



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:56 PM UTC

PDB ID : 6IQS / pdb_00006iqs
Title : Crystal structure of Prc with L245A and L340G mutations in complex with NlpI
Authors : Chuech, C.K.; Chang, C.I.
Deposited on : 2018-11-08
Resolution : 2.69 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

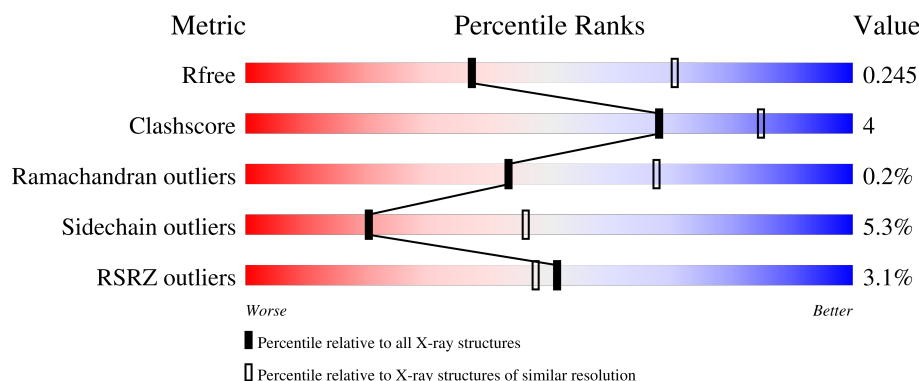
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.69 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	296	77% 7% 14%
1	B	296	78% 8% 12%
2	C	688	3% 69% 8% 21%
2	D	688	3% 67% 9% 22%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12892 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 Lipoprotein NlpI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	256	Total	C	N	O	S	0	0	0
			2071	1325	342	401	3			
1	B	259	Total	C	N	O	S	0	0	0
			2097	1340	347	407	3			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP P0AFB1
A	2	GLY	-	expression tag	UNP P0AFB1
A	3	SER	-	expression tag	UNP P0AFB1
A	4	SER	-	expression tag	UNP P0AFB1
A	5	HIS	-	expression tag	UNP P0AFB1
A	6	HIS	-	expression tag	UNP P0AFB1
A	7	HIS	-	expression tag	UNP P0AFB1
A	8	HIS	-	expression tag	UNP P0AFB1
A	9	HIS	-	expression tag	UNP P0AFB1
A	10	HIS	-	expression tag	UNP P0AFB1
A	11	SER	-	expression tag	UNP P0AFB1
A	12	SER	-	expression tag	UNP P0AFB1
A	13	GLY	-	expression tag	UNP P0AFB1
A	14	LEU	-	expression tag	UNP P0AFB1
A	15	VAL	-	expression tag	UNP P0AFB1
A	16	PRO	-	expression tag	UNP P0AFB1
A	17	ARG	-	expression tag	UNP P0AFB1
A	18	GLY	-	expression tag	UNP P0AFB1
A	19	SER	-	expression tag	UNP P0AFB1
A	20	HIS	-	expression tag	UNP P0AFB1
A	21	MET	-	expression tag	UNP P0AFB1
B	1	MET	-	initiating methionine	UNP P0AFB1
B	2	GLY	-	expression tag	UNP P0AFB1
B	3	SER	-	expression tag	UNP P0AFB1
B	4	SER	-	expression tag	UNP P0AFB1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	5	HIS	-	expression tag	UNP P0AFB1
B	6	HIS	-	expression tag	UNP P0AFB1
B	7	HIS	-	expression tag	UNP P0AFB1
B	8	HIS	-	expression tag	UNP P0AFB1
B	9	HIS	-	expression tag	UNP P0AFB1
B	10	HIS	-	expression tag	UNP P0AFB1
B	11	SER	-	expression tag	UNP P0AFB1
B	12	SER	-	expression tag	UNP P0AFB1
B	13	GLY	-	expression tag	UNP P0AFB1
B	14	LEU	-	expression tag	UNP P0AFB1
B	15	VAL	-	expression tag	UNP P0AFB1
B	16	PRO	-	expression tag	UNP P0AFB1
B	17	ARG	-	expression tag	UNP P0AFB1
B	18	GLY	-	expression tag	UNP P0AFB1
B	19	SER	-	expression tag	UNP P0AFB1
B	20	HIS	-	expression tag	UNP P0AFB1
B	21	MET	-	expression tag	UNP P0AFB1

- Molecule 2 is a protein called Tail-specific protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	541	Total	C	N	O	S	0	0	0
			4348	2739	760	841	8			
2	D	538	Total	C	N	O	S	0	0	0
			4325	2725	754	838	8			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	245	ALA	LEU	engineered mutation	UNP P23865
C	340	GLY	LEU	engineered mutation	UNP P23865
C	683	HIS	-	expression tag	UNP P23865
C	684	HIS	-	expression tag	UNP P23865
C	685	HIS	-	expression tag	UNP P23865
C	686	HIS	-	expression tag	UNP P23865
C	687	HIS	-	expression tag	UNP P23865
C	688	HIS	-	expression tag	UNP P23865
D	245	ALA	LEU	engineered mutation	UNP P23865
D	340	GLY	LEU	engineered mutation	UNP P23865
D	683	HIS	-	expression tag	UNP P23865
D	684	HIS	-	expression tag	UNP P23865
D	685	HIS	-	expression tag	UNP P23865

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Chain	Residue	Modelled	Actual	Comment	Reference
D	686	HIS	-	expression tag	UNP P23865
D	687	HIS	-	expression tag	UNP P23865
D	688	HIS	-	expression tag	UNP P23865

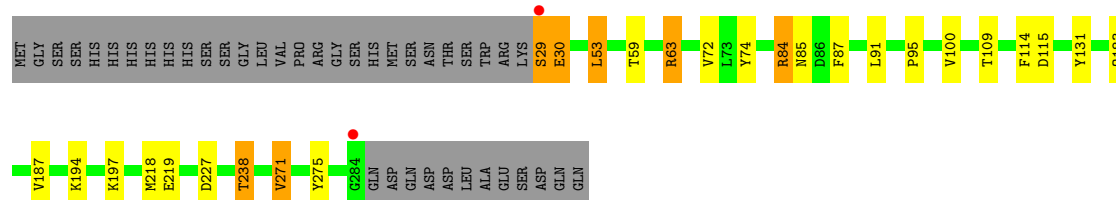
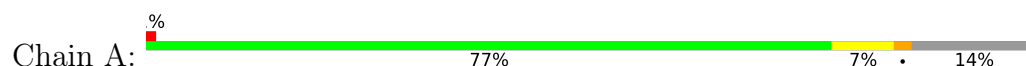
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total 11	O 11	0	0
3	B	5	Total 5	O 5	0	0
3	C	13	Total 13	O 13	0	0
3	D	22	Total 22	O 22	0	0

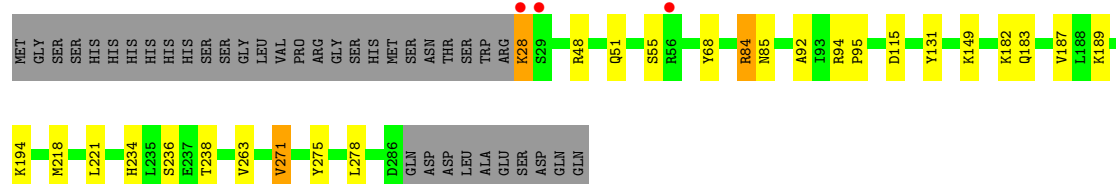
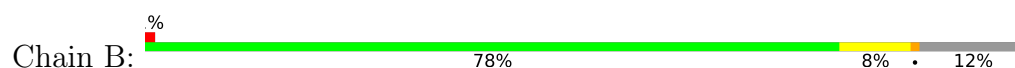
3 Residue-property plots [i](#)

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.

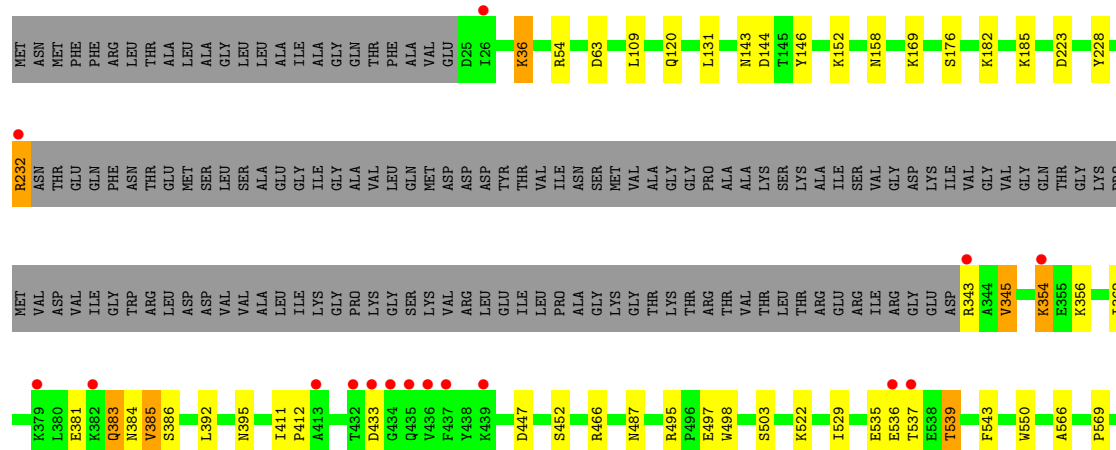
• Molecule 1: Lipoprotein NlpI



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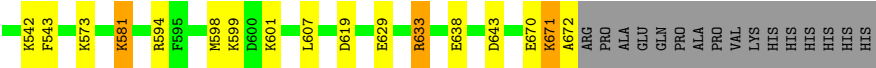
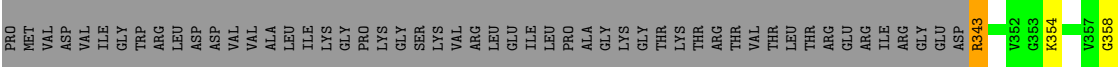
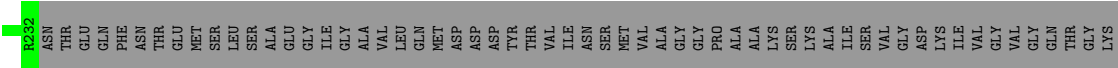


• Molecule 2: Tail-specific protease





● Molecule 2: Tail-specific protease



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	119.34Å 144.41Å 152.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.91 – 2.69 29.91 – 2.69	Depositor EDS
% Data completeness (in resolution range)	93.2 (29.91-2.69) 93.4 (29.91-2.69)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.68Å)	Xtrriage
Refinement program	REFMAC 5.8.0230	Depositor
R, R_{free}	0.205 , 0.245 0.208 , 0.245	Depositor DCC
R_{free} test set	3435 reflections (4.64%)	wwPDB-VP
Wilson B-factor (Å ²)	45.3	Xtrriage
Anisotropy	0.265	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 24.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12892	wwPDB-VP
Average B, all atoms (Å ²)	54.0	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 41.09 % 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 2.4836e-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

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.27	10/2115 (0.5%)	1.29	5/2871 (0.2%)
1	B	1.22	7/2141 (0.3%)	1.28	5/2905 (0.2%)
2	C	1.15	10/4427 (0.2%)	1.25	6/5977 (0.1%)
2	D	1.14	9/4403 (0.2%)	1.27	14/5944 (0.2%)
All	All	1.18	36/13086 (0.3%)	1.27	30/17697 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
2	D	0	1
All	All	0	4

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	53	LEU	C-O	-8.61	1.13	1.24
1	B	84	ARG	NE-CZ	8.45	1.42	1.33
1	A	115	ASP	N-CA	-7.96	1.37	1.46
2	D	158	ASN	C-O	-7.61	1.16	1.24
2	D	503	SER	C-O	-7.10	1.15	1.23
2	D	426	VAL	C-O	-7.00	1.16	1.24
1	B	85	ASN	C-O	-6.83	1.16	1.24
1	A	109	THR	C-O	-6.77	1.16	1.24
1	A	74	TYR	C-O	-6.75	1.16	1.24
1	A	95	PRO	C-O	-6.57	1.16	1.24
2	C	158	ASN	C-O	-6.57	1.16	1.23
1	B	263	VAL	C-O	-6.55	1.16	1.24
2	C	487	ASN	C-O	-6.54	1.15	1.23
2	C	638	GLU	C-O	-6.20	1.16	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	529	ILE	C-O	-6.13	1.17	1.24
1	A	114	PHE	C-O	-6.03	1.17	1.24
2	D	638	GLU	C-O	-6.03	1.16	1.23
1	A	72	VAL	N-CA	-5.92	1.39	1.46
2	C	152	LYS	N-CA	5.88	1.52	1.46
2	D	88	SER	C-O	5.65	1.30	1.24
1	B	236	SER	C-O	-5.64	1.17	1.24
2	D	39	THR	N-CA	5.57	1.53	1.46
2	C	539	THR	N-CA	5.52	1.52	1.45
2	C	228	TYR	C-O	-5.50	1.17	1.24
2	C	503	SER	C-O	-5.50	1.17	1.23
2	D	506	TYR	C-O	-5.42	1.17	1.23
2	C	345	VAL	C-O	-5.39	1.17	1.24
1	B	92	ALA	C-O	-5.32	1.17	1.24
1	B	131	TYR	N-CA	5.25	1.53	1.46
1	A	59	THR	C-O	-5.25	1.16	1.23
1	A	131	TYR	N-CA	5.24	1.53	1.46
2	D	599	LYS	N-CA	5.23	1.53	1.46
2	C	232	ARG	N-CA	5.22	1.56	1.46
1	A	238	THR	N-CA	-5.19	1.39	1.46
1	B	95	PRO	C-O	-5.04	1.17	1.24
2	D	228	TYR	C-O	-5.03	1.17	1.24

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	115	ASP	CB-CA-C	11.53	129.25	110.92
1	A	115	ASP	N-CA-CB	-9.75	95.01	109.82
2	C	54	ARG	CG-CD-NE	7.72	129.00	112.00
2	D	414	GLY	CA-C-N	7.43	127.19	119.76
2	D	414	GLY	C-N-CA	7.43	127.19	119.76
2	D	495	ARG	CG-CD-NE	-7.08	96.42	112.00
2	D	416	ILE	CB-CA-C	-6.20	105.61	111.06
1	A	85	ASN	CB-CA-C	-6.08	101.30	110.90
1	B	84	ARG	NE-CZ-NH2	6.03	124.63	119.20
2	D	495	ARG	CB-CG-CD	6.02	125.14	111.30
2	D	223	ASP	CB-CA-C	5.81	117.85	110.76
1	A	227	ASP	CA-CB-CG	5.81	118.41	112.60
1	B	182	LYS	CB-CG-CD	5.76	124.56	111.30
2	C	223	ASP	CB-CA-C	5.42	117.37	110.76
2	D	619	ASP	CA-CB-CG	5.39	117.99	112.60
2	D	672	ALA	CA-C-O	-5.35	111.70	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	222	ILE	N-CA-C	-5.34	105.88	110.74
1	A	84	ARG	CB-CG-CD	5.30	123.50	111.30
2	C	447	ASP	CA-CB-CG	5.28	117.88	112.60
2	D	638	GLU	CB-CG-CD	5.25	121.52	112.60
2	D	672	ALA	N-CA-C	-5.24	96.34	111.00
2	D	581	LYS	CB-CG-CD	5.18	123.22	111.30
2	C	36	LYS	CB-CG-CD	5.10	123.02	111.30
1	B	68	TYR	N-CA-C	-5.10	105.62	111.07
2	C	63	ASP	CA-CB-CG	5.05	117.65	112.60
2	C	619	ASP	CA-CB-CG	5.04	117.64	112.60
2	D	633	ARG	CB-CA-C	5.04	119.15	110.79
1	B	194	LYS	CB-CA-C	-5.04	101.00	110.67
2	D	495	ARG	CD-NE-CZ	5.04	131.45	124.40
1	B	115	ASP	CB-CA-C	-5.01	102.53	110.84

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	29	SER	Peptide
1	A	63	ARG	Sidechain
1	A	84	ARG	Sidechain
2	D	25	ASP	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	2071	0	2007	26	0
1	B	2097	0	2032	26	0
2	C	4348	0	4324	24	0
2	D	4325	0	4297	35	0
3	A	11	0	0	0	0
3	B	5	0	0	0	0
3	C	13	0	0	0	0
3	D	22	0	0	1	0
All	All	12892	0	12660	92	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 4.

All (92) 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:A:30:GLU:CD	1:B:28:LYS:NZ	1.82	1.36
1:A:30:GLU:CD	1:B:28:LYS:HZ1	1.36	1.28
1:B:48:ARG:NH1	1:B:51:GLN:OE1	1.77	1.17
2:D:537:THR:O	2:D:543:PHE:CZ	2.14	1.00
1:A:30:GLU:CD	1:B:28:LYS:HZ2	1.72	0.96
2:C:602:ARG:O	2:C:602:ARG:NH2	2.11	0.82
2:C:537:THR:O	2:C:543:PHE:CZ	2.33	0.81
2:C:497:GLU:OE1	2:C:497:GLU:N	2.13	0.79
2:D:375:VAL:O	2:D:378:GLN:HG2	1.84	0.78
1:A:30:GLU:OE2	1:B:28:LYS:NZ	2.17	0.76
1:A:30:GLU:CD	1:B:28:LYS:CE	2.61	0.74
2:C:362:ILE:O	2:C:395:ASN:ND2	2.21	0.73
2:C:539:THR:O	2:C:543:PHE:CE2	2.41	0.73
1:A:30:GLU:CG	1:B:28:LYS:NZ	2.51	0.72
2:D:537:THR:O	2:D:543:PHE:CE2	2.43	0.72
2:C:566:ALA:O	2:C:569:PRO:HD2	1.87	0.72
1:A:30:GLU:CG	1:B:28:LYS:CE	2.68	0.71
2:D:411:ILE:O	3:D:701:HOH:O	2.09	0.71
2:D:362:ILE:O	2:D:395:ASN:ND2	2.24	0.70
2:C:146:TYR:HB2	2:C:607:LEU:HD21	1.73	0.70
1:A:30:GLU:HG3	1:B:28:LYS:CD	2.22	0.69
2:D:146:TYR:HB2	2:D:607:LEU:HD21	1.74	0.69
2:D:497:GLU:H	2:D:497:GLU:CD	1.99	0.69
1:A:30:GLU:OE1	1:B:28:LYS:NZ	2.10	0.66
2:D:496:PRO:N	2:D:497:GLU:OE1	2.29	0.66
2:D:496:PRO:CD	2:D:497:GLU:OE1	2.45	0.65
1:A:30:GLU:CG	1:B:28:LYS:HZ2	2.08	0.64
1:B:218:MET:HE1	1:B:238:THR:HG21	1.79	0.64
2:D:496:PRO:HD2	2:D:497:GLU:OE1	1.99	0.63
2:D:539:THR:O	2:D:543:PHE:CE2	2.52	0.63
1:A:30:GLU:HG3	1:B:28:LYS:CE	2.29	0.63
2:C:535:GLU:OE2	2:C:596:ASN:ND2	2.31	0.62
1:A:30:GLU:HB2	1:B:28:LYS:HD2	1.81	0.62
2:D:358:GLY:HA3	2:D:380:LEU:CD2	2.30	0.62
2:C:356:LYS:C	2:C:385:VAL:HG12	2.25	0.61
2:C:539:THR:O	2:C:543:PHE:HE2	1.84	0.60
1:A:218:MET:HE1	1:A:238:THR:HG21	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:539:THR:O	2:C:543:PHE:CD2	2.56	0.58
2:D:539:THR:O	2:D:543:PHE:CD2	2.58	0.57
1:A:271:VAL:HG22	1:A:275:TYR:HE2	1.69	0.57
1:A:30:GLU:OE1	1:B:28:LYS:CE	2.52	0.57
2:C:381:GLU:O	2:C:383:GLN:O	2.21	0.57
1:A:30:GLU:HG3	1:B:28:LYS:HZ2	1.69	0.57
2:D:413:ALA:HB3	2:D:464:TYR:CE2	2.42	0.55
1:B:271:VAL:HG22	1:B:275:TYR:HE2	1.72	0.54
2:D:437:PHE:CD1	2:D:437:PHE:N	2.75	0.54
2:D:437:PHE:N	2:D:437:PHE:HD1	2.04	0.54
1:A:30:GLU:OE1	1:B:28:LYS:HE3	2.08	0.53
2:C:598:MET:HE3	2:C:601:LYS:HE2	1.90	0.53
2:C:566:ALA:O	2:C:569:PRO:CD	2.57	0.52
2:D:78:LEU:O	2:D:123:ARG:NH2	2.43	0.52
2:D:435:GLN:HB3	2:D:437:PHE:HE1	1.75	0.52
1:A:30:GLU:HG3	1:B:28:LYS:NZ	2.24	0.51
2:C:535:GLU:CD	2:C:596:ASN:HD21	2.15	0.51
2:D:495:ARG:C	2:D:497:GLU:OE1	2.53	0.51
2:C:601:LYS:O	2:C:601:LYS:HD2	2.11	0.50
1:A:183:GLN:O	1:A:187:VAL:HG23	2.12	0.50
1:A:30:GLU:HG3	1:B:28:LYS:HD2	1.91	0.49
1:A:271:VAL:HG22	1:A:275:TYR:CE2	2.47	0.49
2:D:358:GLY:HA3	2:D:380:LEU:HD21	1.93	0.49
1:A:53:LEU:O	1:A:63:ARG:NH2	2.41	0.48
2:D:539:THR:O	2:D:543:PHE:HE2	1.95	0.47
1:A:87:PHE:HB3	1:A:100:VAL:HG13	1.96	0.47
1:B:183:GLN:O	1:B:187:VAL:HG23	2.13	0.47
2:C:354:LYS:HD3	2:C:354:LYS:H	1.80	0.47
2:D:358:GLY:HA3	2:D:380:LEU:HD22	1.97	0.47
1:B:271:VAL:HG22	1:B:275:TYR:CE2	2.51	0.46
2:C:392:LEU:O	2:C:395:ASN:HB2	2.16	0.46
2:D:435:GLN:HB3	2:D:437:PHE:CE1	2.51	0.46
2:C:495:ARG:HD3	2:C:498:TRP:CH2	2.50	0.46
2:C:466:ARG:O	2:C:466:ARG:HD2	2.16	0.45
2:D:343:ARG:HE	2:D:343:ARG:C	2.25	0.45
1:A:30:GLU:CG	1:B:28:LYS:HE3	2.44	0.45
1:A:30:GLU:CD	1:B:28:LYS:HE3	2.42	0.45
2:D:89:ASP:HB3	2:D:118:LEU:HD21	1.99	0.45
1:B:221:LEU:HG	1:B:234:HIS:HB3	1.99	0.45
2:C:522:LYS:HG3	2:C:550:TRP:CD1	2.52	0.44
2:D:484:ARG:HD3	2:D:484:ARG:HA	1.75	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:392:LEU:O	2:D:395:ASN:HB2	2.18	0.44
1:B:28:LYS:HE2	1:B:28:LYS:HB2	1.67	0.43
2:D:54:ARG:HH11	2:D:54:ARG:HG2	1.84	0.43
1:A:91:LEU:HG	1:A:100:VAL:HG11	2.00	0.43
2:C:356:LYS:O	2:C:385:VAL:HG12	2.18	0.42
2:D:671:LYS:HE3	2:D:671:LYS:HA	2.01	0.42
2:C:411:ILE:HA	2:C:412:PRO:HD3	1.86	0.42
2:D:447:ASP:HA	2:D:472:GLU:HB2	2.02	0.42
2:D:539:THR:C	2:D:543:PHE:HD2	2.27	0.42
2:D:358:GLY:N	2:D:385:VAL:HG21	2.35	0.42
2:D:56:HIS:CD2	2:D:506:TYR:HB2	2.55	0.41
2:D:415:PRO:HD2	2:D:464:TYR:OH	2.21	0.40
2:C:497:GLU:H	2:C:497:GLU:CD	2.16	0.40
2:D:594:ARG:O	2:D:598:MET:HG2	2.21	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	254/296 (86%)	248 (98%)	6 (2%)	0	100	100
1	B	257/296 (87%)	248 (96%)	9 (4%)	0	100	100
2	C	537/688 (78%)	518 (96%)	18 (3%)	1 (0%)	43	68
2	D	534/688 (78%)	511 (96%)	21 (4%)	2 (0%)	30	54
All	All	1582/1968 (80%)	1525 (96%)	54 (3%)	3 (0%)	43	68

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	452	SER

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Mol	Chain	Res	Type
2	D	452	SER
2	D	412	PRO

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	217/253 (86%)	211 (97%)	6 (3%)	38	68
1	B	220/253 (87%)	212 (96%)	8 (4%)	31	60
2	C	471/588 (80%)	440 (93%)	31 (7%)	15	36
2	D	469/588 (80%)	441 (94%)	28 (6%)	17	41
All	All	1377/1682 (82%)	1304 (95%)	73 (5%)	20	46

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	SER
1	A	30	GLU
1	A	194	LYS
1	A	197	LYS
1	A	219	GLU
1	A	271	VAL
1	B	28	LYS
1	B	55	SER
1	B	84	ARG
1	B	94	ARG
1	B	149	LYS
1	B	189	LYS
1	B	271	VAL
1	B	278	LEU
2	C	36	LYS
2	C	109	LEU
2	C	120	GLN
2	C	131	LEU
2	C	143	ASN

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Mol	Chain	Res	Type
2	C	144	ASP
2	C	169	LYS
2	C	176	SER
2	C	182	LYS
2	C	185	LYS
2	C	232	ARG
2	C	343	ARG
2	C	345	VAL
2	C	354	LYS
2	C	383	GLN
2	C	384	ASN
2	C	385	VAL
2	C	386	SER
2	C	433	ASP
2	C	536	GLU
2	C	573	LYS
2	C	589	MET
2	C	598	MET
2	C	601	LYS
2	C	629	GLU
2	C	633	ARG
2	C	637	PRO
2	C	643	ASP
2	C	647	LYS
2	C	650	GLN
2	C	668	LYS
2	D	28	ARG
2	D	120	GLN
2	D	144	ASP
2	D	185	LYS
2	D	193	ARG
2	D	343	ARG
2	D	354	LYS
2	D	380	LEU
2	D	382	LYS
2	D	383	GLN
2	D	385	VAL
2	D	433	ASP
2	D	437	PHE
2	D	439	LYS
2	D	482	GLN
2	D	497	GLU

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Mol	Chain	Res	Type
2	D	532	THR
2	D	538	GLU
2	D	539	THR
2	D	542	LYS
2	D	573	LYS
2	D	581	LYS
2	D	601	LYS
2	D	629	GLU
2	D	633	ARG
2	D	643	ASP
2	D	670	GLU
2	D	671	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	89	GLN
1	B	41	GLN
1	B	132	ASN
1	B	265	ASN
2	C	120	GLN
2	C	204	GLN
2	C	481	GLN
2	C	603	ASN
2	D	31	GLN
2	D	120	GLN
2	D	395	ASN
2	D	418	GLN
2	D	603	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [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	256/296 (86%)	-0.56	2 (0%) 82 81	22, 36, 64, 81	0
1	B	259/296 (87%)	-0.54	3 (1%) 76 75	23, 36, 65, 103	0
2	C	541/688 (78%)	0.04	24 (4%) 39 35	31, 55, 101, 130	0
2	D	538/688 (78%)	0.04	21 (3%) 43 39	35, 57, 98, 141	0
All	All	1594/1968 (80%)	-0.15	50 (3%) 51 48	22, 51, 94, 141	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	672	ALA	4.9
2	C	675	ALA	4.6
2	D	414	GLY	4.5
1	B	28	LYS	4.4
2	C	595	PHE	4.1
2	D	595	PHE	3.9
2	D	413	ALA	3.8
2	D	537	THR	3.8
2	C	354	LYS	3.7
2	C	439	LYS	3.7
1	A	29	SER	3.6
2	D	343	ARG	3.6
2	C	433	ASP	3.1
2	D	412	PRO	3.0
1	A	284	GLY	2.7
2	C	232	ARG	2.7
1	B	56	ARG	2.7
2	C	668	LYS	2.7
2	D	193	ARG	2.6
2	C	434	GLY	2.6
2	C	671	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
2	D	601	LYS	2.5
2	D	598	MET	2.5
2	D	538	GLU	2.5
2	D	599	LYS	2.4
2	C	379	LYS	2.4
2	C	537	THR	2.4
2	C	674	PRO	2.4
2	C	435	GLN	2.4
2	C	672	ALA	2.4
2	D	26	ILE	2.4
2	C	413	ALA	2.3
2	C	598	MET	2.3
2	D	352	VAL	2.3
1	B	29	SER	2.2
2	D	434	GLY	2.2
2	C	382	LYS	2.2
2	D	439	LYS	2.2
2	C	536	GLU	2.1
2	C	343	ARG	2.1
2	C	26	ILE	2.1
2	C	436	VAL	2.1
2	D	357	VAL	2.1
2	C	437	PHE	2.1
2	D	433	ASP	2.1
2	D	354	LYS	2.0
2	D	232	ARG	2.0
2	C	601	LYS	2.0
2	D	671	LYS	2.0
2	C	432	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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.