Full wwPDB X-ray Structure Validation Report ⓘ

Nov 14, 2023 – 10:28 PM JST

PDB ID	:	6IQS
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:

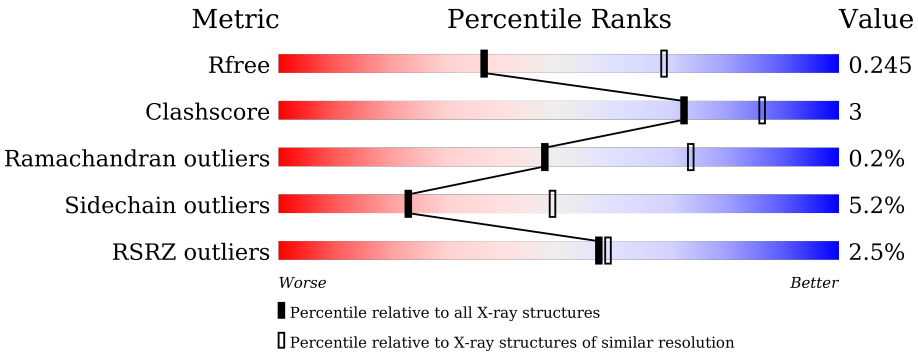
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (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 <=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	80% 5% • 14%
1	B	296	% 80% 7% • 12%
2	C	688	3% 70% 8% • 21%
2	D	688	2% 68% 8% • 22%

2 Entry composition (i)

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.

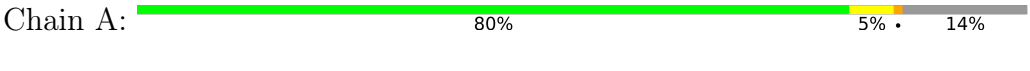
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	11	Total O 11 11	0	0
3	B	5	Total O 5 5	0	0
3	C	13	Total O 13 13	0	0
3	D	22	Total O 22 22	0	0

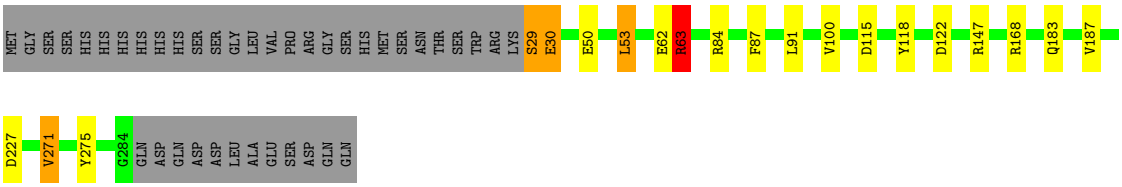
3 Residue-property plots (i)

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

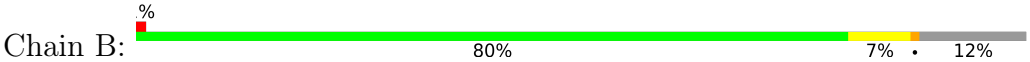
Chain A:

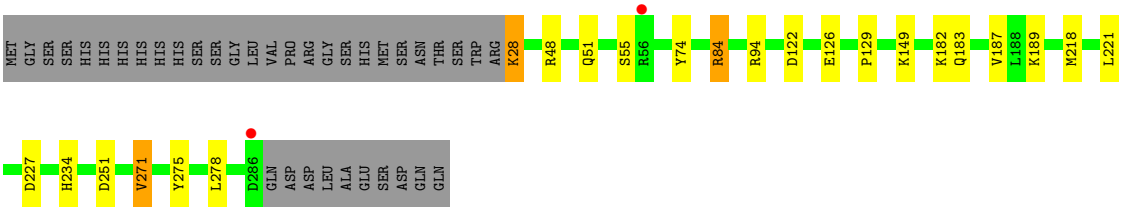




• Molecule 1: Lipoprotein NlpI

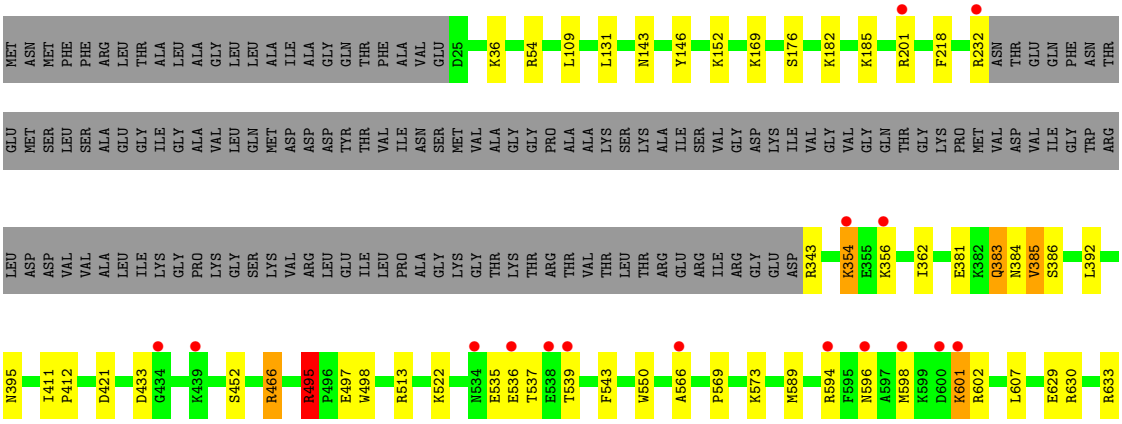
Chain B:

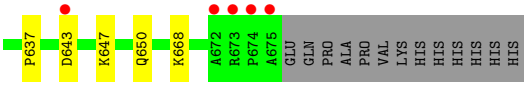




• Molecule 2: Tail-specific protease

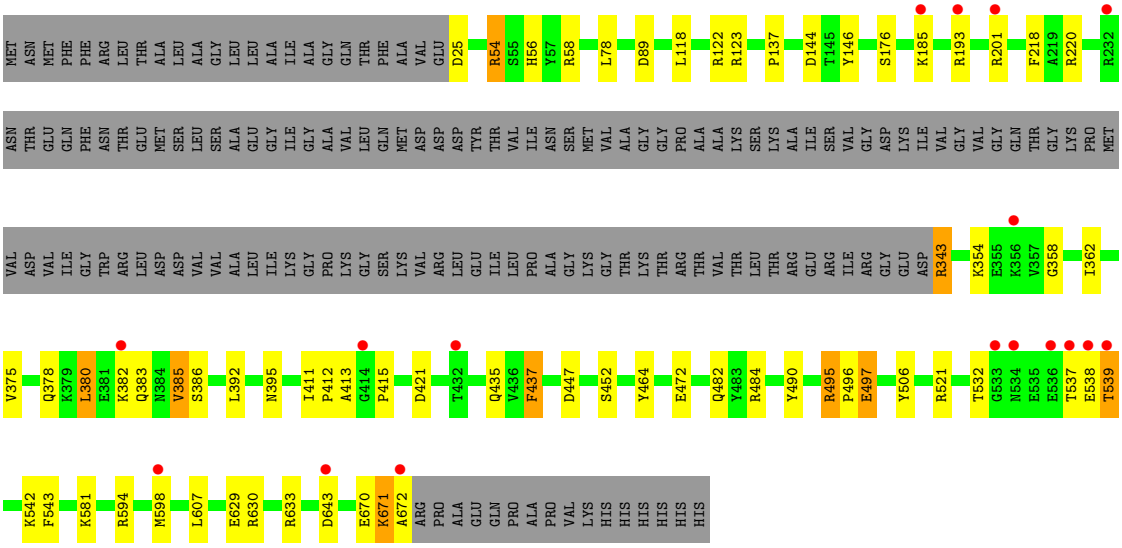
Chain C:





• Molecule 2: Tail-specific protease

Chain D:



4 Data and refinement statistics (i)

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 29.89 – 2.69	Depositor EDS
% Data completeness (in resolution range)	93.2 (29.91-2.69) 93.3 (29.89-2.69)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.02 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5.8.0230	Depositor
R, R_{free}	0.205 , 0.245 0.209 , 0.245	Depositor DCC
R_{free} test set	3435 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å^2)	45.3	Xtriage
Anisotropy	0.265	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 26.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
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

Xtriage'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.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|$ >5	RMSZ	#$\|Z\|$ >5
1	A	0.92	3/2115 (0.1%)	1.10	11/2871 (0.4%)
1	B	0.88	1/2141 (0.0%)	1.06	7/2905 (0.2%)
2	C	0.80	0/4427	1.00	8/5977 (0.1%)
2	D	0.80	0/4403	1.01	12/5944 (0.2%)
All	All	0.83	4/13086 (0.0%)	1.03	38/17697 (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 (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	84	ARG	NE-CZ	7.15	1.42	1.33
1	A	53	LEU	C-O	-5.28	1.13	1.23
1	A	50	GLU	CD-OE2	-5.26	1.19	1.25
1	A	62	GLU	CD-OE2	-5.09	1.20	1.25

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	63	ARG	NE-CZ-NH2	-10.37	115.12	120.30
1	A	115	ASP	CB-CA-C	9.43	129.25	110.40
1	A	63	ARG	NE-CZ-NH1	9.12	124.86	120.30
1	A	115	ASP	N-CA-CB	-8.66	95.01	110.60
1	B	84	ARG	NE-CZ-NH2	8.65	124.63	120.30
1	A	227	ASP	CB-CG-OD1	8.35	125.81	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	54	ARG	NE-CZ-NH2	-8.26	116.17	120.30
2	C	54	ARG	CG-CD-NE	8.19	129.00	111.80
1	B	227	ASP	CB-CG-OD1	7.53	125.07	118.30
1	B	251	ASP	CB-CG-OD2	-7.52	111.53	118.30
2	D	495	ARG	CG-CD-NE	-7.32	96.42	111.80
1	B	218	MET	CG-SD-CE	-6.41	89.94	100.20
1	B	122	ASP	CB-CG-OD2	-6.33	112.61	118.30
2	D	58	ARG	NE-CZ-NH2	-6.30	117.15	120.30
2	D	421	ASP	CB-CG-OD1	6.29	123.96	118.30
2	D	630	ARG	NE-CZ-NH2	-6.28	117.16	120.30
1	A	122	ASP	CB-CG-OD2	-6.25	112.67	118.30
2	D	521	ARG	NE-CZ-NH2	-6.20	117.20	120.30
2	C	201	ARG	NE-CZ-NH1	6.15	123.38	120.30
1	A	227	ASP	CB-CG-OD2	-6.10	112.81	118.30
1	A	168	ARG	NE-CZ-NH1	6.00	123.30	120.30
2	D	54	ARG	NE-CZ-NH2	-5.97	117.31	120.30
2	D	201	ARG	NE-CZ-NH1	5.94	123.27	120.30
2	C	421	ASP	CB-CG-OD1	5.81	123.53	118.30
1	B	122	ASP	CB-CG-OD1	5.76	123.49	118.30
2	D	495	ARG	CD-NE-CZ	5.61	131.45	123.60
1	B	74	TYR	CB-CG-CD2	-5.54	117.68	121.00
2	C	495	ARG	NE-CZ-NH2	-5.51	117.54	120.30
2	D	672	ALA	N-CA-C	-5.43	96.34	111.00
1	A	122	ASP	CB-CG-OD1	5.41	123.17	118.30
1	A	118	TYR	CB-CG-CD1	5.38	124.23	121.00
2	C	630	ARG	NE-CZ-NH2	-5.37	117.61	120.30
1	A	147	ARG	NE-CZ-NH1	5.30	122.95	120.30
2	C	513	ARG	NE-CZ-NH2	-5.26	117.67	120.30
2	D	495	ARG	CB-CG-CD	5.21	125.14	111.60
2	C	466	ARG	NE-CZ-NH2	-5.09	117.76	120.30
2	D	220	ARG	NE-CZ-NH1	-5.06	117.77	120.30
2	D	122	ARG	NE-CZ-NH2	-5.04	117.78	120.30

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 (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2071	0	2007	24	0
1	B	2097	0	2032	25	0
2	C	4348	0	4324	22	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	87	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 (87) 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	1.32
1:A:30:GLU:CD	1:B:28:LYS:HZ1	1.36	1.23
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	0.91
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:GLU:CD	1:B:28:LYS:CE	2.61	0.69
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
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
2:D:497:GLU:H	2:D:497:GLU:CD	1.99	0.64
1:A:30:GLU:CG	1:B:28:LYS:HZ2	2.08	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:539:THR:O	2:C:543:PHE:HE2	1.84	0.60
2:C:539:THR:O	2:C:543:PHE:CD2	2.56	0.58
2:C:356:LYS:C	2:C:385:VAL:HG12	2.25	0.57
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: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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:535:GLU:CD	2:C:596:ASN:HD21	2.15	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:D:495:ARG:C	2:D:497:GLU:OE1	2.53	0.45
2:C:466:ARG:O	2:C:466:ARG:HD2	2.16	0.45
1:A:30:GLU:CG	1:B:28:LYS:HE3	2.44	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
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
1:B:126:GLU:OE1	2:D:490:TYR:OH	2.30	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: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:D:343:ARG:HE	2:D:343:ARG:C	2.25	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%)	47	73
2	D	534/688 (78%)	511 (96%)	21 (4%)	2 (0%)	34	60
All	All	1582/1968 (80%)	1525 (96%)	54 (3%)	3 (0%)	47	73

All (3) Ramachandran outliers are listed below:

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

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	217/253 (86%)	213 (98%)	4 (2%)	59	83
1	B	220/253 (87%)	210 (96%)	10 (4%)	27	55
2	C	471/588 (80%)	439 (93%)	32 (7%)	16	36
2	D	469/588 (80%)	443 (94%)	26 (6%)	21	46
All	All	1377/1682 (82%)	1305 (95%)	72 (5%)	23	49

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	SER
1	A	30	GLU
1	A	63	ARG
1	A	271	VAL
1	B	28	LYS
1	B	55	SER
1	B	84	ARG
1	B	94	ARG

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Mol	Chain	Res	Type
1	B	129	PRO
1	B	149	LYS
1	B	182	LYS
1	B	189	LYS
1	B	271	VAL
1	B	278	LEU
2	C	36	LYS
2	C	109	LEU
2	C	131	LEU
2	C	143	ASN
2	C	152	LYS
2	C	169	LYS
2	C	176	SER
2	C	182	LYS
2	C	185	LYS
2	C	218	PHE
2	C	232	ARG
2	C	343	ARG
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	495	ARG
2	C	536	GLU
2	C	573	LYS
2	C	589	MET
2	C	594	ARG
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	137	PRO
2	D	144	ASP
2	D	176	SER
2	D	185	LYS

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Mol	Chain	Res	Type
2	D	193	ARG
2	D	218	PHE
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	386	SER
2	D	437	PHE
2	D	482	GLN
2	D	497	GLU
2	D	532	THR
2	D	538	GLU
2	D	539	THR
2	D	542	LYS
2	D	581	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 (7) such sidechains are listed below:

Mol	Chain	Res	Type
2	C	204	GLN
2	C	481	GLN
2	C	586	GLN
2	C	603	ASN
2	D	31	GLN
2	D	395	ASN
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 (i)

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

5.5 Carbohydrates (i)

There are no monosaccharides 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, 95^{th} 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%)	-0.42	0 100 100	22, 36, 64, 81	0
1	B	259/296 (87%)	-0.45	2 (0%) 86 87	23, 36, 65, 103	0
2	C	541/688 (78%)	-0.09	21 (3%) 39 38	31, 55, 101, 130	0
2	D	538/688 (78%)	-0.11	17 (3%) 47 48	35, 57, 98, 141	0
All	All	1594/1968 (80%)	-0.21	40 (2%) 57 59	22, 51, 94, 141	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	232	ARG	4.0
2	D	414	GLY	3.8
1	B	286	ASP	3.8
2	C	674	PRO	3.8
2	C	675	ALA	3.7
2	C	539	THR	3.5
2	D	539	THR	3.5
2	C	232	ARG	3.3
2	D	538	GLU	3.3
2	D	536	GLU	3.2
2	C	538	GLU	3.2
2	C	536	GLU	3.0
2	D	537	THR	3.0
2	D	598	MET	3.0
2	C	673	ARG	2.9
2	C	600	ASP	2.9
2	D	672	ALA	2.8
2	C	439	LYS	2.8
2	C	598	MET	2.8
2	D	534	ASN	2.7
2	C	356	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
2	C	672	ALA	2.4
2	D	432	THR	2.3
2	C	534	ASN	2.3
2	C	434	GLY	2.3
2	C	201	ARG	2.3
2	C	566	ALA	2.3
2	D	356	LYS	2.3
2	C	594	ARG	2.2
2	C	643	ASP	2.2
2	D	533	GLY	2.2
2	D	643	ASP	2.2
2	C	601	LYS	2.2
2	D	193	ARG	2.2
2	C	596	ASN	2.1
2	D	382	LYS	2.1
2	C	354	LYS	2.1
1	B	56	ARG	2.1
2	D	201	ARG	2.1
2	D	185	LYS	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides 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.