



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:19 AM UTC

PDB ID : 3ZHA / pdb_00003zha
Title : Molecular basis for the action of the collagen-specific chaperone Hsp47 SER-PINH1 and its structure-specific client recognition.
Authors : Widmer, C.; Gebauer, J.M.; Baumann, U.
Deposited on : 2012-12-20
Resolution : 2.55 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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

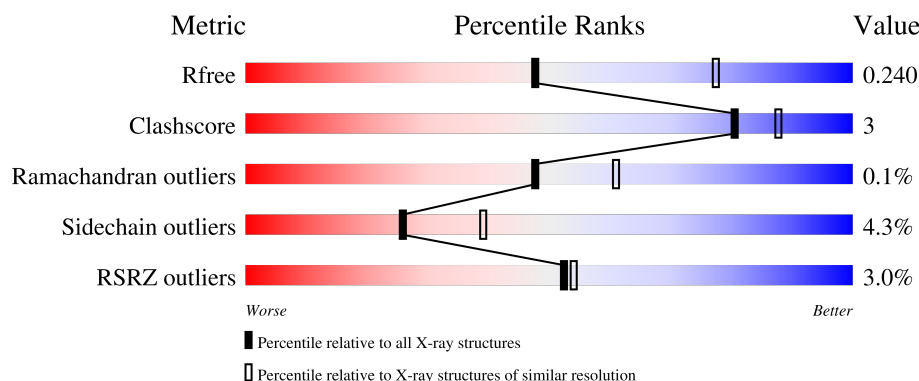
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	392	
1	B	392	
1	C	392	
1	D	392	
1	K	392	

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Mol	Chain	Length	Quality of chain
1	L	392	
1	P	392	
1	Q	392	
2	E	19	
2	F	19	
2	G	19	
2	H	19	
2	I	19	
2	J	19	
2	M	19	
2	N	19	
2	O	19	
2	R	19	
2	S	19	
2	T	19	

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
3	SIN	A	1421	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 24911 atoms, of which 8 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 HSP47.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	380	Total	C	N	O	S	0	0	0
			2981	1898	517	552	14			
1	B	364	Total	C	N	O	S	0	0	0
			2855	1820	496	526	13			
1	C	376	Total	C	N	O	S	0	0	0
			2961	1884	516	548	13			
1	D	368	Total	C	N	O	S	0	0	0
			2894	1840	506	535	13			
1	K	374	Total	C	N	O	S	0	0	0
			2955	1879	515	547	14			
1	L	369	Total	C	N	O	S	0	0	0
			2902	1847	505	537	13			
1	P	373	Total	C	N	O	S	0	0	0
			2938	1868	511	546	13			
1	Q	365	Total	C	N	O	S	0	0	0
			2872	1828	500	531	13			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	35	MET	-	expression tag	UNP C7C419
A	419	LEU	-	expression tag	UNP C7C419
A	420	GLU	-	expression tag	UNP C7C419
A	421	HIS	-	expression tag	UNP C7C419
A	422	HIS	-	expression tag	UNP C7C419
A	423	HIS	-	expression tag	UNP C7C419
A	424	HIS	-	expression tag	UNP C7C419
A	425	HIS	-	expression tag	UNP C7C419
A	426	HIS	-	expression tag	UNP C7C419
B	35	MET	-	expression tag	UNP C7C419
B	419	LEU	-	expression tag	UNP C7C419
B	420	GLU	-	expression tag	UNP C7C419
B	421	HIS	-	expression tag	UNP C7C419

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Chain	Residue	Modelled	Actual	Comment	Reference
B	422	HIS	-	expression tag	UNP C7C419
B	423	HIS	-	expression tag	UNP C7C419
B	424	HIS	-	expression tag	UNP C7C419
B	425	HIS	-	expression tag	UNP C7C419
B	426	HIS	-	expression tag	UNP C7C419
C	35	MET	-	expression tag	UNP C7C419
C	419	LEU	-	expression tag	UNP C7C419
C	420	GLU	-	expression tag	UNP C7C419
C	421	HIS	-	expression tag	UNP C7C419
C	422	HIS	-	expression tag	UNP C7C419
C	423	HIS	-	expression tag	UNP C7C419
C	424	HIS	-	expression tag	UNP C7C419
C	425	HIS	-	expression tag	UNP C7C419
C	426	HIS	-	expression tag	UNP C7C419
D	35	MET	-	expression tag	UNP C7C419
D	419	LEU	-	expression tag	UNP C7C419
D	420	GLU	-	expression tag	UNP C7C419
D	421	HIS	-	expression tag	UNP C7C419
D	422	HIS	-	expression tag	UNP C7C419
D	423	HIS	-	expression tag	UNP C7C419
D	424	HIS	-	expression tag	UNP C7C419
D	425	HIS	-	expression tag	UNP C7C419
D	426	HIS	-	expression tag	UNP C7C419
K	35	MET	-	expression tag	UNP C7C419
K	419	LEU	-	expression tag	UNP C7C419
K	420	GLU	-	expression tag	UNP C7C419
K	421	HIS	-	expression tag	UNP C7C419
K	422	HIS	-	expression tag	UNP C7C419
K	423	HIS	-	expression tag	UNP C7C419
K	424	HIS	-	expression tag	UNP C7C419
K	425	HIS	-	expression tag	UNP C7C419
K	426	HIS	-	expression tag	UNP C7C419
L	35	MET	-	expression tag	UNP C7C419
L	419	LEU	-	expression tag	UNP C7C419
L	420	GLU	-	expression tag	UNP C7C419
L	421	HIS	-	expression tag	UNP C7C419
L	422	HIS	-	expression tag	UNP C7C419
L	423	HIS	-	expression tag	UNP C7C419
L	424	HIS	-	expression tag	UNP C7C419
L	425	HIS	-	expression tag	UNP C7C419
L	426	HIS	-	expression tag	UNP C7C419
P	35	MET	-	expression tag	UNP C7C419

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Chain	Residue	Modelled	Actual	Comment	Reference
P	419	LEU	-	expression tag	UNP C7C419
P	420	GLU	-	expression tag	UNP C7C419
P	421	HIS	-	expression tag	UNP C7C419
P	422	HIS	-	expression tag	UNP C7C419
P	423	HIS	-	expression tag	UNP C7C419
P	424	HIS	-	expression tag	UNP C7C419
P	425	HIS	-	expression tag	UNP C7C419
P	426	HIS	-	expression tag	UNP C7C419
Q	35	MET	-	expression tag	UNP C7C419
Q	419	LEU	-	expression tag	UNP C7C419
Q	420	GLU	-	expression tag	UNP C7C419
Q	421	HIS	-	expression tag	UNP C7C419
Q	422	HIS	-	expression tag	UNP C7C419
Q	423	HIS	-	expression tag	UNP C7C419
Q	424	HIS	-	expression tag	UNP C7C419
Q	425	HIS	-	expression tag	UNP C7C419
Q	426	HIS	-	expression tag	UNP C7C419

- Molecule 2 is a protein called COLLAGEN MODEL PEPTIDE 18-T8R11.

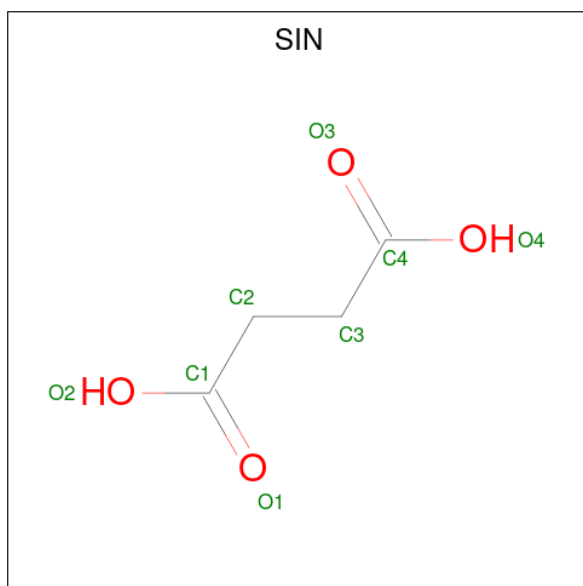
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	18	Total	C	N	O	0	0	0
			111	72	20	19			
2	F	17	Total	C	N	O	0	0	0
			104	67	19	18			
2	G	16	Total	C	N	O	0	0	0
			97	62	18	17			
2	H	18	Total	C	N	O	0	0	0
			111	72	20	19			
2	I	17	Total	C	N	O	0	0	0
			104	67	19	18			
2	J	16	Total	C	N	O	0	0	0
			97	62	18	17			
2	M	18	Total	C	N	O	0	0	0
			111	72	20	19			
2	N	18	Total	C	N	O	0	0	0
			111	72	20	19			
2	O	18	Total	C	N	O	0	0	0
			111	72	20	19			
2	R	18	Total	C	N	O	0	0	0
			111	72	20	19			
2	S	18	Total	C	N	O	0	0	0
			111	72	20	19			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	T	17	Total	C	N	O	0	0	0
			104	67	19	18			

- Molecule 3 is SUCCINIC ACID (CCD ID: SIN) (formula: C₄H₆O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O		0	0
			8	4	4			
3	C	1	Total	C	O		0	0
			8	4	4			
3	K	1	Total	C	O		0	0
			8	4	4			
3	P	1	Total	C	H	O	0	0
			12	4	4	4		
3	Q	1	Total	C	H	O	0	0
			12	4	4	4		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	23	Total	O	0	0
			23	23		
4	B	21	Total	O	0	0
			21	21		
4	C	33	Total	O	0	0
			33	33		

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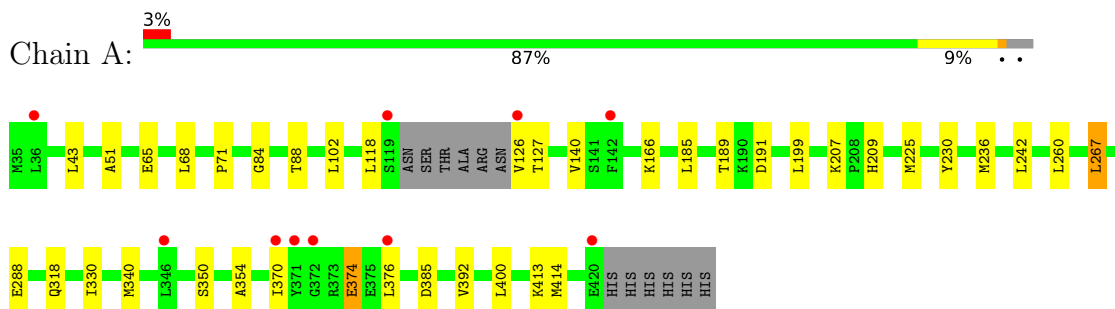
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	37	Total 37	O 37	0	0
4	E	2	Total 2	O 2	0	0
4	F	1	Total 1	O 1	0	0
4	I	2	Total 2	O 2	0	0
4	K	24	Total 24	O 24	0	0
4	L	27	Total 27	O 27	0	0
4	N	3	Total 3	O 3	0	0
4	P	18	Total 18	O 18	0	0
4	Q	27	Total 27	O 27	0	0
4	R	1	Total 1	O 1	0	0
4	S	2	Total 2	O 2	0	0
4	T	1	Total 1	O 1	0	0

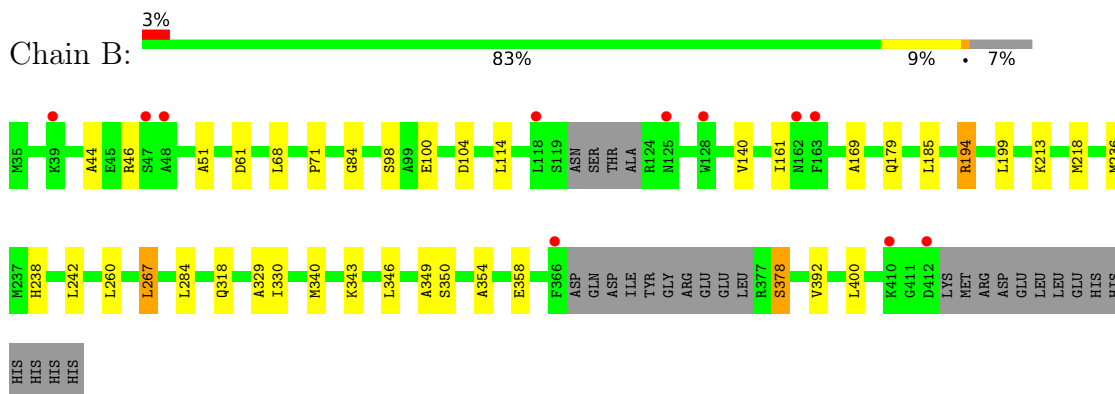
3 Residue-property plots [i](#)

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.

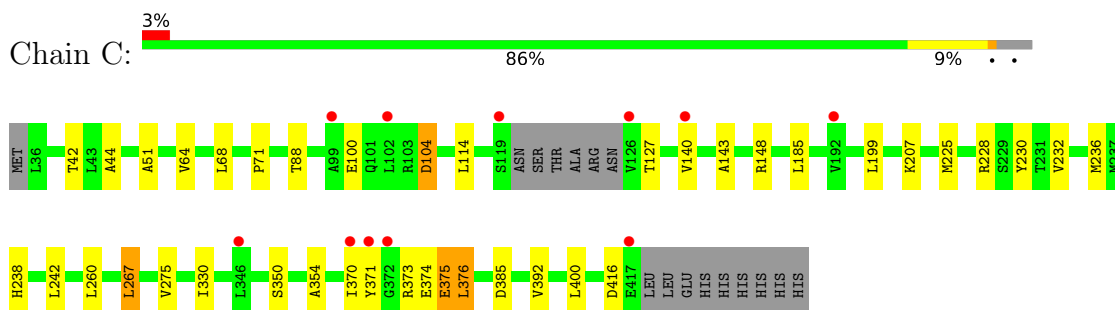
- Molecule 1: HSP47



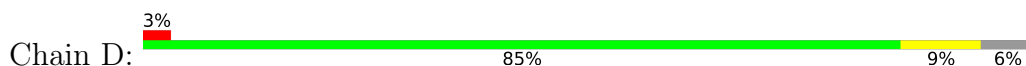
- Molecule 1: HSP47

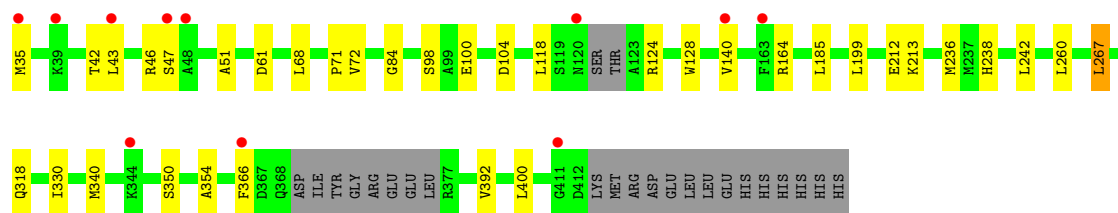


- Molecule 1: HSP47

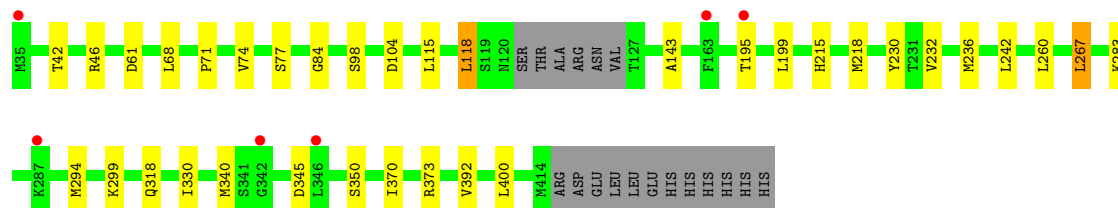
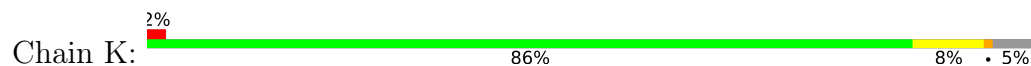


- Molecule 1: HSP47

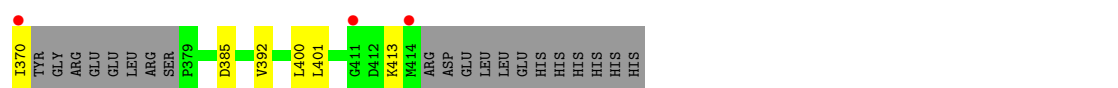
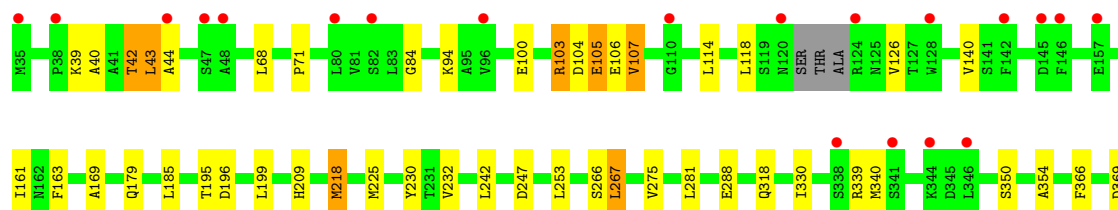
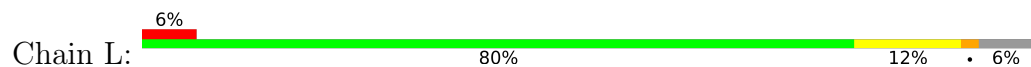




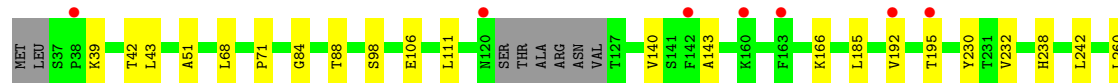
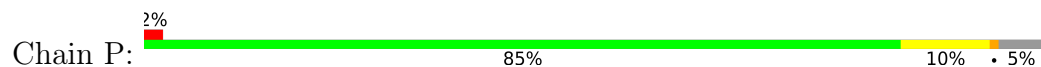
• Molecule 1: HSP47



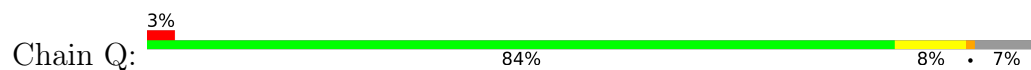
• Molecule 1: HSP47

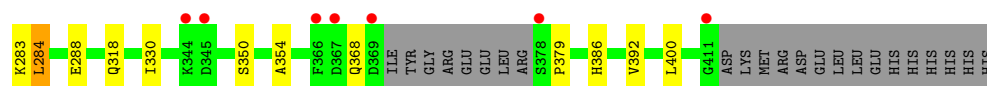


• Molecule 1: HSP47



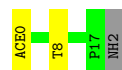
• Molecule 1: HSP47





- Molecule 2: COLLAGEN MODEL PEPTIDE 18-T8R11

Chain E: 84% 11% 5%



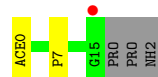
- Molecule 2: COLLAGEN MODEL PEPTIDE 18-T8R11

Chain F: 79% 11% 11%



- Molecule 2: COLLAGEN MODEL PEPTIDE 18-T8R11

Chain G: 5% 74% 11% 16%



- Molecule 2: COLLAGEN MODEL PEPTIDE 18-T8R11

Chain H: 79% 16% 5%



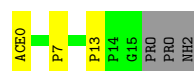
- Molecule 2: COLLAGEN MODEL PEPTIDE 18-T8R11

Chain I: 74% 16% 11%



- Molecule 2: COLLAGEN MODEL PEPTIDE 18-T8R11

Chain J: 68% 16% 16%



- Molecule 2: COLLAGEN MODEL PEPTIDE 18-T8R11

Chain M: 89% 5% 5%



- Molecule 2: COLLAGEN MODEL PEPTIDE 18-T8R11

Chain N:  89% 5% 5%



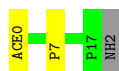
- Molecule 2: COLLAGEN MODEL PEPTIDE 18-T8R11

Chain O:  89% 5% 5%

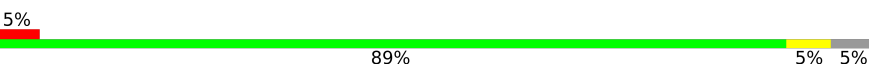


- Molecule 2: COLLAGEN MODEL PEPTIDE 18-T8R11

Chain R:  84% 11% 5%



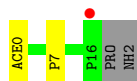
- Molecule 2: COLLAGEN MODEL PEPTIDE 18-T8R11

Chain S:  5% 89% 5% 5%



- Molecule 2: COLLAGEN MODEL PEPTIDE 18-T8R11

Chain T:  5% 79% 11% 11%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	102.71Å 104.91Å 171.84Å 90.00° 103.71° 90.00°	Depositor
Resolution (Å)	49.16 – 2.55 49.16 – 2.55	Depositor EDS
% Data completeness (in resolution range)	98.0 (49.16-2.55) 98.1 (49.16-2.55)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 2.54Å)	Xtrriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.195 , 0.218 0.218 , 0.240	Depositor DCC
R_{free} test set	1780 reflections (1.57%)	wwPDB-VP
Wilson B-factor (Å ²)	40.1	Xtrriage
Anisotropy	0.190	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 63.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.37$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	24911	wwPDB-VP
Average B, all atoms (Å ²)	58.0	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 47.19 % 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 1.0277e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: ACE, SIN

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	0.72	1/3041 (0.0%)	1.29	0/4101
1	B	0.73	1/2913 (0.0%)	1.31	4/3928 (0.1%)
1	C	0.73	1/3021 (0.0%)	1.32	10/4073 (0.2%)
1	D	0.74	1/2952 (0.0%)	1.31	5/3980 (0.1%)
1	K	0.73	1/3015 (0.0%)	1.29	3/4063 (0.1%)
1	L	0.75	1/2960 (0.0%)	1.32	8/3990 (0.2%)
1	P	0.72	0/2998	1.31	6/4042 (0.1%)
1	Q	0.76	1/2930 (0.0%)	1.30	5/3951 (0.1%)
2	E	1.17	1/118 (0.8%)	0.85	0/168
2	F	1.11	1/110 (0.9%)	0.80	0/156
2	G	1.12	1/102 (1.0%)	0.85	0/144
2	H	1.15	1/118 (0.8%)	0.90	0/168
2	I	1.09	1/110 (0.9%)	0.74	0/156
2	J	1.14	1/102 (1.0%)	0.88	0/144
2	M	1.08	1/118 (0.8%)	0.88	0/168
2	N	1.19	1/118 (0.8%)	0.81	0/168
2	O	1.14	1/118 (0.8%)	0.85	0/168
2	R	1.13	1/118 (0.8%)	0.80	0/168
2	S	1.18	1/118 (0.8%)	0.74	0/168
2	T	1.18	1/110 (0.9%)	0.76	0/156
All	All	0.76	19/25190 (0.1%)	1.28	41/34060 (0.1%)

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	N	0	ACE	C-N	-9.35	1.34	1.43
2	T	0	ACE	C-N	-9.24	1.34	1.43
2	O	0	ACE	C-N	-9.03	1.34	1.43
2	S	0	ACE	C-N	-9.03	1.34	1.43
2	F	0	ACE	C-N	-8.85	1.34	1.43
2	I	0	ACE	C-N	-8.73	1.34	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	0	ACE	C-N	-8.73	1.34	1.43
2	R	0	ACE	C-N	-8.72	1.34	1.43
2	E	0	ACE	C-N	-8.69	1.34	1.43
2	G	0	ACE	C-N	-8.56	1.34	1.43
2	J	0	ACE	C-N	-8.54	1.34	1.43
2	M	0	ACE	C-N	-8.06	1.35	1.43
1	L	218	MET	SD-CE	-7.38	1.61	1.79
1	D	236	MET	SD-CE	-6.99	1.62	1.79
1	B	236	MET	SD-CE	-6.74	1.62	1.79
1	A	236	MET	SD-CE	-5.90	1.64	1.79
1	Q	386	HIS	CA-C	5.87	1.57	1.52
1	K	236	MET	SD-CE	-5.60	1.65	1.79
1	C	236	MET	SD-CE	-5.22	1.66	1.79

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	416	ASP	N-CA-C	7.84	121.25	109.95
1	P	343	LYS	CA-C-N	7.15	129.86	120.28
1	P	343	LYS	C-N-CA	7.15	129.86	120.28
1	C	104	ASP	CA-C-N	6.46	129.77	120.79
1	C	104	ASP	C-N-CA	6.46	129.77	120.79
1	C	376	LEU	N-CA-C	-6.37	104.81	112.59
1	Q	275	VAL	CB-CA-C	6.30	118.03	111.23
1	Q	179	GLN	CA-C-N	5.97	128.28	120.28
1	Q	179	GLN	C-N-CA	5.97	128.28	120.28
1	L	196	ASP	N-CA-C	5.83	118.58	111.82
1	B	329	ALA	CA-C-N	5.65	129.91	120.64
1	B	329	ALA	C-N-CA	5.65	129.91	120.64
1	C	376	LEU	CA-C-N	5.58	127.70	120.44
1	C	376	LEU	C-N-CA	5.58	127.70	120.44
1	B	358	GLU	N-CA-C	5.47	117.64	108.73
1	L	105	GLU	CA-C-N	5.46	128.14	120.28
1	L	105	GLU	C-N-CA	5.46	128.14	120.28
1	C	373	ARG	N-CA-C	-5.42	104.00	111.81
1	D	61	ASP	CA-CB-CG	5.42	118.02	112.60
1	L	140	VAL	N-CA-C	5.34	116.47	109.58
1	L	179	GLN	CA-C-N	5.34	127.44	120.28
1	L	179	GLN	C-N-CA	5.34	127.44	120.28
1	P	143	ALA	CA-C-N	5.31	127.65	120.38
1	P	143	ALA	C-N-CA	5.31	127.65	120.38
1	K	143	ALA	CA-C-N	5.27	127.60	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	143	ALA	C-N-CA	5.27	127.60	120.38
1	P	373	ARG	CA-C-N	5.27	128.08	120.38
1	P	373	ARG	C-N-CA	5.27	128.08	120.38
1	D	100	GLU	CA-C-N	5.26	127.86	120.28
1	D	100	GLU	C-N-CA	5.26	127.86	120.28
1	D	35	MET	CA-C-N	5.23	128.13	120.71
1	D	35	MET	C-N-CA	5.23	128.13	120.71
1	L	107	VAL	N-CA-C	-5.22	105.62	110.53
1	Q	100	GLU	CA-C-N	5.09	128.05	120.31
1	Q	100	GLU	C-N-CA	5.09	128.05	120.31
1	K	61	ASP	CA-CB-CG	5.09	117.69	112.60
1	C	143	ALA	CA-C-N	5.06	127.32	120.38
1	C	143	ALA	C-N-CA	5.06	127.32	120.38
1	B	61	ASP	CA-CB-CG	5.05	117.65	112.60
1	L	40	ALA	N-CA-C	-5.04	105.87	111.36
1	C	275	VAL	N-CA-CB	5.03	116.84	111.46

There are no chirality outliers.

There are no planarity outliers.

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	2981	0	2995	17	0
1	B	2855	0	2880	16	0
1	C	2961	0	2982	18	0
1	D	2894	0	2918	14	0
1	K	2955	0	2982	12	0
1	L	2902	0	2926	25	0
1	P	2938	0	2953	16	0
1	Q	2872	0	2895	12	0
2	E	111	0	108	1	0
2	F	104	0	101	1	0
2	G	97	0	94	1	0
2	H	111	0	108	3	0
2	I	104	0	101	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	J	97	0	94	2	0
2	M	111	0	108	0	0
2	N	111	0	108	0	0
2	O	111	0	108	0	0
2	R	111	0	108	2	0
2	S	111	0	108	0	0
2	T	104	0	101	1	0
3	A	8	0	4	1	0
3	C	8	0	4	1	0
3	K	8	0	4	0	0
3	P	8	4	4	0	0
3	Q	8	4	4	0	0
4	A	23	0	0	0	0
4	B	21	0	0	1	0
4	C	33	0	0	0	0
4	D	37	0	0	0	0
4	E	2	0	0	0	0
4	F	1	0	0	0	0
4	I	2	0	0	0	0
4	K	24	0	0	0	0
4	L	27	0	0	1	0
4	N	3	0	0	0	0
4	P	18	0	0	0	0
4	Q	27	0	0	0	0
4	R	1	0	0	0	0
4	S	2	0	0	0	0
4	T	1	0	0	0	0
All	All	24903	8	24798	130	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (130) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:218:MET:HE3	4:B:2008:HOH:O	1.50	1.09
1:C:127:THR:HG23	1:C:207:LYS:HB3	1.48	0.94
1:A:414:MET:HE1	1:B:378:SER:HB3	1.60	0.84
1:L:218:MET:HE3	4:L:2013:HOH:O	1.80	0.80
1:A:126:VAL:HG12	1:A:209:HIS:H	1.46	0.79
1:C:374:GLU:HG3	1:C:376:LEU:HB2	1.63	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:42:THR:HG22	1:D:46:ARG:HH12	1.53	0.73
1:C:185:LEU:HD21	1:C:354:ALA:HB1	1.78	0.66
1:L:104:ASP:O	1:L:106:GLU:N	2.33	0.62
1:B:194:ARG:HG3	1:B:349:ALA:HB1	1.82	0.61
1:L:163:PHE:HB2	1:L:195:THR:HG23	1.82	0.61
1:D:71:PRO:HG2	1:D:400:LEU:HB3	1.83	0.61
1:D:42:THR:HG22	1:D:46:ARG:NH1	2.17	0.59
1:B:71:PRO:HG2	1:B:400:LEU:HB3	1.84	0.59
1:P:230:TYR:CE1	1:P:232:VAL:HG23	2.36	0.59
1:P:230:TYR:HE1	1:P:232:VAL:HG23	1.66	0.59
1:A:230:TYR:HE1	1:A:413:LYS:HG3	1.69	0.58
1:A:414:MET:HE1	1:B:378:SER:CB	2.34	0.58
1:K:230:TYR:CE1	1:K:232:VAL:HG23	2.38	0.58
1:A:65:GLU:H	3:A:1421:SIN:H22	1.68	0.57
1:A:185:LEU:HD21	1:A:354:ALA:HB1	1.87	0.57
1:C:230:TYR:CE1	1:C:232:VAL:HG23	2.39	0.57
1:A:230:TYR:CE1	1:A:413:LYS:HG3	2.40	0.57
1:Q:71:PRO:HG2	1:Q:400:LEU:HB3	1.87	0.57
1:C:230:TYR:HE1	1:C:232:VAL:HG23	1.71	0.56
1:K:230:TYR:HE1	1:K:232:VAL:HG23	1.70	0.56
1:L:39:LYS:HA	1:L:42:THR:HG22	1.87	0.56
1:P:71:PRO:HG2	1:P:400:LEU:HB3	1.88	0.56
1:L:366:PHE:CZ	1:L:370:ILE:HG12	2.40	0.56
1:L:366:PHE:HZ	1:L:370:ILE:HG12	1.71	0.56
1:A:230:TYR:HE1	1:A:413:LYS:CG	2.19	0.55
1:P:43:LEU:HD23	1:P:111:LEU:HG	1.88	0.54
1:C:71:PRO:HG2	1:C:400:LEU:HB3	1.89	0.54
1:Q:185:LEU:HD21	1:Q:354:ALA:HB1	1.90	0.54
1:L:230:TYR:CE1	1:L:413:LYS:HG3	2.43	0.53
1:K:370:ILE:HA	1:K:373:ARG:HH11	1.73	0.53
1:C:371:TYR:HA	1:C:374:GLU:HB2	1.90	0.53
1:B:318:GLN:HG2	1:B:330:ILE:HG13	1.91	0.53
1:B:185:LEU:HD21	1:B:354:ALA:HB1	1.89	0.53
1:D:185:LEU:HD21	1:D:354:ALA:HB1	1.89	0.52
1:A:71:PRO:HG2	1:A:400:LEU:HB3	1.91	0.52
1:D:212:GLU:HG2	1:D:366:PHE:HB2	1.92	0.52
1:Q:267:LEU:HD12	1:Q:392:VAL:HG22	1.92	0.52
1:K:267:LEU:HD12	1:K:392:VAL:HG22	1.91	0.52
1:B:84:GLY:HA3	1:B:340:MET:HE3	1.90	0.51
1:Q:199:LEU:HD22	1:Q:350:SER:HB2	1.92	0.51
1:L:185:LEU:HD21	1:L:354:ALA:HB1	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:84:GLY:HA3	1:K:340:MET:HE3	1.91	0.51
1:L:230:TYR:HE1	1:L:413:LYS:HG3	1.76	0.50
1:C:375:GLU:OE1	1:C:375:GLU:HA	2.10	0.50
1:B:44:ALA:HB2	1:B:114:LEU:HD21	1.94	0.50
1:L:71:PRO:HG2	1:L:400:LEU:HB3	1.94	0.50
1:L:230:TYR:HE1	1:L:413:LYS:CG	2.24	0.50
1:P:84:GLY:HA3	1:P:340:MET:HE3	1.92	0.50
1:D:267:LEU:HD12	1:D:392:VAL:HG22	1.94	0.49
1:P:230:TYR:HE1	1:P:232:VAL:CG2	2.24	0.49
1:P:414:MET:HE1	1:Q:379:PRO:HD2	1.93	0.49
1:P:166:LYS:HE3	1:P:192:VAL:HG12	1.94	0.49
1:D:51:ALA:HB2	1:D:71:PRO:HG3	1.94	0.49
1:P:343:LYS:HE3	1:P:345:ASP:HB3	1.95	0.49
1:L:199:LEU:HD22	1:L:350:SER:HB2	1.95	0.49
1:L:267:LEU:HD12	1:L:392:VAL:HG22	1.94	0.49
1:A:84:GLY:HA3	1:A:340:MET:HE3	1.95	0.49
1:D:72:VAL:HG21	1:D:118:LEU:HD22	1.94	0.48
1:D:124:ARG:HA	1:D:124:ARG:HE	1.78	0.48
1:A:267:LEU:HD12	1:A:392:VAL:HG22	1.94	0.48
1:P:267:LEU:HD12	1:P:392:VAL:HG22	1.96	0.47
1:B:46:ARG:HD2	1:B:98:SER:HB3	1.96	0.47
1:C:64:VAL:HA	3:C:1418:SIN:H31	1.96	0.47
1:Q:254:GLN:HG2	1:Q:273:HIS:CD2	2.49	0.47
2:E:8:THR:HA	2:F:7:PRO:HD2	1.95	0.47
1:K:46:ARG:HD2	1:K:98:SER:HB3	1.95	0.47
1:C:238:HIS:CD2	2:H:7:PRO:HB3	2.49	0.47
1:L:161:ILE:HD11	1:L:169:ALA:HA	1.97	0.47
1:B:267:LEU:HD12	1:B:392:VAL:HG22	1.97	0.47
1:P:185:LEU:HD21	1:P:354:ALA:HB1	1.96	0.47
1:D:199:LEU:HD22	1:D:350:SER:HB2	1.97	0.47
1:K:199:LEU:HD22	1:K:350:SER:HB2	1.97	0.47
1:C:230:TYR:HE1	1:C:232:VAL:CG2	2.28	0.46
1:C:238:HIS:CG	2:H:7:PRO:HB3	2.50	0.46
1:K:71:PRO:HG2	1:K:400:LEU:HB3	1.96	0.46
1:L:44:ALA:HB2	1:L:114:LEU:HD21	1.97	0.46
1:L:84:GLY:HA3	1:L:340:MET:HE3	1.97	0.46
1:D:84:GLY:HA3	1:D:340:MET:HE3	1.98	0.46
1:D:238:HIS:CG	2:J:7:PRO:HB3	2.51	0.46
1:Q:253:LEU:HD22	1:Q:281:LEU:HD13	1.98	0.46
1:B:51:ALA:HB2	1:B:71:PRO:HG3	1.98	0.46
1:K:230:TYR:HE1	1:K:232:VAL:CG2	2.28	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:267:LEU:HD12	1:C:392:VAL:HG22	1.97	0.45
1:C:185:LEU:HD21	1:C:354:ALA:CB	2.46	0.45
2:H:8:THR:HA	2:I:7:PRO:HD2	1.98	0.45
1:P:238:HIS:CD2	2:R:7:PRO:HB3	2.51	0.45
1:A:225:MET:HE2	1:A:385:ASP:OD2	2.17	0.45
1:Q:253:LEU:HD22	1:Q:281:LEU:CD1	2.47	0.45
1:P:294:MET:HE1	1:P:393:ARG:NH2	2.31	0.45
1:P:51:ALA:HB2	1:P:71:PRO:HG3	2.00	0.44
1:A:51:ALA:HB2	1:A:71:PRO:HG3	1.98	0.44
1:C:51:ALA:HB2	1:C:71:PRO:HG3	1.98	0.44
1:K:115:LEU:HA	1:K:118:LEU:HD22	2.00	0.44
1:A:127:THR:HG23	1:A:207:LYS:HB3	1.99	0.44
1:L:118:LEU:HD23	1:L:401:LEU:CD1	2.48	0.44
1:A:166:LYS:HE2	1:A:191:ASP:OD1	2.18	0.43
1:A:199:LEU:HD22	1:A:350:SER:HB2	1.99	0.43
1:B:161:ILE:HD11	1:B:169:ALA:HA	2.00	0.43
1:B:238:HIS:CG	2:G:7:PRO:HB3	2.54	0.43
1:B:199:LEU:HD22	1:B:350:SER:HB2	2.00	0.43
1:L:253:LEU:HD22	1:L:281:LEU:CD1	2.49	0.43
1:Q:51:ALA:HB2	1:Q:71:PRO:HG3	2.00	0.43
1:Q:103:ARG:HB2	1:Q:106:GLU:HG3	2.01	0.43
1:K:74:VAL:O	1:K:77:SER:OG	2.26	0.42
1:C:225:MET:HE2	1:C:385:ASP:OD2	2.19	0.42
1:L:94:LYS:HB3	1:L:100:GLU:HA	2.02	0.42
1:C:44:ALA:HB2	1:C:114:LEU:HD21	2.02	0.41
1:L:225:MET:CE	1:L:275:VAL:HG21	2.50	0.41
1:A:370:ILE:O	1:A:376:LEU:HD13	2.19	0.41
1:P:238:HIS:CG	2:R:7:PRO:HB3	2.55	0.41
1:L:126:VAL:HG12	1:L:209:HIS:H	1.85	0.41
1:L:225:MET:HE2	1:L:385:ASP:OD2	2.21	0.41
1:B:343:LYS:HG2	1:B:346:LEU:H	1.85	0.41
1:C:199:LEU:HD22	1:C:350:SER:HB2	2.01	0.41
1:L:253:LEU:HD22	1:L:281:LEU:HD13	2.03	0.41
1:K:215:HIS:HB3	1:K:218:MET:HG3	2.02	0.41
1:P:39:LYS:O	1:P:43:LEU:HD13	2.21	0.41
1:D:46:ARG:HG2	1:D:98:SER:HB3	2.03	0.41
1:L:225:MET:HE3	1:L:275:VAL:HG21	2.03	0.41
1:Q:238:HIS:CG	2:T:7:PRO:HB3	2.56	0.41
1:Q:281:LEU:HA	1:Q:284:LEU:HD22	2.03	0.40
1:D:51:ALA:CB	1:D:71:PRO:HG3	2.51	0.40
2:I:14:PRO:HA	2:J:13:PRO:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:43:LEU:HD11	1:L:107:VAL:HA	2.02	0.40

There are no symmetry-related clashes.

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	376/392 (96%)	363 (96%)	12 (3%)	1 (0%)	36	45
1	B	358/392 (91%)	350 (98%)	8 (2%)	0	100	100
1	C	372/392 (95%)	358 (96%)	14 (4%)	0	100	100
1	D	362/392 (92%)	351 (97%)	11 (3%)	0	100	100
1	K	370/392 (94%)	360 (97%)	10 (3%)	0	100	100
1	L	363/392 (93%)	348 (96%)	13 (4%)	2 (1%)	21	29
1	P	369/392 (94%)	359 (97%)	10 (3%)	0	100	100
1	Q	359/392 (92%)	348 (97%)	10 (3%)	1 (0%)	36	45
2	E	16/19 (84%)	16 (100%)	0	0	100	100
2	F	15/19 (79%)	15 (100%)	0	0	100	100
2	G	14/19 (74%)	14 (100%)	0	0	100	100
2	H	16/19 (84%)	16 (100%)	0	0	100	100
2	I	15/19 (79%)	15 (100%)	0	0	100	100
2	J	14/19 (74%)	14 (100%)	0	0	100	100
2	M	16/19 (84%)	16 (100%)	0	0	100	100
2	N	16/19 (84%)	16 (100%)	0	0	100	100
2	O	16/19 (84%)	16 (100%)	0	0	100	100
2	R	16/19 (84%)	16 (100%)	0	0	100	100
2	S	16/19 (84%)	16 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	T	15/19 (79%)	15 (100%)	0	0	100	100
All	All	3114/3364 (93%)	3022 (97%)	88 (3%)	4 (0%)	48	61

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	103	ARG
1	L	105	GLU
1	A	374	GLU
1	Q	368	GLN

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	320/337 (95%)	306 (96%)	14 (4%)	25	38
1	B	308/337 (91%)	296 (96%)	12 (4%)	28	43
1	C	319/337 (95%)	305 (96%)	14 (4%)	25	38
1	D	313/337 (93%)	300 (96%)	13 (4%)	26	40
1	K	320/337 (95%)	306 (96%)	14 (4%)	25	38
1	L	314/337 (93%)	300 (96%)	14 (4%)	24	38
1	P	317/337 (94%)	300 (95%)	17 (5%)	20	30
1	Q	311/337 (92%)	295 (95%)	16 (5%)	21	33
2	E	12/12 (100%)	12 (100%)	0	100	100
2	F	11/12 (92%)	11 (100%)	0	100	100
2	G	10/12 (83%)	10 (100%)	0	100	100
2	H	12/12 (100%)	12 (100%)	0	100	100
2	I	11/12 (92%)	11 (100%)	0	100	100
2	J	10/12 (83%)	10 (100%)	0	100	100
2	M	12/12 (100%)	12 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	N	12/12 (100%)	12 (100%)	0	100	100
2	O	12/12 (100%)	12 (100%)	0	100	100
2	R	12/12 (100%)	12 (100%)	0	100	100
2	S	12/12 (100%)	12 (100%)	0	100	100
2	T	11/12 (92%)	11 (100%)	0	100	100
All	All	2659/2840 (94%)	2545 (96%)	114 (4%)	26	39

All (114) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	LEU
1	A	68	LEU
1	A	88	THR
1	A	102	LEU
1	A	118	LEU
1	A	140	VAL
1	A	189	THR
1	A	242	LEU
1	A	260	LEU
1	A	267	LEU
1	A	288	GLU
1	A	318	GLN
1	A	330	ILE
1	A	374	GLU
1	B	68	LEU
1	B	100	GLU
1	B	104	ASP
1	B	140	VAL
1	B	179	GLN
1	B	194	ARG
1	B	213	LYS
1	B	242	LEU
1	B	260	LEU
1	B	267	LEU
1	B	284	LEU
1	B	378	SER
1	C	42	THR
1	C	68	LEU
1	C	88	THR
1	C	100	GLU

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Mol	Chain	Res	Type
1	C	104	ASP
1	C	140	VAL
1	C	148	ARG
1	C	228	ARG
1	C	242	LEU
1	C	260	LEU
1	C	267	LEU
1	C	330	ILE
1	C	370	ILE
1	C	375	GLU
1	D	43	LEU
1	D	47	SER
1	D	68	LEU
1	D	104	ASP
1	D	128	TRP
1	D	140	VAL
1	D	164	ARG
1	D	213	LYS
1	D	242	LEU
1	D	260	LEU
1	D	267	LEU
1	D	318	GLN
1	D	330	ILE
1	K	42	THR
1	K	68	LEU
1	K	104	ASP
1	K	118	LEU
1	K	195	THR
1	K	242	LEU
1	K	260	LEU
1	K	267	LEU
1	K	283	LYS
1	K	294	MET
1	K	299	LYS
1	K	318	GLN
1	K	330	ILE
1	K	345	ASP
1	L	42	THR
1	L	43	LEU
1	L	68	LEU
1	L	103	ARG
1	L	232	VAL

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Mol	Chain	Res	Type
1	L	242	LEU
1	L	247	ASP
1	L	266	SER
1	L	267	LEU
1	L	288	GLU
1	L	318	GLN
1	L	330	ILE
1	L	339	ARG
1	L	369	ASP
1	P	42	THR
1	P	68	LEU
1	P	88	THR
1	P	98	SER
1	P	106	GLU
1	P	140	VAL
1	P	195	THR
1	P	242	LEU
1	P	260	LEU
1	P	267	LEU
1	P	281	LEU
1	P	283	LYS
1	P	299	LYS
1	P	318	GLN
1	P	330	ILE
1	P	391	LEU
1	P	396	GLN
1	Q	68	LEU
1	Q	98	SER
1	Q	104	ASP
1	Q	140	VAL
1	Q	195	THR
1	Q	213	LYS
1	Q	242	LEU
1	Q	260	LEU
1	Q	266	SER
1	Q	267	LEU
1	Q	275	VAL
1	Q	283	LYS
1	Q	284	LEU
1	Q	288	GLU
1	Q	318	GLN
1	Q	330	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	GLN
1	A	368	GLN
1	B	56	GLN
1	B	155	ASN
1	B	158	HIS
1	B	216	HIS
1	B	273	HIS
1	C	56	GLN
1	C	209	HIS
1	C	216	HIS
1	D	56	GLN
1	D	120	ASN
1	D	155	ASN
1	D	216	HIS
1	D	318	GLN
1	K	155	ASN
1	K	216	HIS
1	K	254	GLN
1	K	320	HIS
1	L	120	ASN
1	L	155	ASN
1	L	216	HIS
1	L	318	GLN
1	P	108	HIS
1	P	171	GLN
1	P	209	HIS
1	P	216	HIS
1	P	254	GLN
1	P	320	HIS
1	P	368	GLN
1	Q	120	ASN
1	Q	209	HIS
1	Q	318	GLN
1	Q	320	HIS
1	Q	368	GLN

5.3.3 RNA ⓘ

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](#)

5 ligands 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	SIN	C	1418	-	7,7,7	1.71	2 (28%)	8,8,8	1.39	2 (25%)
3	SIN	A	1421	-	7,7,7	1.41	2 (28%)	8,8,8	1.49	2 (25%)
3	SIN	K	1415	-	7,7,7	1.41	2 (28%)	8,8,8	1.78	2 (25%)
3	SIN	Q	1412	-	7,7,7	1.44	2 (28%)	8,8,8	1.49	2 (25%)
3	SIN	P	1416	-	7,7,7	1.47	2 (28%)	8,8,8	1.55	2 (25%)

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	SIN	C	1418	-	-	2/5/5/5	-
3	SIN	A	1421	-	-	5/5/5/5	-
3	SIN	K	1415	-	-	3/5/5/5	-
3	SIN	Q	1412	-	-	2/5/5/5	-
3	SIN	P	1416	-	-	2/5/5/5	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1418	SIN	O1-C1	3.66	1.34	1.22
3	P	1416	SIN	O1-C1	3.02	1.32	1.22
3	K	1415	SIN	O1-C1	2.97	1.31	1.22
3	A	1421	SIN	O1-C1	2.90	1.31	1.22
3	Q	1412	SIN	O1-C1	2.81	1.31	1.22
3	C	1418	SIN	O2-C1	-2.60	1.22	1.30
3	Q	1412	SIN	O2-C1	-2.44	1.22	1.30
3	P	1416	SIN	O2-C1	-2.35	1.23	1.30
3	A	1421	SIN	O2-C1	-2.28	1.23	1.30
3	K	1415	SIN	O2-C1	-2.15	1.23	1.30

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	K	1415	SIN	O1-C1-C2	-3.61	111.65	123.09
3	K	1415	SIN	O2-C1-C2	3.42	124.81	114.00
3	P	1416	SIN	O2-C1-C2	3.10	123.79	114.00
3	P	1416	SIN	O1-C1-C2	-3.06	113.38	123.09
3	Q	1412	SIN	O2-C1-C2	2.99	123.44	114.00
3	Q	1412	SIN	O1-C1-C2	-2.94	113.78	123.09
3	A	1421	SIN	O2-C1-C2	2.91	123.19	114.00
3	A	1421	SIN	O1-C1-C2	-2.90	113.89	123.09
3	C	1418	SIN	O1-C1-C2	-2.79	114.24	123.09
3	C	1418	SIN	O2-C1-C2	2.72	122.58	114.00

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	K	1415	SIN	C1-C2-C3-C4
3	A	1421	SIN	C1-C2-C3-C4
3	P	1416	SIN	O2-C1-C2-C3
3	P	1416	SIN	O1-C1-C2-C3
3	C	1418	SIN	C2-C3-C4-O3
3	C	1418	SIN	C2-C3-C4-O4
3	K	1415	SIN	C2-C3-C4-O3
3	K	1415	SIN	C2-C3-C4-O4
3	A	1421	SIN	O1-C1-C2-C3
3	A	1421	SIN	O2-C1-C2-C3
3	A	1421	SIN	C2-C3-C4-O4
3	A	1421	SIN	C2-C3-C4-O3
3	Q	1412	SIN	O1-C1-C2-C3
3	Q	1412	SIN	O2-C1-C2-C3

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1418	SIN	1	0
3	A	1421	SIN	1	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	380/392 (96%)	0.16	10 (2%) 57 58	28, 57, 105, 133	0
1	B	364/392 (92%)	0.16	11 (3%) 52 54	22, 54, 107, 147	0
1	C	376/392 (95%)	0.21	11 (2%) 53 55	24, 57, 118, 148	0
1	D	368/392 (93%)	0.09	11 (2%) 52 54	22, 49, 102, 133	0
1	K	374/392 (95%)	0.08	6 (1%) 70 71	25, 53, 96, 125	0
1	L	369/392 (94%)	0.38	23 (6%) 26 26	22, 59, 119, 140	0
1	P	373/392 (95%)	0.16	8 (2%) 63 64	26, 59, 103, 131	0
1	Q	365/392 (93%)	0.08	11 (3%) 52 54	24, 50, 98, 125	0
2	E	17/19 (89%)	-0.31	0 100 100	24, 30, 66, 75	0
2	F	16/19 (84%)	-0.60	0 100 100	27, 35, 59, 68	0
2	G	15/19 (78%)	-0.44	1 (6%) 24 23	26, 31, 56, 91	0
2	H	17/19 (89%)	-0.37	0 100 100	21, 29, 65, 76	0
2	I	16/19 (84%)	-0.56	0 100 100	22, 32, 63, 71	0
2	J	15/19 (78%)	-0.47	0 100 100	19, 30, 57, 90	0
2	M	17/19 (89%)	-0.45	0 100 100	23, 31, 68, 78	0
2	N	17/19 (89%)	-0.38	0 100 100	27, 32, 71, 78	0
2	O	17/19 (89%)	-0.30	0 100 100	24, 31, 87, 100	0
2	R	17/19 (89%)	-0.57	0 100 100	26, 32, 61, 67	0
2	S	17/19 (89%)	-0.29	1 (5%) 28 27	29, 34, 64, 76	0
2	T	16/19 (84%)	-0.38	1 (6%) 26 25	28, 33, 65, 92	0
All	All	3166/3364 (94%)	0.13	94 (2%) 52 54	19, 53, 106, 148	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	370	ILE	5.1
1	Q	378	SER	5.1
1	B	163	PHE	4.2
1	B	128	TRP	3.6
1	B	48	ALA	3.5
1	C	119	SER	3.4
1	C	371	TYR	3.3
2	T	16	PRO	3.2
1	A	126	VAL	3.2
1	B	412	ASP	3.1
1	C	370	ILE	3.1
1	L	82	SER	3.1
1	L	338	SER	3.1
1	L	344	LYS	3.0
1	P	142	PHE	3.0
1	D	47	SER	3.0
1	P	192	VAL	2.9
1	B	366	PHE	2.9
1	Q	344	LYS	2.9
1	A	420	GLU	2.9
1	K	163	PHE	2.8
1	L	346	LEU	2.8
1	P	163	PHE	2.8
1	P	38	PRO	2.8
1	A	119	SER	2.8
1	P	120	ASN	2.7
1	D	366	PHE	2.7
1	B	47	SER	2.7
1	L	145	ASP	2.7
2	G	15	GLY	2.7
1	A	371	TYR	2.7
1	P	160	LYS	2.7
1	C	126	VAL	2.7
1	L	110	GLY	2.6
1	A	372	GLY	2.6
1	Q	345	ASP	2.6
1	K	287	LYS	2.6
1	K	346	LEU	2.5
1	B	410	LYS	2.5
1	C	417	GLU	2.5
1	D	140	VAL	2.5
1	P	195	THR	2.5
1	C	372	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	346	LEU	2.5
1	B	125	ASN	2.5
1	L	414	MET	2.5
1	Q	142	PHE	2.4
1	A	370	ILE	2.4
1	L	128	TRP	2.4
1	L	38	PRO	2.3
1	Q	162	ASN	2.3
1	A	142	PHE	2.3
1	B	162	ASN	2.3
1	K	342	GLY	2.3
1	K	35	MET	2.3
1	D	120	ASN	2.3
1	L	411	GLY	2.3
1	L	47	SER	2.3
1	Q	369	ASP	2.3
1	D	163	PHE	2.3
1	A	346	LEU	2.2
1	L	146	PHE	2.2
1	L	124	ARG	2.2
1	L	44	ALA	2.2
1	Q	125	ASN	2.2
1	D	35	MET	2.2
1	C	192	VAL	2.2
1	L	120	ASN	2.2
1	A	376	LEU	2.2
1	C	102	LEU	2.2
1	L	341	SER	2.2
1	B	39	LYS	2.2
1	Q	366	PHE	2.1
1	Q	367	ASP	2.1
1	D	48	ALA	2.1
1	L	48	ALA	2.1
1	A	36	LEU	2.1
1	B	118	LEU	2.1
1	D	43	LEU	2.1
1	D	411	GLY	2.1
1	Q	411	GLY	2.1
2	S	17	PRO	2.1
1	C	140	VAL	2.1
1	L	80	LEU	2.1
1	K	195	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	L	157	GLU	2.1
1	D	39	LYS	2.1
1	P	345	ASP	2.1
1	L	142	PHE	2.0
1	Q	192	VAL	2.0
1	C	99	ALA	2.0
1	D	344	LYS	2.0
1	L	96	VAL	2.0
1	L	35	MET	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($\text{\AA}^2$)	Q<0.9
3	SIN	K	1415	8/8	0.76	0.13	56,60,63,64	0
3	SIN	P	1416	8/8	0.76	0.15	71,73,76,77	0
3	SIN	A	1421	8/8	0.77	0.14	59,62,64,65	0
3	SIN	C	1418	8/8	0.77	0.13	55,58,62,64	0
3	SIN	Q	1412	8/8	0.84	0.10	50,60,68,71	0

6.5 Other polymers [i](#)

There are no such residues in this entry.