



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 11:43 AM UTC

PDB ID : 5MU0 / pdb_00005mu0
Title : ACC1 Fab fragment in complex with citrullinated C1 epitope of CII (IA03)
Authors : Dobritsch, D.; Holmdahl, R.; Ge, C.
Deposited on : 2017-01-11
Resolution : 2.70 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

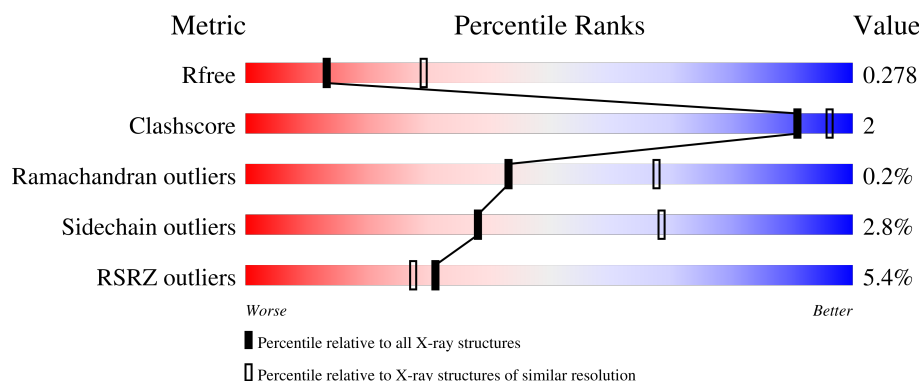
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	218	2% 93% 5%
1	C	218	2% 89% 8%
1	E	218	4% 98% .
1	G	218	% 94% 5%
1	I	218	8% 94% 5%

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Mol	Chain	Length	Quality of chain
1	K	218	
1	M	218	
1	O	218	
2	B	218	
2	D	218	
2	F	218	
2	H	218	
2	J	218	
2	L	218	
2	N	218	
2	P	218	
3	Q	18	
3	R	18	
3	S	18	
3	T	18	
3	U	18	
3	V	18	
3	W	18	
3	X	18	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 26780 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 heavy chain of ACC1 antibody Fab fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	0	0	0
			1618	1021	269	320	8			
1	C	212	Total	C	N	O	S	0	0	0
			1600	1011	264	317	8			
1	E	218	Total	C	N	O	S	0	0	0
			1640	1034	273	325	8			
1	G	215	Total	C	N	O	S	0	0	0
			1622	1023	270	321	8			
1	I	216	Total	C	N	O	S	0	0	0
			1629	1028	271	322	8			
1	K	215	Total	C	N	O	S	0	0	0
			1625	1026	270	321	8			
1	M	212	Total	C	N	O	S	0	0	0
			1600	1011	264	317	8			
1	O	211	Total	C	N	O	S	0	0	0
			1589	1005	262	314	8			

- Molecule 2 is a protein called light chain of ACC1 antibody Fab fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	217	Total	C	N	O	S	0	0	0
			1660	1035	281	338	6			
2	D	217	Total	C	N	O	S	0	0	0
			1660	1035	281	338	6			
2	F	216	Total	C	N	O	S	0	0	0
			1651	1030	280	335	6			
2	H	217	Total	C	N	O	S	0	0	0
			1660	1035	281	338	6			
2	J	217	Total	C	N	O	S	0	0	0
			1660	1035	281	338	6			
2	L	217	Total	C	N	O	S	0	0	0
			1660	1035	281	338	6			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	N	215	Total	C	N	O	S	0	0	0
			1643	1026	278	333	6			
2	P	216	Total	C	N	O	S	0	0	0
			1651	1030	280	335	6			

- Molecule 3 is a protein called IA03 peptide containing the citrullinated C1 epitope of collagen type II, Collagen alpha-1(II) chain, IA03 peptide containing the citrullinated C1 epitope of collagen type II.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	Q	11	Total	C	N	O	0	0	0
			74	43	16	15			
3	R	10	Total	C	N	O	0	0	0
			66	38	15	13			
3	S	10	Total	C	N	O	0	0	0
			66	38	15	13			
3	T	10	Total	C	N	O	0	0	0
			66	38	15	13			
3	U	10	Total	C	N	O	0	0	0
			66	38	15	13			
3	V	10	Total	C	N	O	0	0	0
			66	38	15	13			
3	W	9	Total	C	N	O	0	0	0
			62	36	14	12			
3	X	10	Total	C	N	O	0	0	0
			66	38	15	13			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	7	Total	O	0	0
			7	7		
4	B	6	Total	O	0	0
			6	6		
4	C	9	Total	O	0	0
			9	9		
4	D	20	Total	O	0	0
			20	20		
4	E	5	Total	O	0	0
			5	5		
4	F	3	Total	O	0	0
			3	3		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	5	Total 5	O 5	0	0
4	H	7	Total 7	O 7	0	0
4	I	4	Total 4	O 4	0	0
4	J	6	Total 6	O 6	0	0
4	K	1	Total 1	O 1	0	0
4	L	2	Total 2	O 2	0	0
4	M	2	Total 2	O 2	0	0
4	N	1	Total 1	O 1	0	0
4	T	1	Total 1	O 1	0	0
4	V	1	Total 1	O 1	0	0

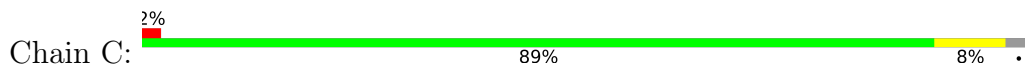
3 Residue-property plots [i](#)

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.

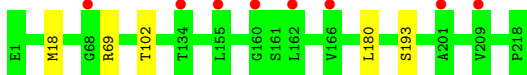
- Molecule 1: heavy chain of ACC1 antibody Fab fragment



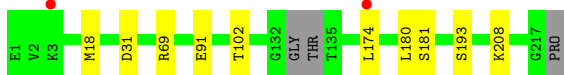
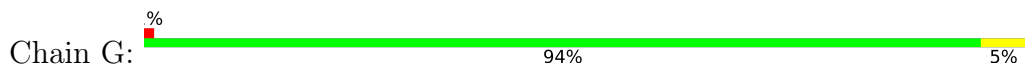
- Molecule 1: heavy chain of ACC1 antibody Fab fragment



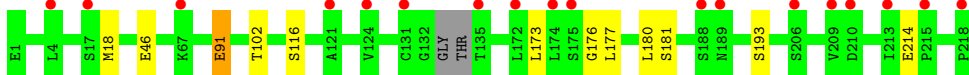
- Molecule 1: heavy chain of ACC1 antibody Fab fragment



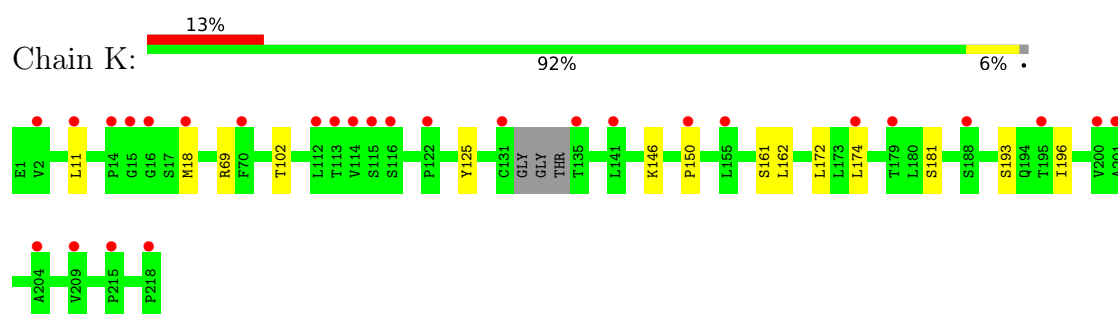
- Molecule 1: heavy chain of ACC1 antibody Fab fragment



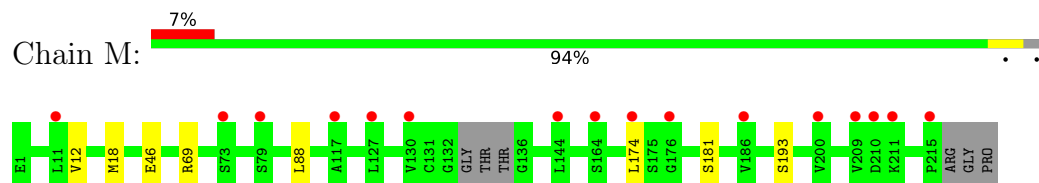
- Molecule 1: heavy chain of ACC1 antibody Fab fragment



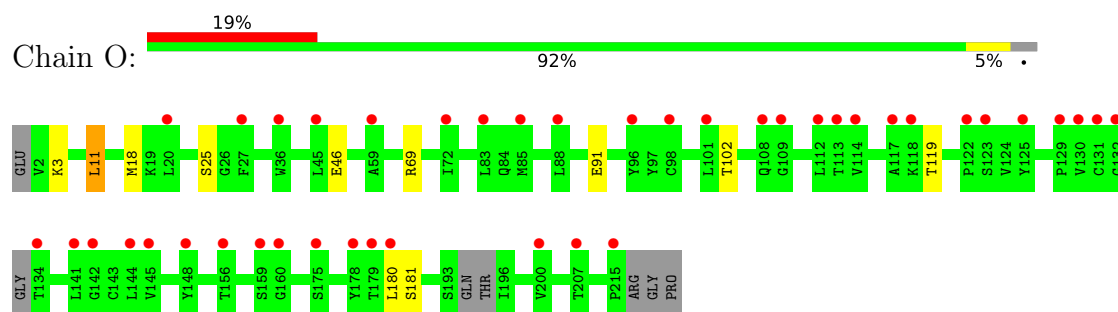
- Molecule 1: heavy chain of ACC1 antibody Fab fragment



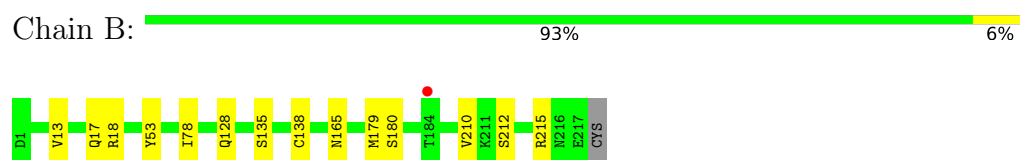
- Molecule 1: heavy chain of ACC1 antibody Fab fragment



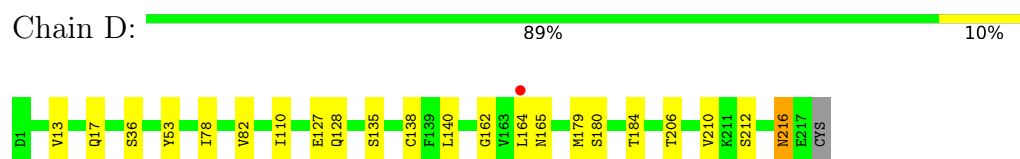
- Molecule 1: heavy chain of ACC1 antibody Fab fragment



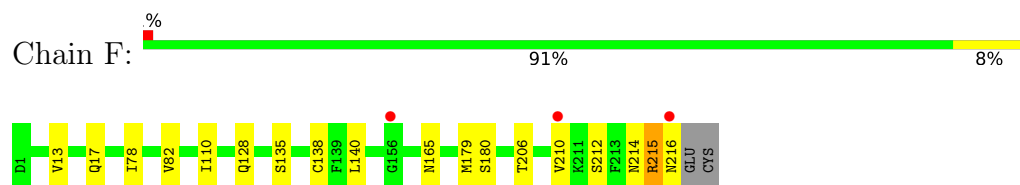
- Molecule 2: light chain of ACC1 antibody Fab fragment



- Molecule 2: light chain of ACC1 antibody Fab fragment

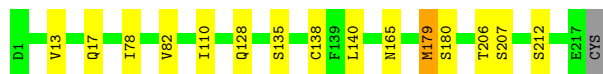


- Molecule 2: light chain of ACC1 antibody Fab fragment



- Molecule 2: light chain of ACC1 antibody Fab fragment

Chain H:  93% 6%



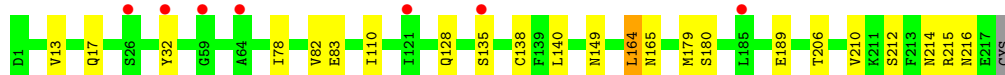
- Molecule 2: light chain of ACC1 antibody Fab fragment

Chain J:  2% 90% 10%

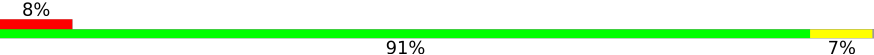


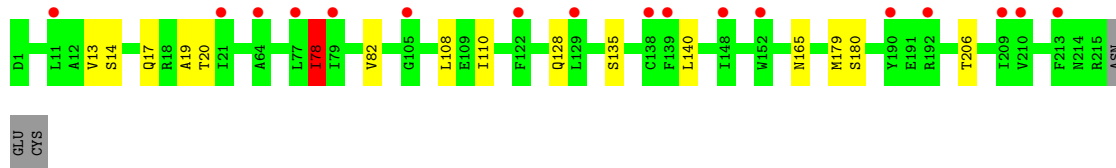
- Molecule 2: light chain of ACC1 antibody Fab fragment

Chain L:  3% 89% 10%



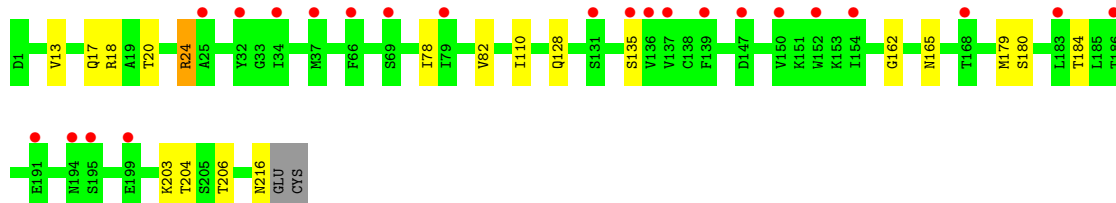
- Molecule 2: light chain of ACC1 antibody Fab fragment

Chain N:  8% 91% 7%

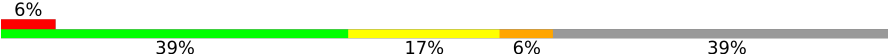


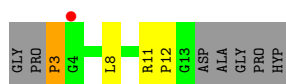
- Molecule 2: light chain of ACC1 antibody Fab fragment

Chain P:  11% 90% 8%

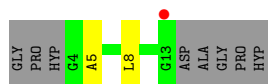
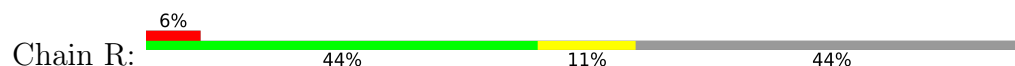


- Molecule 3: IA03 peptide containing the citrullinated C1 epitope of collagen type II, Collagen alpha-1(II) chain, IA03 peptide containing the citrullinated C1 epitope of collagen type II

Chain Q:  6% 39% 17% 6% 39%



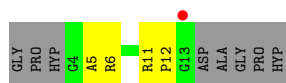
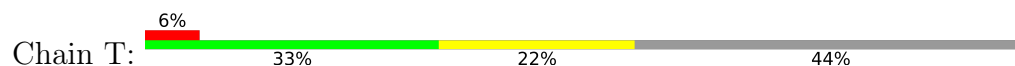
- Molecule 3: IA03 peptide containing the citrullinated C1 epitope of collagen type II,Collagen alpha-1(II) chain,IA03 peptide containing the citrullinated C1 epitope of collagen type II



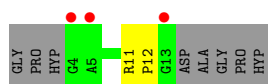
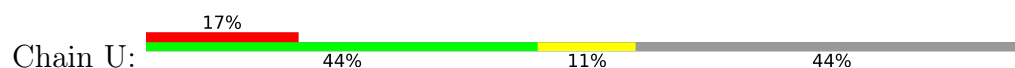
- Molecule 3: IA03 peptide containing the citrullinated C1 epitope of collagen type II,Collagen alpha-1(II) chain,IA03 peptide containing the citrullinated C1 epitope of collagen type II



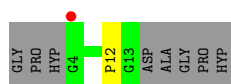
- Molecule 3: IA03 peptide containing the citrullinated C1 epitope of collagen type II,Collagen alpha-1(II) chain,IA03 peptide containing the citrullinated C1 epitope of collagen type II



- Molecule 3: IA03 peptide containing the citrullinated C1 epitope of collagen type II,Collagen alpha-1(II) chain,IA03 peptide containing the citrullinated C1 epitope of collagen type II



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- Molecule 3: IA03 peptide containing the citrullinated C1 epitope of collagen type II,Collagen alpha-1(II) chain,IA03 peptide containing the citrullinated C1 epitope of collagen type II





- Molecule 3: IA03 peptide containing the citrullinated C1 epitope of collagen type II,Collagen alpha-1(II) chain,IA03 peptide containing the citrullinated C1 epitope of collagen type II



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	87.15Å 93.05Å 121.75Å 92.75° 106.45° 104.37°	Depositor
Resolution (Å)	78.15 – 2.70 78.15 – 2.70	Depositor EDS
% Data completeness (in resolution range)	89.2 (78.15-2.70) 89.2 (78.15-2.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.61 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.252 , 0.279 0.253 , 0.278	Depositor DCC
R_{free} test set	4409 reflections (4.55%)	wwPDB-VP
Wilson B-factor (Å ²)	54.8	Xtriage
Anisotropy	0.338	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 36.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	26780	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.82% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HYP, CIR

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	0.65	0/1657	0.79	2/2260 (0.1%)
1	C	0.66	0/1639	0.80	0/2236
1	E	0.65	0/1681	0.79	1/2295 (0.0%)
1	G	0.66	0/1661	0.78	1/2265 (0.0%)
1	I	0.63	0/1669	0.78	1/2277 (0.0%)
1	K	0.63	0/1665	0.78	1/2272 (0.0%)
1	M	0.61	0/1639	0.76	1/2236 (0.0%)
1	O	0.63	0/1627	0.90	4/2219 (0.2%)
2	B	0.66	0/1697	0.82	0/2302
2	D	0.70	1/1697 (0.1%)	0.82	0/2302
2	F	0.62	0/1688	0.81	1/2290 (0.0%)
2	H	0.75	0/1697	0.86	1/2302 (0.0%)
2	J	0.70	0/1697	0.85	1/2302 (0.0%)
2	L	0.67	0/1697	0.85	2/2302 (0.1%)
2	N	0.62	0/1680	0.80	1/2279 (0.0%)
2	P	0.62	0/1688	0.80	2/2290 (0.1%)
3	Q	1.04	0/45	1.25	0/56
3	R	1.10	0/45	1.43	0/56
3	S	0.79	0/45	1.07	0/56
3	T	1.02	0/45	1.57	1/56 (1.8%)
3	U	1.12	0/45	1.43	0/56
3	V	0.97	0/45	1.31	0/56
3	W	0.94	0/41	1.58	1/51 (2.0%)
3	X	1.21	0/45	1.34	1/56 (1.8%)
All	All	0.66	1/27135 (0.0%)	0.82	22/36872 (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
3	Q	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	216	ASN	C-N	5.21	1.40	1.33

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	O	69	ARG	NE-CZ-NH2	-12.61	107.85	119.20
1	O	69	ARG	NE-CZ-NH1	11.13	132.63	121.50
1	O	69	ARG	CD-NE-CZ	8.39	136.15	124.40
2	J	207	SER	N-CA-C	-7.61	100.43	110.40
2	N	78	ILE	CB-CG1-CD1	7.28	129.09	113.80
3	W	6	ARG	N-CA-C	-6.26	104.38	111.14
2	H	212	SER	CA-CB-OG	6.07	123.25	111.10
3	X	7	GLY	N-CA-C	5.90	118.06	112.04
2	F	215	ARG	N-CA-C	5.90	119.30	111.75
1	I	91	GLU	CA-CB-CG	5.80	125.70	114.10
3	T	6	ARG	NE-CZ-NH2	5.76	124.38	119.20
2	L	215	ARG	CA-CB-CG	5.73	125.56	114.10
1	K	69	ARG	NE-CZ-NH2	5.29	123.96	119.20
1	E	69	ARG	NE-CZ-NH2	5.14	123.83	119.20
1	A	69	ARG	NE-CZ-NH2	5.12	123.81	119.20
2	P	24	ARG	CA-CB-CG	5.09	124.28	114.10
1	O	69	ARG	CG-CD-NE	-5.06	100.88	112.00
2	L	83	GLU	CA-CB-CG	5.05	124.20	114.10
1	M	69	ARG	NE-CZ-NH2	5.04	123.74	119.20
2	P	204	THR	N-CA-C	5.04	117.51	111.71
1	A	174	LEU	CA-CB-CG	5.04	133.93	116.30
1	G	69	ARG	NE-CZ-NH2	5.01	123.71	119.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	Q	3	HYP	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	1618	0	1580	2	2
1	C	1600	0	1560	9	0
1	E	1640	0	1601	1	0
1	G	1622	0	1583	2	0
1	I	1629	0	1590	2	1
1	K	1625	0	1587	6	1
1	M	1600	0	1560	2	0
1	O	1589	0	1549	6	0
2	B	1660	0	1593	7	0
2	D	1660	0	1593	10	0
2	F	1651	0	1587	6	1
2	H	1660	0	1593	6	1
2	J	1660	0	1593	8	1
2	L	1660	0	1593	13	2
2	N	1643	0	1581	12	0
2	P	1651	0	1587	8	1
3	Q	74	0	69	1	0
3	R	66	0	61	2	0
3	S	66	0	61	0	0
3	T	66	0	61	1	0
3	U	66	0	61	0	0
3	V	66	0	60	0	0
3	W	62	0	57	0	0
3	X	66	0	60	0	0
4	A	7	0	0	0	0
4	B	6	0	0	0	0
4	C	9	0	0	0	0
4	D	20	0	0	0	0
4	E	5	0	0	0	0
4	F	3	0	0	0	0
4	G	5	0	0	0	0
4	H	7	0	0	0	0
4	I	4	0	0	0	0
4	J	6	0	0	0	0
4	K	1	0	0	0	0
4	L	2	0	0	0	0
4	M	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	N	1	0	0	0	0
4	T	1	0	0	0	0
4	V	1	0	0	0	0
All	All	26780	0	25820	91	5

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 2.

All (91) 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:O:11:LEU:HD12	1:O:11:LEU:O	1.87	0.74
1:O:3:LYS:HG2	1:O:25:SER:OG	1.88	0.73
1:O:11:LEU:HD12	1:O:11:LEU:C	2.21	0.66
1:O:11:LEU:HD21	1:O:119:THR:N	2.14	0.63
1:K:172:LEU:HD11	2:L:165:ASN:O	2.01	0.60
1:M:12:VAL:HG11	1:M:88:LEU:CD1	2.34	0.58
1:C:12:VAL:HG11	1:C:88:LEU:CD1	2.35	0.57
2:N:20:THR:HG23	2:N:78:ILE:CD1	2.36	0.56
1:M:12:VAL:HG11	1:M:88:LEU:HD13	1.87	0.56
2:N:20:THR:HG23	2:N:78:ILE:HD12	1.88	0.54
2:B:53:TYR:CE1	3:Q:8:LEU:HD13	2.44	0.53
1:C:12:VAL:HG11	1:C:88:LEU:HD13	1.90	0.52
2:J:98:VAL:O	2:P:24:ARG:NH1	2.41	0.52
2:J:82:VAL:CG1	2:J:110:ILE:HD12	2.40	0.52
2:L:164:LEU:C	2:L:164:LEU:HD23	2.35	0.52
2:B:18:ARG:HH21	2:L:32:TYR:HB3	1.75	0.51
2:H:165:ASN:HB3	2:H:179:MET:HE3	1.92	0.51
2:N:165:ASN:HB3	2:N:179:MET:HE3	1.94	0.50
2:D:165:ASN:HB3	2:D:179:MET:HE3	1.94	0.50
2:L:165:ASN:HB3	2:L:179:MET:HE3	1.93	0.49
2:N:13:VAL:HG22	2:N:108:LEU:HD11	1.93	0.49
2:P:165:ASN:HB3	2:P:179:MET:HE3	1.94	0.48
1:C:31:ASP:O	3:R:5:ALA:HB1	2.13	0.48
2:D:82:VAL:CG1	2:D:110:ILE:HD12	2.43	0.48
2:N:13:VAL:HG21	2:N:82:VAL:HG21	1.96	0.48
2:B:165:ASN:HB3	2:B:179:MET:HE3	1.96	0.47
1:K:162:LEU:CD2	1:K:196:ILE:HD12	2.43	0.47
2:B:18:ARG:NH2	2:L:32:TYR:HB3	2.28	0.47
2:F:82:VAL:CG1	2:F:110:ILE:HD12	2.43	0.47
2:P:162:GLY:O	2:P:184:THR:HG22	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:162:GLY:O	2:D:184:THR:HG22	2.15	0.47
1:K:11:LEU:HB2	1:K:150:PRO:HG3	1.96	0.47
2:L:82:VAL:CG1	2:L:110:ILE:HD12	2.45	0.47
2:N:82:VAL:CG1	2:N:110:ILE:HD12	2.45	0.47
2:J:165:ASN:HB3	2:J:179:MET:HE3	1.97	0.47
2:F:165:ASN:HB3	2:F:179:MET:HE3	1.97	0.46
2:B:78:ILE:HD12	2:B:78:ILE:N	2.31	0.46
2:N:13:VAL:O	2:N:13:VAL:HG23	2.15	0.46
2:H:78:ILE:HD12	2:H:78:ILE:N	2.31	0.46
2:D:78:ILE:N	2:D:78:ILE:HD12	2.32	0.45
2:J:78:ILE:HD12	2:J:78:ILE:N	2.31	0.45
2:J:34:ILE:HD11	2:J:54:ALA:HB1	1.99	0.45
2:F:78:ILE:HD12	2:F:78:ILE:N	2.31	0.45
1:C:11:LEU:HB2	1:C:150:PRO:HG3	1.99	0.44
2:N:13:VAL:HG11	2:N:19:ALA:HB2	2.00	0.44
2:H:82:VAL:CG1	2:H:110:ILE:HD12	2.48	0.44
2:L:78:ILE:HD12	2:L:78:ILE:N	2.32	0.44
2:P:13:VAL:CG1	2:P:17:GLN:HB2	2.48	0.43
2:P:78:ILE:HD12	2:P:78:ILE:N	2.32	0.43
2:N:13:VAL:CG2	2:N:108:LEU:HD11	2.48	0.43
1:A:146:LYS:NZ	1:A:174:LEU:HD21	2.34	0.43
2:H:140:LEU:N	2:H:140:LEU:HD12	2.34	0.43
2:L:13:VAL:CG1	2:L:17:GLN:HB2	2.48	0.42
2:F:13:VAL:CG1	2:F:17:GLN:HB2	2.49	0.42
2:H:13:VAL:CG1	2:H:17:GLN:HB2	2.49	0.42
2:P:82:VAL:CG1	2:P:110:ILE:HD12	2.49	0.42
2:F:128:GLN:HE22	2:F:135:SER:CB	2.33	0.42
2:H:128:GLN:HE22	2:H:135:SER:CB	2.33	0.42
1:O:11:LEU:C	1:O:11:LEU:CD1	2.91	0.42
2:D:13:VAL:CG1	2:D:17:GLN:HB2	2.50	0.42
2:J:13:VAL:CG1	2:J:17:GLN:HB2	2.49	0.42
2:B:13:VAL:CG1	2:B:17:GLN:HB2	2.50	0.41
2:D:128:GLN:HE22	2:D:135:SER:CB	2.33	0.41
1:I:176:GLY:C	1:I:177:LEU:HD12	2.44	0.41
1:I:180:LEU:C	1:I:180:LEU:HD12	2.45	0.41
1:C:195:THR:HG22	1:C:212:LYS:HE3	2.02	0.41
1:C:211:LYS:HE3	2:D:127:GLU:OE1	2.19	0.41
2:D:140:LEU:N	2:D:140:LEU:HD12	2.36	0.41
2:N:140:LEU:HD12	2:N:140:LEU:N	2.36	0.41
1:A:176:GLY:C	1:A:177:LEU:HD12	2.45	0.41
1:C:174:LEU:HD12	2:D:164:LEU:HD13	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:31:ASP:O	3:T:5:ALA:HB1	2.20	0.41
2:J:128:GLN:HE22	2:J:135:SER:CB	2.34	0.41
1:K:174:LEU:HB2	2:L:164:LEU:HD12	2.02	0.41
2:L:128:GLN:HE22	2:L:135:SER:CB	2.33	0.41
2:N:14:SER:HB2	2:N:17:GLN:HG3	2.03	0.41
1:E:180:LEU:C	1:E:180:LEU:HD12	2.46	0.41
2:J:140:LEU:HD12	2:J:140:LEU:N	2.36	0.41
2:L:140:LEU:N	2:L:140:LEU:HD12	2.35	0.41
2:P:128:GLN:HE22	2:P:135:SER:CB	2.33	0.41
2:B:128:GLN:HE22	2:B:135:SER:CB	2.33	0.41
1:K:146:LYS:HE2	2:L:135:SER:OG	2.21	0.41
1:G:180:LEU:C	1:G:180:LEU:HD12	2.45	0.40
2:N:128:GLN:HE22	2:N:135:SER:CB	2.33	0.40
1:O:180:LEU:HD12	1:O:180:LEU:C	2.47	0.40
1:C:12:VAL:HG11	1:C:88:LEU:HD12	2.03	0.40
1:C:176:GLY:C	1:C:177:LEU:HD12	2.45	0.40
2:D:53:TYR:CZ	3:R:8:LEU:HD22	2.56	0.40
1:K:125:TYR:CZ	2:L:128:GLN:HG3	2.56	0.40
2:F:140:LEU:HD12	2:F:140:LEU:N	2.35	0.40
2:P:20:THR:O	2:P:20:THR:HG23	2.21	0.40

All (5) 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:I:116:SER:OG	2:L:189:GLU:OE2[1_556]	1.60	0.60
1:A:116:SER:OG	2:L:214:ASN:ND2[1_556]	1.88	0.32
1:A:78:ARG:O	2:H:207:SER:CB[1_455]	2.00	0.20
2:F:214:ASN:ND2	1:K:161:SER:OG[1_545]	2.11	0.09
2:J:192:ARG:O	2:P:18:ARG:NH2[1_455]	2.19	0.01

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	210/218 (96%)	206 (98%)	3 (1%)	1 (0%)	24	48
1	C	208/218 (95%)	203 (98%)	4 (2%)	1 (0%)	24	48
1	E	216/218 (99%)	211 (98%)	4 (2%)	1 (0%)	24	48
1	G	211/218 (97%)	207 (98%)	3 (1%)	1 (0%)	24	48
1	I	212/218 (97%)	208 (98%)	3 (1%)	1 (0%)	24	48
1	K	211/218 (97%)	207 (98%)	3 (1%)	1 (0%)	24	48
1	M	208/218 (95%)	203 (98%)	5 (2%)	0	100	100
1	O	205/218 (94%)	200 (98%)	4 (2%)	1 (0%)	24	48
2	B	215/218 (99%)	209 (97%)	6 (3%)	0	100	100
2	D	215/218 (99%)	209 (97%)	6 (3%)	0	100	100
2	F	214/218 (98%)	208 (97%)	6 (3%)	0	100	100
2	H	215/218 (99%)	211 (98%)	4 (2%)	0	100	100
2	J	215/218 (99%)	210 (98%)	5 (2%)	0	100	100
2	L	215/218 (99%)	209 (97%)	6 (3%)	0	100	100
2	N	213/218 (98%)	208 (98%)	5 (2%)	0	100	100
2	P	214/218 (98%)	209 (98%)	5 (2%)	0	100	100
3	Q	7/18 (39%)	7 (100%)	0	0	100	100
3	R	6/18 (33%)	6 (100%)	0	0	100	100
3	S	6/18 (33%)	4 (67%)	2 (33%)	0	100	100
3	T	6/18 (33%)	6 (100%)	0	0	100	100
3	U	6/18 (33%)	6 (100%)	0	0	100	100
3	V	6/18 (33%)	6 (100%)	0	0	100	100
3	W	5/18 (28%)	4 (80%)	1 (20%)	0	100	100
3	X	6/18 (33%)	6 (100%)	0	0	100	100
All	All	3445/3632 (95%)	3363 (98%)	75 (2%)	7 (0%)	43	68

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	102	THR
1	A	102	THR
1	E	102	THR
1	G	102	THR

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Mol	Chain	Res	Type
1	I	102	THR
1	K	102	THR
1	O	102	THR

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	183/185 (99%)	178 (97%)	5 (3%)	39	69
1	C	181/185 (98%)	176 (97%)	5 (3%)	38	68
1	E	185/185 (100%)	183 (99%)	2 (1%)	65	85
1	G	183/185 (99%)	177 (97%)	6 (3%)	33	63
1	I	184/185 (100%)	177 (96%)	7 (4%)	29	58
1	K	184/185 (100%)	181 (98%)	3 (2%)	55	80
1	M	181/185 (98%)	176 (97%)	5 (3%)	38	68
1	O	180/185 (97%)	175 (97%)	5 (3%)	38	68
2	B	187/188 (100%)	182 (97%)	5 (3%)	39	69
2	D	187/188 (100%)	180 (96%)	7 (4%)	30	59
2	F	186/188 (99%)	179 (96%)	7 (4%)	29	58
2	H	187/188 (100%)	183 (98%)	4 (2%)	47	75
2	J	187/188 (100%)	181 (97%)	6 (3%)	34	64
2	L	187/188 (100%)	179 (96%)	8 (4%)	26	54
2	N	185/188 (98%)	182 (98%)	3 (2%)	55	80
2	P	186/188 (99%)	182 (98%)	4 (2%)	45	74
3	Q	3/6 (50%)	3 (100%)	0	100	100
3	R	3/6 (50%)	3 (100%)	0	100	100
3	S	3/6 (50%)	3 (100%)	0	100	100
3	T	3/6 (50%)	3 (100%)	0	100	100
3	U	3/6 (50%)	3 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	V	3/6 (50%)	3 (100%)	0	100	100
3	W	3/6 (50%)	3 (100%)	0	100	100
3	X	3/6 (50%)	3 (100%)	0	100	100
All	All	2977/3032 (98%)	2895 (97%)	82 (3%)	38	68

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	MET
1	A	91	GLU
1	A	174	LEU
1	A	181	SER
1	A	193	SER
2	B	138	CYS
2	B	180	SER
2	B	210	VAL
2	B	212	SER
2	B	215	ARG
1	C	1	GLU
1	C	18	MET
1	C	46	GLU
1	C	181	SER
1	C	193	SER
2	D	36	SER
2	D	138	CYS
2	D	180	SER
2	D	206	THR
2	D	210	VAL
2	D	212	SER
2	D	216	ASN
1	E	18	MET
1	E	193	SER
2	F	138	CYS
2	F	180	SER
2	F	206	THR
2	F	210	VAL
2	F	212	SER
2	F	215	ARG
2	F	216	ASN
1	G	18	MET
1	G	91	GLU

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Mol	Chain	Res	Type
1	G	174	LEU
1	G	181	SER
1	G	193	SER
1	G	208	LYS
2	H	138	CYS
2	H	179	MET
2	H	180	SER
2	H	206	THR
1	I	18	MET
1	I	46	GLU
1	I	91	GLU
1	I	173	LEU
1	I	181	SER
1	I	193	SER
1	I	214	GLU
2	J	147	ASP
2	J	149	ASN
2	J	180	SER
2	J	206	THR
2	J	210	VAL
2	J	212	SER
1	K	18	MET
1	K	181	SER
1	K	193	SER
2	L	138	CYS
2	L	149	ASN
2	L	164	LEU
2	L	180	SER
2	L	206	THR
2	L	210	VAL
2	L	212	SER
2	L	216	ASN
1	M	18	MET
1	M	46	GLU
1	M	174	LEU
1	M	181	SER
1	M	193	SER
2	N	78	ILE
2	N	180	SER
2	N	206	THR
1	O	11	LEU
1	O	18	MET

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Mol	Chain	Res	Type
1	O	46	GLU
1	O	91	GLU
1	O	181	SER
2	P	180	SER
2	P	203	LYS
2	P	206	THR
2	P	216	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	84	GLN
1	A	194	GLN
2	B	128	GLN
2	B	194	ASN
1	C	57	ASN
1	C	84	GLN
1	C	86	ASN
1	C	194	GLN
2	D	128	GLN
2	D	194	ASN
2	D	214	ASN
1	E	39	GLN
1	E	84	GLN
1	E	194	GLN
2	F	42	GLN
2	F	94	GLN
2	F	128	GLN
2	F	194	ASN
2	F	214	ASN
1	G	84	GLN
1	G	167	HIS
2	H	80	HIS
2	H	94	GLN
2	H	128	GLN
2	H	194	ASN
2	H	214	ASN
2	H	216	ASN
2	J	128	GLN
2	J	194	ASN
2	J	214	ASN
2	L	149	ASN

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Mol	Chain	Res	Type
2	L	194	ASN
2	L	214	ASN
2	L	216	ASN
1	M	84	GLN
1	M	108	GLN
1	M	194	GLN
2	N	94	GLN
2	N	128	GLN
2	N	194	ASN
2	N	214	ASN
1	O	84	GLN
1	O	86	ASN
2	P	128	GLN
2	P	149	ASN
2	P	194	ASN
2	P	214	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

17 non-standard protein/DNA/RNA residues are modelled in this entry.

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	$\# Z > 2$
3	CIR	S	11	3	9,10,11	0.62	0	6,11,13	2.61	3 (50%)
3	HYP	R	12	3	7,8,9	0.55	0	5,10,12	1.52	0
3	CIR	U	11	3	9,10,11	0.60	0	6,11,13	2.09	3 (50%)
3	CIR	X	11	3	9,10,11	0.86	0	6,11,13	1.12	0
3	CIR	T	11	3	9,10,11	1.35	2 (22%)	6,11,13	0.83	0
3	CIR	Q	11	3	9,10,11	0.72	0	6,11,13	2.05	3 (50%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	HYP	Q	3	3	7,8,9	0.72	0	5,10,12	1.31	1 (20%)
3	HYP	V	12	3	7,8,9	0.64	0	5,10,12	2.12	1 (20%)
3	HYP	S	12	3	7,8,9	0.55	0	5,10,12	1.76	2 (40%)
3	HYP	U	12	3	7,8,9	0.56	0	5,10,12	1.67	2 (40%)
3	CIR	V	11	3	9,10,11	0.47	0	6,11,13	1.01	0
3	CIR	R	11	3	9,10,11	0.95	0	6,11,13	0.90	0
3	HYP	Q	12	3	7,8,9	1.41	2 (28%)	5,10,12	2.18	2 (40%)
3	HYP	T	12	3	7,8,9	1.53	1 (14%)	5,10,12	0.88	0
3	CIR	W	11	3	9,10,11	0.97	0	6,11,13	1.72	1 (16%)
3	HYP	W	12	3	7,8,9	0.54	0	5,10,12	1.45	0
3	HYP	X	12	3	7,8,9	0.47	0	5,10,12	1.97	2 (40%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	CIR	S	11	3	-	2/8/9/11	-
3	HYP	R	12	3	-	0/0/11/13	0/1/1/1
3	CIR	U	11	3	-	4/8/9/11	-
3	CIR	X	11	3	-	2/8/9/11	-
3	CIR	T	11	3	-	5/8/9/11	-
3	CIR	Q	11	3	-	4/8/9/11	-
3	HYP	Q	3	3	-	0/0/11/13	0/1/1/1
3	HYP	V	12	3	-	0/0/11/13	0/1/1/1
3	HYP	S	12	3	-	0/0/11/13	0/1/1/1
3	HYP	U	12	3	-	0/0/11/13	0/1/1/1
3	CIR	V	11	3	-	1/8/9/11	-
3	CIR	R	11	3	-	2/8/9/11	-
3	HYP	Q	12	3	-	0/0/11/13	0/1/1/1
3	HYP	T	12	3	-	0/0/11/13	0/1/1/1
3	CIR	W	11	3	-	2/8/9/11	-
3	HYP	W	12	3	-	0/0/11/13	0/1/1/1
3	HYP	X	12	3	-	0/0/11/13	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	Q	12	HYP	CB-CA	-2.89	1.48	1.54
3	T	12	HYP	CA-N	-2.80	1.43	1.48
3	T	11	CIR	O7-C7	-2.29	1.20	1.24
3	T	11	CIR	C7-N6	-2.12	1.31	1.34
3	Q	12	HYP	CA-N	-2.01	1.45	1.48

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	V	12	HYP	CB-CG-CD	4.21	107.85	103.16
3	Q	12	HYP	CB-CG-CD	4.19	107.82	103.16
3	S	11	CIR	N8-C7-N6	4.06	123.75	116.64
3	X	12	HYP	CB-CG-CD	3.49	107.04	103.16
3	S	11	CIR	C5-N6-C7	3.42	127.10	122.64
3	U	11	CIR	N8-C7-N6	3.10	122.08	116.64
3	Q	11	CIR	N8-C7-N6	2.97	121.84	116.64
3	U	11	CIR	C5-N6-C7	2.96	126.50	122.64
3	W	11	CIR	O7-C7-N8	-2.92	117.30	123.18
3	S	11	CIR	O7-C7-N8	-2.92	117.30	123.18
3	S	12	HYP	CB-CG-CD	2.82	106.29	103.16
3	Q	11	CIR	C5-N6-C7	2.66	126.10	122.64
3	U	12	HYP	CG-CB-CA	2.57	106.73	103.75
3	U	11	CIR	O7-C7-N8	-2.44	118.27	123.18
3	Q	12	HYP	O-C-CA	-2.23	119.04	124.77
3	U	12	HYP	O-C-CA	-2.23	119.04	124.77
3	Q	3	HYP	OD1-CG-CD	2.17	114.74	110.30
3	S	12	HYP	O-C-CA	-2.14	119.28	124.77
3	X	12	HYP	O-C-CA	-2.06	119.47	124.77
3	Q	11	CIR	O7-C7-N6	-2.05	118.44	121.69

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	Q	11	CIR	C4-C3-CA-C
3	Q	11	CIR	C4-C3-CA-N
3	T	11	CIR	C4-C3-CA-C
3	T	11	CIR	C4-C3-CA-N
3	T	11	CIR	O7-C7-N6-C5
3	T	11	CIR	N8-C7-N6-C5
3	U	11	CIR	C4-C3-CA-C
3	V	11	CIR	C4-C3-CA-C

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Mol	Chain	Res	Type	Atoms
3	T	11	CIR	C3-C4-C5-N6
3	X	11	CIR	C3-C4-C5-N6
3	R	11	CIR	CA-C3-C4-C5
3	Q	11	CIR	O7-C7-N6-C5
3	Q	11	CIR	N8-C7-N6-C5
3	S	11	CIR	O7-C7-N6-C5
3	S	11	CIR	N8-C7-N6-C5
3	U	11	CIR	O7-C7-N6-C5
3	U	11	CIR	N8-C7-N6-C5
3	U	11	CIR	CA-C3-C4-C5
3	X	11	CIR	C4-C3-CA-C
3	W	11	CIR	C3-C4-C5-N6
3	R	11	CIR	C4-C3-CA-N
3	W	11	CIR	C4-C3-CA-N

There are no ring outliers.

No monomer is involved in short contacts.

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	214/218 (98%)	0.37	5 (2%) 61 58	30, 56, 88, 102	0
1	C	212/218 (97%)	0.43	4 (1%) 66 63	32, 61, 96, 115	0
1	E	218/218 (100%)	0.62	8 (3%) 45 41	47, 70, 104, 121	0
1	G	215/218 (98%)	0.14	2 (0%) 81 80	36, 53, 76, 91	0
1	I	216/218 (99%)	0.84	18 (8%) 17 14	45, 78, 101, 117	0
1	K	215/218 (98%)	1.01	28 (13%) 7 6	41, 78, 106, 124	0
1	M	212/218 (97%)	0.84	16 (7%) 20 17	40, 83, 125, 139	0
1	O	211/218 (96%)	1.52	42 (19%) 3 2	77, 119, 163, 192	0
2	B	217/218 (99%)	0.37	1 (0%) 87 86	33, 60, 94, 113	0
2	D	217/218 (99%)	0.29	1 (0%) 87 86	27, 51, 93, 109	0
2	F	216/218 (99%)	0.42	3 (1%) 73 72	42, 64, 103, 117	0
2	H	217/218 (99%)	0.18	0 100 100	33, 53, 79, 104	0
2	J	217/218 (99%)	0.53	5 (2%) 61 58	31, 66, 111, 126	0
2	L	217/218 (99%)	0.65	7 (3%) 50 46	32, 68, 107, 127	0
2	N	215/218 (98%)	1.00	17 (7%) 18 16	52, 93, 125, 138	0
2	P	216/218 (99%)	1.17	23 (10%) 11 9	55, 98, 141, 165	0
3	Q	8/18 (44%)	0.25	1 (12%) 8 7	41, 44, 59, 71	0
3	R	8/18 (44%)	0.89	1 (12%) 8 7	58, 63, 66, 79	0
3	S	8/18 (44%)	0.59	0 100 100	67, 71, 78, 89	0
3	T	8/18 (44%)	1.00	1 (12%) 8 7	61, 66, 78, 85	0
3	U	8/18 (44%)	1.46	3 (37%) 1 0	63, 70, 73, 80	0
3	V	8/18 (44%)	0.70	1 (12%) 8 7	62, 69, 76, 92	0
3	W	7/18 (38%)	0.15	0 100 100	60, 62, 70, 75	0
3	X	8/18 (44%)	1.72	3 (37%) 1 0	94, 101, 104, 107	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
All	All	3508/3632 (96%)	0.65	190 (5%)	31	27	27, 69, 121, 192	0

All (190) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	174	LEU	4.3
1	O	122	PRO	4.3
1	O	113	THR	4.1
1	O	117	ALA	4.0
1	O	132	GLY	3.8
3	U	4	GLY	3.6
1	O	112	LEU	3.6
3	R	13	GLY	3.6
1	I	172	LEU	3.5
1	K	135	THR	3.5
1	A	217	GLY	3.4
3	X	4	GLY	3.4
1	O	148	TYR	3.4
1	K	16	GLY	3.4
1	M	215	PRO	3.3
1	C	174	LEU	3.3
1	O	114	VAL	3.3
2	P	137	VAL	3.2
1	M	176	GLY	3.2
2	J	156	GLY	3.2
1	K	174	LEU	3.2
1	O	142	GLY	3.1
1	K	215	PRO	3.1
1	O	215	PRO	3.0
1	O	36	TRP	3.0
1	O	123	SER	3.0
1	O	144	LEU	3.0
1	I	189	ASN	3.0
1	O	109	GLY	3.0
2	N	105	GLY	3.0
2	P	152	TRP	2.9
1	M	73	SER	2.9
1	O	134	THR	2.9
1	O	125	TYR	2.9
2	P	131	SER	2.9
1	K	70	PHE	2.9
2	P	34	ILE	2.9

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Mol	Chain	Res	Type	RSRZ
2	J	157	SER	2.8
1	I	215	PRO	2.8
1	K	131	CYS	2.8
2	P	32	TYR	2.8
1	M	164	SER	2.8
2	P	194	ASN	2.8
1	C	215	PRO	2.8
1	O	88	LEU	2.8
1	O	72	ILE	2.7
1	O	159	SER	2.7
1	O	130	VAL	2.7
2	N	213	PHE	2.7
1	O	108	GLN	2.7
2	N	152	TRP	2.7
2	P	69	SER	2.7
1	O	20	LEU	2.7
1	A	175	SER	2.7
1	K	116	SER	2.7
1	K	209	VAL	2.7
1	I	210	ASP	2.7
2	L	59	GLY	2.6
1	O	129	PRO	2.6
1	K	15	GLY	2.6
1	K	201	ALA	2.6
1	O	27	PHE	2.6
1	K	195	THR	2.5
2	P	183	LEU	2.5
1	M	210	ASP	2.5
2	N	209	ILE	2.5
1	I	135	THR	2.5
1	K	155	LEU	2.5
1	O	141	LEU	2.5
1	O	178	TYR	2.5
1	I	209	VAL	2.5
1	O	85	MET	2.5
1	I	131	CYS	2.5
1	O	118	LYS	2.5
1	M	11	LEU	2.5
1	I	218	PRO	2.5
1	K	18	MET	2.5
2	P	37	MET	2.5
1	M	174	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
2	L	135	SER	2.4
1	M	200	VAL	2.4
2	N	79	ILE	2.4
2	P	150	VAL	2.4
2	F	156	GLY	2.4
2	N	210	VAL	2.4
2	N	64	ALA	2.4
1	E	155	LEU	2.3
1	I	4	LEU	2.3
1	O	101	LEU	2.3
2	N	129	LEU	2.3
2	F	216	ASN	2.3
1	K	115	SER	2.3
1	E	68	GLY	2.3
2	N	11	LEU	2.3
2	B	184	THR	2.3
1	G	3	LYS	2.3
1	M	79	SER	2.3
2	N	138	CYS	2.3
1	M	186	VAL	2.3
1	O	200	VAL	2.3
2	P	136	VAL	2.3
1	K	141	LEU	2.3
2	N	122	PHE	2.3
3	U	5	ALA	2.3
1	O	207	THR	2.3
1	I	175	SER	2.3
1	K	150	PRO	2.3
3	X	10	GLY	2.3
1	K	200	VAL	2.3
1	O	98	CYS	2.3
1	O	131	CYS	2.3
1	O	145	VAL	2.3
2	P	139	PHE	2.3
2	P	79	ILE	2.3
3	X	5	ALA	2.3
1	E	162	LEU	2.3
1	I	121	ALA	2.2
1	M	117	ALA	2.2
2	L	121	ILE	2.2
2	P	25	ALA	2.2
2	N	190	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	177	LEU	2.2
1	E	166	VAL	2.2
1	C	196	ILE	2.2
1	I	213	ILE	2.2
1	O	59	ALA	2.2
2	P	154	ILE	2.2
2	P	186	THR	2.2
2	L	32	TYR	2.2
2	J	161	ASN	2.2
1	K	188	SER	2.2
1	E	209	VAL	2.2
1	K	114	VAL	2.2
2	N	148	ILE	2.2
2	P	199	GLU	2.2
1	K	122	PRO	2.2
1	I	67	LYS	2.2
1	O	83	LEU	2.2
3	U	13	GLY	2.2
1	G	174	LEU	2.2
1	M	127	LEU	2.2
2	P	147	ASP	2.2
1	I	188	SER	2.2
2	L	26	SER	2.2
2	N	21	ILE	2.2
1	K	14	PRO	2.1
1	K	218	PRO	2.1
1	I	174	LEU	2.1
2	L	185	LEU	2.1
1	E	160	GLY	2.1
1	E	201	ALA	2.1
1	O	180	LEU	2.1
3	T	13	GLY	2.1
1	I	206	SER	2.1
2	N	139	PHE	2.1
2	P	191	GLU	2.1
2	J	24	ARG	2.1
1	E	134	THR	2.1
1	K	113	THR	2.1
1	A	215	PRO	2.1
1	K	11	LEU	2.1
2	D	164	LEU	2.1
2	N	77	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
3	V	4	GLY	2.1
1	M	130	VAL	2.1
2	P	66	PHE	2.1
1	O	175	SER	2.1
2	P	135	SER	2.1
2	P	195	SER	2.1
1	M	211	LYS	2.1
1	K	179	THR	2.1
1	O	179	THR	2.1
2	P	168	THR	2.1
1	K	112	LEU	2.1
1	M	144	LEU	2.1
1	I	124	VAL	2.1
1	K	2	VAL	2.1
1	M	209	VAL	2.1
1	C	131	CYS	2.0
1	O	96	TYR	2.0
2	F	210	VAL	2.0
2	J	196	TYR	2.0
1	K	204	ALA	2.0
2	L	64	ALA	2.0
1	I	17	SER	2.0
1	O	156	THR	2.0
1	O	45	LEU	2.0
1	O	160	GLY	2.0
3	Q	4	GLY	2.0
2	N	192	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
3	HYP	Q	3	8/9	0.73	0.15	77,82,88,90	0
3	CIR	X	11	11/12	0.74	0.12	92,93,95,99	0
3	CIR	V	11	11/12	0.77	0.15	70,72,76,78	0
3	HYP	V	12	8/9	0.80	0.15	84,89,93,93	0
3	HYP	W	12	8/9	0.81	0.11	74,76,79,80	0
3	CIR	T	11	11/12	0.81	0.13	65,69,74,77	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	HYP	X	12	8/9	0.82	0.12	95,96,98,99	0
3	HYP	U	12	8/9	0.84	0.16	82,83,83,83	0
3	HYP	T	12	8/9	0.85	0.11	73,76,78,79	0
3	HYP	R	12	8/9	0.86	0.13	71,73,75,75	0
3	HYP	S	12	8/9	0.86	0.15	83,86,87,87	0
3	CIR	R	11	11/12	0.88	0.12	63,64,69,70	0
3	CIR	W	11	11/12	0.89	0.12	71,75,91,97	0
3	HYP	Q	12	8/9	0.89	0.09	49,50,52,52	0
3	CIR	U	11	11/12	0.90	0.12	77,80,83,83	0
3	CIR	S	11	11/12	0.91	0.11	76,80,87,88	0
3	CIR	Q	11	11/12	0.91	0.10	44,47,60,65	0

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.