LEU	C-N	-5.25	1.26	1.33
3	I	64	GLU	C-O	-5.20	1.17	1.24
2	C	240	ALA	CA-CB	5.10	1.61	1.53
2	E	65	ALA	CA-CB	-5.02	1.45	1.53

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	80	ASP	N-CA-CB	14.36	131.24	110.12
2	K	79	VAL	CA-CB-CG2	-8.01	96.78	110.40
2	K	225	ARG	NE-CZ-NH2	7.50	125.95	119.20
2	D	225	ARG	NE-CZ-NH2	6.92	125.43	119.20
2	K	219	ALA	CA-C-N	6.92	129.78	120.65
2	K	219	ALA	C-N-CA	6.92	129.78	120.65
2	K	225	ARG	NE-CZ-NH1	-6.80	114.69	121.50
2	C	225	ARG	NE-CZ-NH2	6.78	125.30	119.20
2	L	219	ALA	CA-C-N	6.62	129.39	120.65
2	L	219	ALA	C-N-CA	6.62	129.39	120.65
2	G	225	ARG	NE-CZ-NH2	6.48	125.03	119.20
2	H	148	ILE	CA-C-N	-6.29	113.90	120.38
2	H	148	ILE	C-N-CA	-6.29	113.90	120.38
2	G	225	ARG	NE-CZ-NH1	-6.25	115.25	121.50
2	G	206	ARG	CD-NE-CZ	-6.18	115.74	124.40
2	L	225	ARG	NE-CZ-NH2	6.17	124.76	119.20
2	F	225	ARG	NE-CZ-NH2	6.14	124.73	119.20
2	H	225	ARG	NE-CZ-NH2	6.11	124.70	119.20
2	B	225	ARG	NE-CZ-NH2	6.09	124.68	119.20
2	G	60	THR	CA-C-N	-6.01	118.03	122.18
2	G	60	THR	C-N-CA	-6.01	118.03	122.18
2	B	220	LYS	CB-CA-C	-6.01	101.71	110.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	82	GLY	CA-C-N	6.00	125.97	120.03
2	F	82	GLY	C-N-CA	6.00	125.97	120.03
2	B	192	VAL	CA-C-N	-5.84	113.77	119.78
2	B	192	VAL	C-N-CA	-5.84	113.77	119.78
2	H	225	ARG	NE-CZ-NH1	-5.80	115.70	121.50
1	A	225	ARG	NE-CZ-NH2	5.74	124.36	119.20
2	L	225	ARG	NE-CZ-NH1	-5.72	115.78	121.50
2	B	219	ALA	CA-C-N	5.60	128.04	120.65
2	B	219	ALA	C-N-CA	5.60	128.04	120.65
2	B	225	ARG	NE-CZ-NH1	-5.59	115.91	121.50
2	L	212	ASN	CA-C-N	5.54	125.20	119.05
2	L	212	ASN	C-N-CA	5.54	125.20	119.05
3	I	225	ARG	NE-CZ-NH2	5.51	124.16	119.20
2	F	225	ARG	NE-CZ-NH1	-5.40	116.10	121.50
2	L	220	LYS	CB-CA-C	-5.38	102.68	110.96
2	K	80	ASP	CA-CB-CG	5.30	117.91	112.60
2	J	185	MET	N-CA-C	5.23	118.92	112.54
2	C	235	GLU	N-CA-C	-5.22	105.59	111.28
2	L	148	ILE	CA-C-N	-5.12	115.10	120.38
2	L	148	ILE	C-N-CA	-5.12	115.10	120.38
2	D	79	VAL	N-CA-C	5.08	115.51	110.23
2	K	220	LYS	CB-CA-C	-5.07	103.15	110.96
2	B	132	ILE	N-CA-C	-5.05	100.85	108.12
2	H	92	ARG	NE-CZ-NH2	-5.04	114.66	119.20
1	A	148	ILE	CA-C-N	-5.01	115.22	120.38
1	A	148	ILE	C-N-CA	-5.01	115.22	120.38

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	G	230	THR	CB

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	E	248	LEU	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	1958	0	1962	20	0
2	B	1934	0	1952	41	0
2	C	1959	0	1961	18	1
2	D	1968	0	1973	28	0
2	E	1959	0	1963	18	0
2	F	1946	0	1959	22	0
2	G	1960	0	1977	25	0
2	H	1946	0	1948	25	0
2	J	1941	0	1939	25	0
2	K	1952	0	1964	20	1
2	L	1936	0	1940	29	0
3	I	1944	0	1949	27	0
4	A	163	0	0	4	0
4	B	115	0	0	8	0
4	C	149	0	0	2	0
4	D	149	0	0	5	0
4	E	169	0	0	3	0
4	F	175	0	0	5	0
4	G	198	0	0	5	0
4	H	152	0	0	6	0
4	I	148	0	0	2	0
4	J	146	0	0	2	0
4	K	151	0	0	5	0
4	L	146	0	0	5	0
All	All	25264	0	23487	261	1

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 6.

All (261) 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)
2:B:11:VAL:CG2	2:B:41:MET:HE2	1.64	1.27
2:E:71:ASP:CG	2:E:74:GLU:HG3	1.71	1.15
2:B:56[B]:VAL:HG12	2:B:108:PRO:HG2	1.31	1.10
2:B:24:LEU:HD23	2:B:41:MET:HE3	1.30	1.10
2:B:11:VAL:HG21	2:B:41:MET:HE2	1.29	1.07
2:B:24:LEU:CD2	2:B:41:MET:HE3	1.86	1.05
2:E:71:ASP:OD2	2:E:74:GLU:HG3	1.62	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:56[B]:VAL:HG12	2:K:108:PRO:HG2	1.42	0.98
2:B:11:VAL:HG22	2:B:41:MET:HE2	1.45	0.98
2:L:100:LYS:HE3	4:L:2063:HOH:O	1.65	0.96
3:I:33:MET:HE3	3:I:67:THR:HG21	1.49	0.94
3:I:232:GLN:HG2	2:K:232:GLU:HG2	1.47	0.93
1:A:232:GLU:HG2	2:D:232:GLU:HG2	1.52	0.92
2:H:4:TYR:N	4:H:2003:HOH:O	2.04	0.89
2:B:232:GLU:HG2	2:F:232:GLU:HG2	1.54	0.89
2:F:199:GLU:HG3	4:F:2135:HOH:O	1.73	0.87
2:D:71:ASP:CG	2:D:74:GLU:HG3	2.00	0.86
2:C:184:GLU:HG2	4:C:2106:HOH:O	1.75	0.85
2:C:17:ASP:O	2:C:206:ARG:NH1	2.09	0.85
2:K:203:GLU:HG2	4:K:2130:HOH:O	1.76	0.85
2:G:173:THR:OG1	2:G:175:LYS:HG2	1.77	0.85
3:I:56:VAL:HG12	3:I:108:PRO:HG2	1.60	0.83
2:C:56:VAL:HG22	2:C:108:PRO:HG2	1.60	0.83
2:B:138:THR:CG2	2:B:176[A]:THR:CG2	2.58	0.80
2:H:92:ARG:NH2	2:J:49:GLU:OE1	2.13	0.80
2:B:138:THR:HG23	2:B:176[A]:THR:CG2	2.13	0.78
2:B:11:VAL:CG2	2:B:41:MET:CE	2.58	0.77
2:D:148:ILE:HD12	4:D:2075:HOH:O	1.84	0.77
2:B:138:THR:CG2	2:B:176[A]:THR:HG21	2.14	0.76
2:D:92:ARG:HG3	4:D:2054:HOH:O	1.86	0.76
2:H:232:GLU:HG2	2:L:232:GLU:CG	2.16	0.75
2:C:232:GLU:HG2	2:E:232:GLU:HG2	1.68	0.74
1:A:70:MET:HG3	4:A:2081:HOH:O	1.86	0.74
2:K:199:GLU:HG3	4:K:2124:HOH:O	1.86	0.74
2:K:152:ASN:HD22	2:K:153:LEU:H	1.36	0.74
2:B:24:LEU:HD23	2:B:41:MET:CE	2.14	0.73
2:K:16:GLU:HG2	2:K:16:GLU:O	1.87	0.72
2:H:236:ASP:OD2	2:K:225:ARG:HD2	1.89	0.71
2:G:232:GLU:HG3	4:G:2189:HOH:O	1.90	0.71
2:H:184:GLU:HG2	4:H:2106:HOH:O	1.89	0.70
1:A:96[A]:GLN:OE1	4:A:2079:HOH:O	2.09	0.70
2:B:92:ARG:NH2	2:D:49:GLU:OE1	2.24	0.70
2:G:232:GLU:HG2	2:J:232:GLU:HG2	1.73	0.70
3:I:33:MET:CE	3:I:67:THR:HG21	2.20	0.70
2:H:152:ASN:HD22	2:H:153:LEU:H	1.40	0.70
2:J:99:TRP:HE1	2:J:152:ASN:HD21	1.40	0.69
2:J:56:VAL:HG22	2:J:108:PRO:HG2	1.74	0.69
2:H:99:TRP:HE1	2:H:152:ASN:HD21	1.42	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:99:TRP:HE1	3:I:152:ASN:HD21	1.40	0.68
2:L:152:ASN:HD22	2:L:153:LEU:H	1.41	0.67
2:B:248:LEU:O	2:B:249:ASP:CB	2.42	0.66
2:H:232:GLU:HG2	2:L:232:GLU:HG2	1.76	0.66
1:A:236:ASP:OD2	2:F:225:ARG:HD2	1.95	0.66
2:B:11:VAL:HG22	2:B:41:MET:CE	2.23	0.66
2:H:225:ARG:HD2	2:K:236:ASP:OD2	1.95	0.66
2:B:11:VAL:CG1	2:B:41:MET:HE1	2.26	0.65
2:B:225:ARG:HD2	2:E:236:ASP:OD2	1.97	0.65
2:D:199:GLU:O	2:D:203:GLU:HG3	1.96	0.65
2:F:56:VAL:HG22	2:F:108:PRO:HG2	1.79	0.64
2:G:236:ASP:OD2	2:L:225:ARG:HD2	1.97	0.64
2:G:152:ASN:HD22	2:G:153:LEU:H	1.45	0.63
2:G:225:ARG:HD2	2:L:236:ASP:OD2	1.98	0.63
2:K:99:TRP:HE1	2:K:152:ASN:HD21	1.47	0.63
2:D:230:THR:H	2:D:233:GLN:HE21	1.47	0.62
2:L:199:GLU:O	2:L:203:GLU:HG3	1.98	0.62
2:G:99:TRP:HE1	2:G:152:ASN:HD21	1.48	0.61
3:I:175:LYS:NZ	4:I:2102:HOH:O	2.33	0.61
2:B:248:LEU:O	2:B:249:ASP:HB3	1.99	0.61
3:I:33:MET:HE3	3:I:67:THR:CG2	2.29	0.61
2:B:11:VAL:HG13	2:B:41:MET:HE1	1.82	0.61
2:G:166:GLN:NE2	4:G:2136:HOH:O	2.34	0.61
2:H:232:GLU:HG3	4:H:2142:HOH:O	2.01	0.61
2:J:74:GLU:O	2:J:77:ARG:HB3	2.00	0.60
1:A:30:ARG:HD2	1:A:65:ALA:HB2	1.81	0.60
2:G:166:GLN:OE1	2:G:185[A]:MET:HG2	2.00	0.60
2:H:232:GLU:HG2	2:L:232:GLU:CB	2.32	0.60
2:J:152:ASN:HD22	2:J:153:LEU:H	1.47	0.60
2:D:26:ARG:NH2	2:D:29:LYS:HD2	2.17	0.60
3:I:152:ASN:HD22	3:I:153:LEU:H	1.51	0.59
2:L:99:TRP:HE1	2:L:152:ASN:HD21	1.50	0.59
2:G:230:THR:HG21	2:G:232:GLU:OE1	2.03	0.59
2:G:232:GLU:HG2	2:J:232:GLU:CG	2.32	0.59
2:H:56:VAL:HG22	2:H:108:PRO:HG2	1.85	0.59
2:H:199:GLU:O	2:H:203:GLU:HG3	2.03	0.59
2:F:99:TRP:HE1	2:F:152:ASN:HD21	1.52	0.58
2:F:152:ASN:HD22	2:F:153:LEU:H	1.49	0.58
2:G:230:THR:CG2	2:G:232:GLU:OE1	2.52	0.58
1:A:225:ARG:HD2	2:F:236:ASP:OD2	2.04	0.57
2:G:242:LEU:HD22	4:G:2195:HOH:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:143:GLU:HB3	2:C:148:ILE:O	2.04	0.57
2:H:184:GLU:CG	4:H:2106:HOH:O	2.49	0.57
3:I:56:VAL:HG21	3:I:205:ALA:HB1	1.85	0.57
2:D:180:GLN:HG3	4:D:2098:HOH:O	2.05	0.56
2:B:138:THR:HG23	2:B:176[A]:THR:HG21	1.82	0.56
2:L:28:GLU:CD	2:L:28:GLU:H	2.13	0.56
2:B:28:GLU:HB2	4:B:2015:HOH:O	2.06	0.55
2:B:11:VAL:CG1	2:B:41:MET:CE	2.85	0.54
2:B:181:LYS:HD3	4:B:2073:HOH:O	2.07	0.54
2:E:149:PRO:HD2	2:F:242:LEU:HD22	1.89	0.54
2:J:166:GLN:NE2	4:J:2099:HOH:O	2.40	0.54
2:C:17:ASP:HB3	2:C:206:ARG:NH1	2.23	0.54
2:L:80:ASP:HA	4:L:2054:HOH:O	2.06	0.54
1:A:199:GLU:OE2	1:A:203:GLU:OE2	2.25	0.54
2:E:92:ARG:HG2	2:E:92:ARG:HH11	1.72	0.54
3:I:166:GLN:NE2	4:I:2098:HOH:O	2.42	0.53
2:L:56:VAL:HG22	2:L:108:PRO:HG2	1.91	0.53
2:B:16:GLU:O	2:B:17:ASP:HB2	2.09	0.53
2:D:74:GLU:O	2:D:78:GLU:HG2	2.09	0.52
2:F:92[B]:ARG:HG3	4:F:2079:HOH:O	2.09	0.52
2:C:236:ASP:OD2	2:D:225:ARG:HD2	2.09	0.52
2:F:100:LYS:HE3	4:F:2079:HOH:O	2.08	0.52
3:I:236:ASP:OD2	2:J:225:ARG:HD2	2.09	0.52
2:G:173:THR:OG1	2:G:175:LYS:CG	2.54	0.52
2:L:27:PRO:HD2	2:L:28:GLU:OE1	2.10	0.52
2:H:30:ARG:HD2	2:H:65:ALA:HB2	1.92	0.52
2:E:71:ASP:OD2	2:E:74:GLU:CG	2.50	0.51
2:F:242:LEU:HD13	2:F:242:LEU:C	2.35	0.51
2:D:75[B]:TYR:CD1	2:D:90:ILE:HG21	2.45	0.51
2:G:30:ARG:HD3	4:G:2070:HOH:O	2.09	0.51
2:K:181:LYS:HD2	4:K:2104:HOH:O	2.11	0.51
2:D:71:ASP:OD2	2:D:74:GLU:HG3	2.11	0.51
2:G:70:MET:HB2	2:G:72:LEU:HD13	1.92	0.51
2:E:71:ASP:CG	2:E:74:GLU:CG	2.64	0.51
2:G:56:VAL:HG11	2:G:205:ALA:HB1	1.93	0.50
2:F:166:GLN:NE2	4:F:2113:HOH:O	2.43	0.50
2:E:166:GLN:NE2	4:E:2115:HOH:O	2.44	0.50
2:H:149:PRO:HD2	3:I:242:LEU:HD22	1.93	0.50
2:H:28:GLU:CD	2:H:28:GLU:H	2.20	0.50
2:K:148:ILE:HD12	4:K:2087:HOH:O	2.10	0.50
2:K:242:LEU:C	2:K:242:LEU:HD13	2.37	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:70:MET:HB2	4:H:2068:HOH:O	2.12	0.50
2:B:11:VAL:HG21	2:B:41:MET:CE	2.20	0.50
2:C:225:ARG:HD2	2:D:236:ASP:OD2	2.12	0.50
2:E:181:LYS:HD2	4:E:2116:HOH:O	2.12	0.49
2:B:143:GLU:HB3	2:B:148:ILE:O	2.12	0.49
2:L:246:ARG:HG2	4:L:2145:HOH:O	2.11	0.49
2:C:17:ASP:HB3	2:C:206:ARG:HH12	1.77	0.49
2:D:72:LEU:HA	2:D:75[A]:TYR:HB2	1.93	0.49
2:L:56:VAL:HG23	2:L:209:LEU:CD2	2.42	0.49
2:L:242:LEU:C	2:L:242:LEU:HD13	2.36	0.49
2:G:173:THR:CB	2:G:175:LYS:HG2	2.42	0.49
2:G:230:THR:HG23	2:G:232:GLU:CD	2.38	0.49
2:D:70:MET:HB2	2:D:72:LEU:HD13	1.94	0.49
2:D:166:GLN:NE2	4:D:2089:HOH:O	2.46	0.49
2:K:152:ASN:ND2	2:K:153:LEU:H	2.08	0.49
1:A:88:GLU:HG2	4:A:2068:HOH:O	2.12	0.48
2:K:99:TRP:NE1	2:K:152:ASN:HD21	2.11	0.48
2:D:66:TRP:CD1	2:D:113:VAL:HA	2.48	0.48
2:B:166:GLN:NE2	4:B:2075:HOH:O	2.46	0.48
1:A:166:GLN:NE2	4:A:2116:HOH:O	2.46	0.48
2:J:66:TRP:CD1	2:J:113:VAL:HA	2.49	0.48
2:L:66:TRP:CD1	2:L:113:VAL:HA	2.48	0.48
2:K:166:GLN:NE2	4:K:2103:HOH:O	2.46	0.48
1:A:14:GLU:OE1	1:A:16:GLU:OE2	2.32	0.48
2:G:92[B]:ARG:HG3	4:G:2098:HOH:O	2.14	0.48
2:B:56[A]:VAL:HG22	2:B:108:PRO:HG2	1.96	0.48
2:F:92[A]:ARG:HG2	2:F:92[A]:ARG:HH11	1.79	0.47
2:H:236:ASP:OD1	2:K:241:LYS:HE3	2.13	0.47
2:J:152:ASN:HD22	2:J:153:LEU:N	2.13	0.47
1:A:232:GLU:HG2	2:D:232:GLU:CG	2.35	0.47
2:B:56[A]:VAL:HG23	2:B:209:LEU:CD2	2.46	0.46
2:B:56[B]:VAL:HG21	2:B:205:ALA:HB1	1.98	0.46
1:A:248:LEU:O	1:A:249:ASP:HB2	2.15	0.46
2:D:230:THR:H	2:D:233:GLN:NE2	2.12	0.46
2:B:138:THR:HG22	2:B:176[A]:THR:CG2	2.45	0.46
2:G:232:GLU:HG2	2:J:232:GLU:CB	2.46	0.46
2:B:181:LYS:CD	4:B:2073:HOH:O	2.63	0.46
2:G:75:TYR:OH	2:G:87:GLN:HG3	2.16	0.46
2:C:56:VAL:HG22	2:C:108:PRO:CG	2.38	0.46
2:E:71:ASP:OD1	2:E:74:GLU:HG3	2.09	0.45
2:D:100:LYS:HE3	4:D:2054:HOH:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:144:ILE:HD12	2:F:215:VAL:HG12	1.97	0.45
2:C:181:LYS:HE3	2:C:181:LYS:HB2	1.35	0.45
1:A:117:CYS:O	1:A:139:PHE:HA	2.17	0.45
2:D:144:ILE:HD11	2:E:216:LEU:HD23	1.98	0.45
2:J:199:GLU:O	2:J:203:GLU:HG3	2.17	0.45
3:I:199:GLU:O	3:I:203:GLU:HG3	2.16	0.45
2:K:181:LYS:HE3	2:K:181:LYS:HB2	1.75	0.45
3:I:99:TRP:NE1	3:I:152:ASN:HD21	2.13	0.45
2:B:181:LYS:HE3	4:B:2080:HOH:O	2.16	0.44
2:G:230:THR:O	2:G:231:TRP:C	2.60	0.44
3:I:232:GLN:H	3:I:232:GLN:NE2	2.15	0.44
1:A:56:VAL:HG11	1:A:205:ALA:HB1	1.98	0.44
2:B:50:GLN:NE2	4:B:2023:HOH:O	2.50	0.44
2:J:37:LEU:C	2:J:37:LEU:HD23	2.43	0.44
2:F:132[B]:ILE:HD13	2:F:200:VAL:HG12	2.00	0.44
2:L:95:SER:HB3	2:L:100:LYS:HE2	1.99	0.44
1:A:215:VAL:HG12	2:C:144:ILE:HD12	2.00	0.44
2:G:206:ARG:HH21	2:G:206:ARG:HD3	1.47	0.44
3:I:75:TYR:OH	3:I:87:GLN:HG3	2.18	0.44
2:G:232:GLU:CD	2:G:232:GLU:H	2.26	0.43
2:B:138:THR:HG22	2:B:176[A]:THR:HG23	1.99	0.43
2:D:70:MET:CB	2:D:72:LEU:HD13	2.48	0.43
1:A:143:GLU:HB3	1:A:148:ILE:O	2.18	0.43
2:B:138:THR:HG23	2:B:176[A]:THR:HG22	1.97	0.43
2:C:115:GLY:O	2:C:116:TRP:C	2.61	0.43
2:C:225:ARG:HG2	2:C:237:TYR:CZ	2.54	0.43
2:L:99:TRP:NE1	2:L:152:ASN:HD21	2.15	0.43
2:D:72:LEU:HA	2:D:75[B]:TYR:HB3	1.99	0.43
2:D:202:ILE:O	2:D:206:ARG:HG3	2.19	0.43
2:H:152:ASN:ND2	2:H:153:LEU:H	2.12	0.43
3:I:96:GLN:HA	3:I:100:LYS:HB2	2.00	0.43
1:A:242:LEU:HD13	1:A:242:LEU:C	2.43	0.43
2:E:66:TRP:CD1	2:E:113:VAL:HA	2.54	0.43
2:E:134:ALA:HA	2:E:192:VAL:O	2.19	0.43
3:I:225:ARG:HD2	2:J:236:ASP:OD2	2.18	0.43
2:C:185:MET:HG2	4:C:2106:HOH:O	2.18	0.43
2:H:166:GLN:NE2	4:H:2099:HOH:O	2.52	0.43
2:B:242:LEU:HD22	4:B:2114:HOH:O	2.18	0.43
2:L:30:ARG:HD2	2:L:65:ALA:HB2	2.01	0.43
2:D:75[B]:TYR:OH	2:D:87:GLN:HG3	2.19	0.42
3:I:13:VAL:HG11	3:I:48:LEU:HD21	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:70:MET:HG3	4:B:2048:HOH:O	2.19	0.42
2:F:134:ALA:HA	2:F:192:VAL:O	2.18	0.42
2:J:134:ALA:HA	2:J:192:VAL:O	2.18	0.42
2:L:242:LEU:HD22	4:L:2145:HOH:O	2.20	0.42
3:I:241:LYS:HD2	3:I:241:LYS:HA	1.84	0.42
2:H:144:ILE:HD12	3:I:215:VAL:HG12	2.00	0.42
2:E:196:GLN:NE2	4:E:2137:HOH:O	2.52	0.42
2:H:92:ARG:HH22	2:J:49:GLU:CD	2.19	0.42
3:I:49:GLU:OE2	2:L:92:ARG:NH2	2.53	0.42
2:J:22:VAL:HG11	2:J:41:MET:HE1	2.02	0.42
2:B:66:TRP:CD1	2:B:113:VAL:HA	2.55	0.42
2:C:148:ILE:HG22	2:C:149:PRO:O	2.20	0.42
3:I:117:CYS:O	3:I:139:PHE:HA	2.20	0.42
2:L:166:GLN:NE2	4:L:2097:HOH:O	2.53	0.42
1:A:246:ARG:HG2	2:F:247:LEU:HD11	2.02	0.41
2:F:70:MET:HE2	2:F:70:MET:HB3	1.67	0.41
3:I:225:ARG:HG2	3:I:237:TYR:CZ	2.55	0.41
2:K:56[A]:VAL:HG23	2:K:209:LEU:CD2	2.51	0.41
2:C:245:SER:O	2:C:249:ASP:HB2	2.20	0.41
2:F:64:GLU:O	2:F:115:GLY:HA3	2.19	0.41
2:J:99:TRP:NE1	2:J:152:ASN:HD21	2.12	0.41
2:J:238:LEU:HD23	2:J:238:LEU:HA	1.91	0.41
2:F:46:GLU:OE2	4:F:2038:HOH:O	2.22	0.41
2:J:30:ARG:HD2	2:J:65:ALA:HB2	2.02	0.41
2:K:216:LEU:O	2:K:220:LYS:HB2	2.21	0.41
1:A:66:TRP:CD1	1:A:113:VAL:HA	2.56	0.41
2:E:56:VAL:HG11	2:E:205:ALA:HB1	2.02	0.41
2:L:152:ASN:ND2	2:L:153:LEU:H	2.14	0.41
2:B:138:THR:CG2	2:B:176[A]:THR:HG23	2.48	0.41
2:D:75[B]:TYR:HD1	2:D:90:ILE:HG21	1.84	0.41
2:H:232:GLU:HG2	2:L:232:GLU:HB2	2.03	0.41
3:I:74:GLU:OE1	3:I:77:ARG:NH1	2.54	0.41
2:K:132:ILE:HD12	2:K:190:GLU:HB3	2.03	0.41
2:L:185:MET:HB2	2:L:185:MET:HE2	1.84	0.41
3:I:134:ALA:HA	3:I:192:VAL:O	2.21	0.41
2:J:92:ARG:HG2	2:J:92:ARG:HH11	1.86	0.41
2:L:143:GLU:HB3	2:L:148:ILE:O	2.21	0.41
2:J:117:CYS:O	2:J:139:PHE:HA	2.21	0.40
2:C:92[A]:ARG:NH2	2:F:49:GLU:OE1	2.45	0.40
3:I:143:GLU:HB3	3:I:148:ILE:O	2.21	0.40
2:J:81:ALA:HB3	4:J:2059:HOH:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:212:ASN:HA	2:L:213:PRO:HD3	1.95	0.40
1:A:166:GLN:OE1	1:A:185[A]:MET:HG2	2.22	0.40
2:B:92:ARG:HH22	2:D:49:GLU:CD	2.27	0.40
2:E:89:LYS:HE2	2:E:93:GLU:OE2	2.21	0.40
2:F:241:LYS:HE2	2:F:244:GLN:OE1	2.21	0.40
2:H:143:GLU:HB3	2:H:148:ILE:O	2.22	0.40
2:J:13:VAL:HG11	2:J:48:LEU:HD21	2.03	0.40
2:L:56:VAL:HG23	2:L:209:LEU:HD21	2.04	0.40

All (1) 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)
2:C:29:LYS:NZ	2:K:17:ASP:OD1[2_664]	1.89	0.31

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	248/276 (90%)	238 (96%)	9 (4%)	1 (0%)	30	19
2	B	247/276 (90%)	237 (96%)	9 (4%)	1 (0%)	30	19
2	C	246/276 (89%)	238 (97%)	8 (3%)	0	100	100
2	D	249/276 (90%)	240 (96%)	9 (4%)	0	100	100
2	E	248/276 (90%)	239 (96%)	8 (3%)	1 (0%)	30	19
2	F	246/276 (89%)	240 (98%)	6 (2%)	0	100	100
2	G	247/276 (90%)	241 (98%)	6 (2%)	0	100	100
2	H	246/276 (89%)	236 (96%)	9 (4%)	1 (0%)	30	19
2	J	245/276 (89%)	236 (96%)	9 (4%)	0	100	100
2	K	247/276 (90%)	237 (96%)	9 (4%)	1 (0%)	30	19

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	L	245/276 (89%)	236 (96%)	9 (4%)	0	100	100
3	I	245/276 (89%)	237 (97%)	8 (3%)	0	100	100
All	All	2959/3312 (89%)	2855 (96%)	99 (3%)	5 (0%)	43	31

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	249	ASP
2	E	249	ASP
2	H	249	ASP
1	A	249	ASP
2	K	75	TYR

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	205/227 (90%)	197 (96%)	8 (4%)	28	16
2	B	203/227 (89%)	190 (94%)	13 (6%)	16	6
2	C	204/227 (90%)	194 (95%)	10 (5%)	22	10
2	D	206/227 (91%)	192 (93%)	14 (7%)	14	5
2	E	205/227 (90%)	194 (95%)	11 (5%)	20	8
2	F	203/227 (89%)	191 (94%)	12 (6%)	18	7
2	G	205/227 (90%)	192 (94%)	13 (6%)	16	6
2	H	203/227 (89%)	188 (93%)	15 (7%)	13	4
2	J	202/227 (89%)	181 (90%)	21 (10%)	7	2
2	K	204/227 (90%)	190 (93%)	14 (7%)	14	5
2	L	202/227 (89%)	183 (91%)	19 (9%)	8	2
3	I	202/227 (89%)	190 (94%)	12 (6%)	18	7
All	All	2444/2724 (90%)	2282 (93%)	162 (7%)	15	5

All (162) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	GLU
1	A	30	ARG
1	A	70	MET
1	A	73	LYS
1	A	184	GLU
1	A	216	LEU
1	A	225	ARG
1	A	242	LEU
2	B	28	GLU
2	B	30	ARG
2	B	72	LEU
2	B	73	LYS
2	B	77	ARG
2	B	181	LYS
2	B	184	GLU
2	B	187	LEU
2	B	196	GLN
2	B	199	GLU
2	B	216	LEU
2	B	220	LYS
2	B	242	LEU
2	C	5	GLU
2	C	28	GLU
2	C	29	LYS
2	C	72	LEU
2	C	73	LYS
2	C	181	LYS
2	C	196	GLN
2	C	206	ARG
2	C	216	LEU
2	C	225	ARG
2	D	5	GLU
2	D	9	LYS
2	D	28	GLU
2	D	72	LEU
2	D	74	GLU
2	D	77	ARG
2	D	78	GLU
2	D	175	LYS
2	D	180	GLN
2	D	196	GLN
2	D	216	LEU

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Mol	Chain	Res	Type
2	D	225	ARG
2	D	229	LEU
2	D	242	LEU
2	E	3	THR
2	E	28	GLU
2	E	70	MET
2	E	72	LEU
2	E	73	LYS
2	E	77	ARG
2	E	85	ILE
2	E	184	GLU
2	E	216	LEU
2	E	229	LEU
2	E	246	ARG
2	F	5	GLU
2	F	28	GLU
2	F	30	ARG
2	F	73	LYS
2	F	79	VAL
2	F	136	GLU
2	F	141	LEU
2	F	148	ILE
2	F	190	GLU
2	F	216	LEU
2	F	225	ARG
2	F	242	LEU
2	G	3	THR
2	G	5	GLU
2	G	30	ARG
2	G	72	LEU
2	G	88	GLU
2	G	141	LEU
2	G	143	GLU
2	G	144	ILE
2	G	148	ILE
2	G	181	LYS
2	G	229	LEU
2	G	230	THR
2	G	232	GLU
2	H	5	GLU
2	H	28	GLU
2	H	30	ARG

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Mol	Chain	Res	Type
2	H	70	MET
2	H	72	LEU
2	H	73	LYS
2	H	85	ILE
2	H	136	GLU
2	H	187	LEU
2	H	196	GLN
2	H	199	GLU
2	H	216	LEU
2	H	229	LEU
2	H	232	GLU
2	H	242	LEU
3	I	28	GLU
3	I	30	ARG
3	I	72	LEU
3	I	79	VAL
3	I	136	GLU
3	I	148	ILE
3	I	196	GLN
3	I	216	LEU
3	I	225	ARG
3	I	229	LEU
3	I	232	GLN
3	I	242	LEU
2	J	5	GLU
2	J	9	LYS
2	J	28	GLU
2	J	29	LYS
2	J	30	ARG
2	J	72	LEU
2	J	73	LYS
2	J	74	GLU
2	J	77	ARG
2	J	136	GLU
2	J	141	LEU
2	J	146	TRP
2	J	148	ILE
2	J	181	LYS
2	J	184	GLU
2	J	187	LEU
2	J	196	GLN
2	J	216	LEU

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Mol	Chain	Res	Type
2	J	225	ARG
2	J	229	LEU
2	J	242	LEU
2	K	3	THR
2	K	16	GLU
2	K	28	GLU
2	K	72	LEU
2	K	73	LYS
2	K	74	GLU
2	K	77	ARG
2	K	85	ILE
2	K	184	GLU
2	K	187	LEU
2	K	216	LEU
2	K	220	LYS
2	K	225	ARG
2	K	229	LEU
2	L	5	GLU
2	L	16	GLU
2	L	28	GLU
2	L	30	ARG
2	L	70	MET
2	L	72	LEU
2	L	73	LYS
2	L	85	ILE
2	L	141	LEU
2	L	148	ILE
2	L	181	LYS
2	L	184	GLU
2	L	187	LEU
2	L	199	GLU
2	L	216	LEU
2	L	220	LYS
2	L	225	ARG
2	L	229	LEU
2	L	242	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	50	GLN

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Mol	Chain	Res	Type
1	A	180	GLN
2	B	50	GLN
2	B	166	GLN
2	C	166	GLN
2	D	25	ASN
2	D	50	GLN
2	D	166	GLN
2	D	233	GLN
2	E	25	ASN
2	E	145	ASN
2	E	166	GLN
2	E	207	ASN
2	F	152	ASN
2	F	166	GLN
2	F	207	ASN
2	G	152	ASN
2	G	166	GLN
2	G	196	GLN
2	G	207	ASN
2	H	50	GLN
2	H	152	ASN
2	H	166	GLN
2	H	207	ASN
3	I	25	ASN
3	I	50	GLN
3	I	152	ASN
3	I	166	GLN
3	I	196	GLN
3	I	207	ASN
3	I	232	GLN
2	J	25	ASN
2	J	152	ASN
2	J	166	GLN
2	K	25	ASN
2	K	87	GLN
2	K	152	ASN
2	K	166	GLN
2	K	196	GLN
2	L	25	ASN
2	L	152	ASN
2	L	166	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	247/276 (89%)	-0.30	6 (2%)	59	60	8, 19, 31, 40	10 (4%)
2	B	245/276 (88%)	0.01	7 (2%)	53	53	13, 22, 33, 44	20 (8%)
2	C	246/276 (89%)	-0.24	3 (1%)	76	77	8, 20, 30, 38	14 (5%)
2	D	248/276 (89%)	-0.17	11 (4%)	39	38	11, 19, 33, 45	15 (6%)
2	E	247/276 (89%)	-0.02	12 (4%)	35	33	4, 19, 30, 44	21 (8%)
2	F	246/276 (89%)	-0.27	6 (2%)	59	60	9, 18, 30, 41	14 (5%)
2	G	246/276 (89%)	-0.34	3 (1%)	76	77	10, 18, 30, 38	7 (2%)
2	H	247/276 (89%)	-0.08	5 (2%)	65	65	12, 21, 32, 44	13 (5%)
2	J	246/276 (89%)	-0.11	9 (3%)	45	45	11, 20, 31, 40	17 (6%)
2	K	247/276 (89%)	-0.16	11 (4%)	38	37	11, 19, 31, 39	21 (8%)
2	L	246/276 (89%)	-0.11	11 (4%)	38	37	11, 20, 31, 44	22 (8%)
3	I	246/276 (89%)	-0.23	4 (1%)	70	71	10, 21, 31, 39	9 (3%)
All	All	2957/3312 (89%)	-0.17	88 (2%)	52	52	4, 20, 31, 45	183 (6%)

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	250	THR	10.1
2	B	250	THR	7.1
1	A	3	THR	6.3
2	E	72	LEU	6.3
2	E	81	ALA	6.2
2	C	249	ASP	6.2
1	A	146	TRP	5.6
2	E	75	TYR	5.5
2	D	75[A]	TYR	4.9
2	E	76	PHE	4.9
2	J	146	TRP	4.7

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Mol	Chain	Res	Type	RSRZ
2	G	230	THR	4.7
2	E	146	TRP	4.5
2	K	76	PHE	4.5
2	F	81	ALA	4.4
2	J	76	PHE	4.4
2	D	146	TRP	4.3
2	J	81	ALA	4.2
2	F	76	PHE	4.2
2	D	72	LEU	4.1
2	F	75	TYR	4.0
2	E	79	VAL	3.9
2	K	75	TYR	3.7
2	B	249	ASP	3.7
2	K	79	VAL	3.6
2	D	80	ASP	3.6
2	L	76	PHE	3.5
2	D	76	PHE	3.5
2	J	249	ASP	3.5
1	A	249	ASP	3.3
2	E	80	ASP	3.3
2	K	81	ALA	3.3
2	B	146	TRP	3.3
2	H	146	TRP	3.2
2	F	77	ARG	3.2
2	D	81	ALA	3.1
2	J	80	ASP	3.0
2	D	79	VAL	3.0
2	G	3	THR	3.0
1	A	70	MET	3.0
2	F	72	LEU	2.9
2	K	77	ARG	2.9
2	H	76	PHE	2.9
2	E	74	GLU	2.9
2	B	76	PHE	2.9
2	L	81	ALA	2.8
2	C	146	TRP	2.8
2	E	77	ARG	2.8
2	L	146	TRP	2.8
2	J	79	VAL	2.8
2	B	6	GLY	2.7
2	B	17	ASP	2.7
2	E	249	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
3	I	80	ASP	2.6
2	K	28	GLU	2.6
2	H	72	LEU	2.6
2	D	251	GLU	2.6
2	K	78	GLU	2.6
2	L	249	ASP	2.6
2	E	3	THR	2.6
2	L	72	LEU	2.5
2	L	15	ILE	2.5
2	J	75	TYR	2.4
2	E	71	ASP	2.4
2	J	77	ARG	2.4
2	G	146	TRP	2.4
3	I	249	ASP	2.3
3	I	72	LEU	2.3
2	K	80	ASP	2.3
2	L	80	ASP	2.3
2	D	78	GLU	2.3
3	I	76	PHE	2.3
2	L	79	VAL	2.3
2	H	70	MET	2.3
2	F	78	GLU	2.2
2	L	28	GLU	2.2
2	C	203	GLU	2.1
2	L	74	GLU	2.1
2	K	72	LEU	2.1
1	A	143	GLU	2.1
2	D	250	THR	2.1
2	K	146	TRP	2.1
2	L	71	ASP	2.1
1	A	148	ILE	2.0
2	J	72	LEU	2.0
2	D	77	ARG	2.0
2	B	115	GLY	2.0
2	K	3	THR	2.0

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

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

6.3 Carbohydrates [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.