Full wwPDB X-ray Structure Validation Report (i)

Jun 17, 2024 – 09:41 PM EDT

PDB ID	:	5XFL
Title	:	Crystal structure of the force-sensing device region of alpha N-catenin
Authors	:	Hirano, Y.; Hakoshima, T.
Deposited on	:	2017-04-10
Resolution	:	2.45 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 (i) 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 (1)) were used in the production of this report:

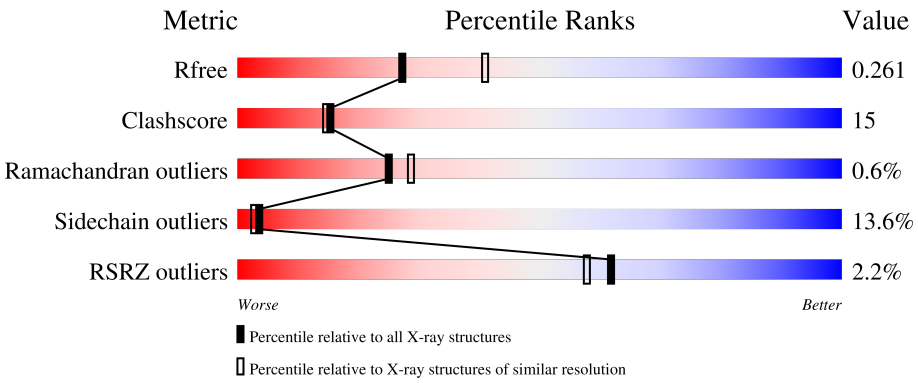
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.45 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	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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	375	69% 22% • 6%
1	B	375	70% 19% 5% 5%
1	C	375	2% 68% 23% 5% • •
1	D	375	6% 63% 26% 5% • 5%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 11346 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 Catenin alpha-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	354	2742	1700	482	545	15	0	0	0
1	B	356	2747	1701	487	544	15	0	0	0
1	C	359	2794	1731	491	557	15	0	0	0
1	D	356	2748	1699	487	547	15	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	258	GLY	-	expression tag	UNP Q61301
A	259	PRO	-	expression tag	UNP Q61301
B	258	GLY	-	expression tag	UNP Q61301
B	259	PRO	-	expression tag	UNP Q61301
C	258	GLY	-	expression tag	UNP Q61301
C	259	PRO	-	expression tag	UNP Q61301
D	258	GLY	-	expression tag	UNP Q61301
D	259	PRO	-	expression tag	UNP Q61301

- Molecule 2 is water.

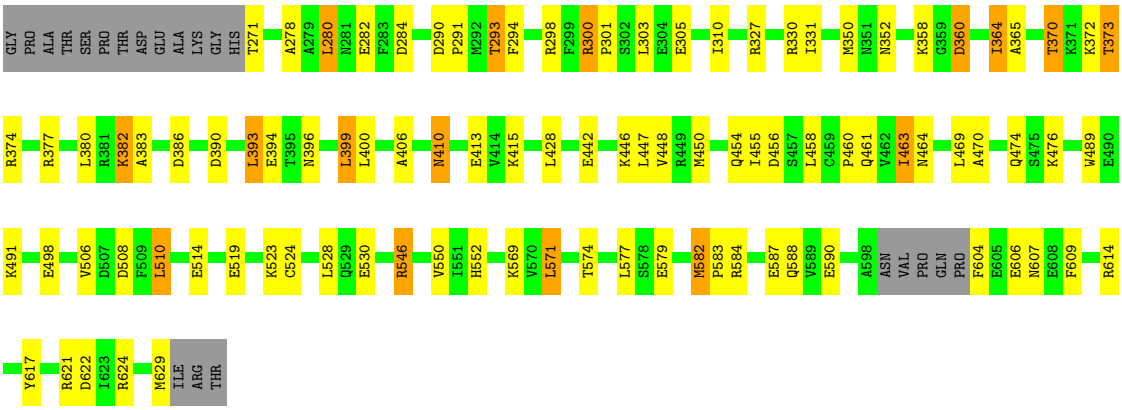
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	99	99	99	0	0
2	B	105	105	105	0	0
2	C	85	85	85	0	0
2	D	26	26	26	0	0

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.

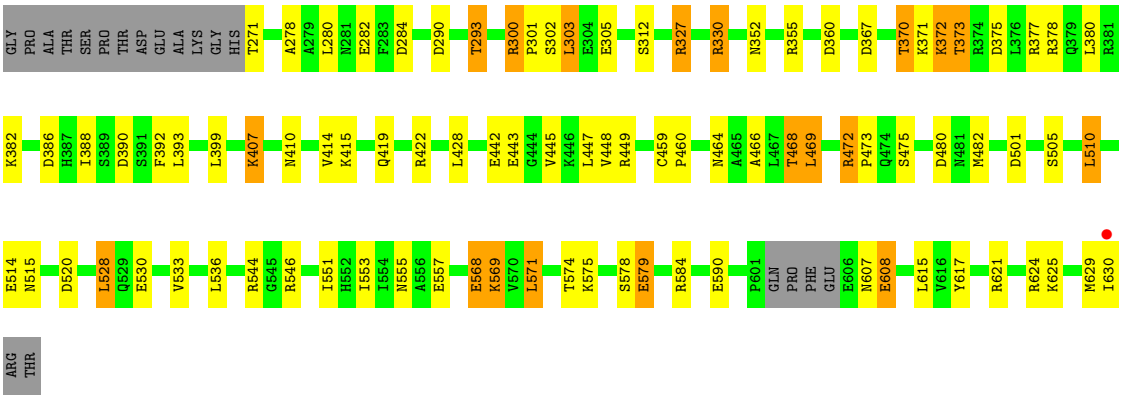
• Molecule 1: Catenin alpha-2

Chain A: 69% 22% • 6%



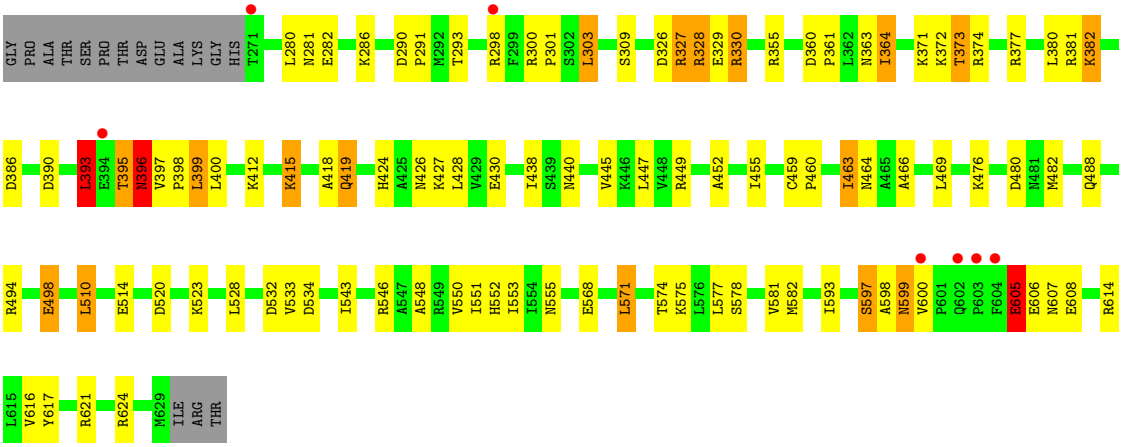
• Molecule 1: Catenin alpha-2

Chain B: 70% 19% 5% 5%



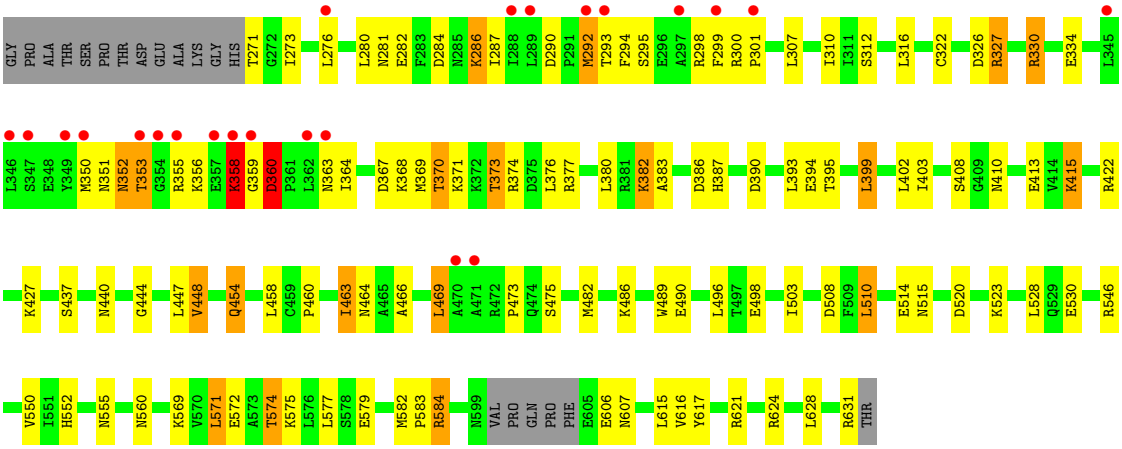
• Molecule 1: Catenin alpha-2

Chain C: 2% 68% 23% 5% • •



- Molecule 1: Catenin alpha-2

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	131.73Å 52.21Å 154.31Å 90.00° 95.05° 90.00°	Depositor
Resolution (Å)	50.00 – 2.45 43.74 – 2.45	Depositor EDS
% Data completeness (in resolution range)	97.8 (50.00-2.45) 97.8 (43.74-2.45)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	3.26 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.224 , 0.267 0.220 , 0.261	Depositor DCC
R_{free} test set	3846 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å^2)	39.3	Xtriage
Anisotropy	0.207	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 38.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	11346	wwPDB-VP
Average B, all atoms (Å^2)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.10% of the height of the origin peak. No significant pseudotranslation is detected.*

[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 (i)

5.1 Standard geometry (i)

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	0.56	0/2768	0.70	3/3739 (0.1%)
1	B	0.54	0/2773	0.64	0/3747
1	C	0.51	0/2823	0.61	0/3816
1	D	0.43	0/2773	0.55	1/3749 (0.0%)
All	All	0.51	0/11137	0.63	4/15051 (0.0%)

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	1
1	C	0	3
All	All	0	4

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	546	ARG	NE-CZ-NH2	-5.93	117.34	120.30
1	D	330	ARG	NE-CZ-NH1	-5.29	117.66	120.30
1	A	582	MET	CG-SD-CE	-5.23	91.83	100.20
1	A	622	ASP	CB-CG-OD1	5.20	122.98	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	350	MET	Peptide
1	C	393	LEU	Peptide
1	C	395	THR	Peptide

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Mol	Chain	Res	Type	Group
1	C	396	ASN	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	2742	0	2731	73	0
1	B	2747	0	2739	87	0
1	C	2794	0	2796	95	0
1	D	2748	0	2734	97	0
2	A	99	0	0	4	0
2	B	105	0	0	10	0
2	C	85	0	0	1	0
2	D	26	0	0	1	0
All	All	11346	0	11000	328	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 15.

All (328) 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:466:ALA:HA	1:B:482:MET:CE	1.66	1.23
1:C:466:ALA:CA	1:C:482:MET:HE1	1.79	1.12
1:B:330:ARG:HG2	1:B:330:ARG:HH11	1.12	1.08
1:C:466:ALA:HA	1:C:482:MET:HE1	1.23	1.08
1:C:466:ALA:HA	1:C:482:MET:CE	1.83	1.08
1:D:300:ARG:HB3	1:D:301:PRO:HD3	1.37	1.06
1:C:605:GLU:OE1	1:C:605:GLU:HA	1.48	1.06
1:C:419:GLN:HE21	1:C:419:GLN:HA	1.17	1.04
1:B:330:ARG:HH11	1:B:330:ARG:CG	1.73	1.02
1:B:466:ALA:HA	1:B:482:MET:HE1	1.01	1.01
1:B:466:ALA:CA	1:B:482:MET:HE1	1.91	1.00
1:A:372:LYS:HE2	1:A:372:LYS:HA	1.39	1.00
1:C:281:ASN:HD21	1:C:440:ASN:HD21	1.06	0.99
1:C:555:ASN:HD21	1:C:574:THR:HG21	1.24	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:419:GLN:HA	1:C:419:GLN:NE2	1.71	0.98
1:B:327:ARG:HH22	1:B:515:ASN:ND2	1.60	0.97
1:D:358:LYS:HD2	1:D:363:ASN:HD21	1.28	0.95
1:A:373:THR:HG23	1:A:377:ARG:HH12	1.32	0.94
1:B:555:ASN:HD21	1:B:574:THR:HG21	1.33	0.92
1:C:605:GLU:OE1	1:C:605:GLU:CA	2.16	0.92
1:D:370:THR:O	1:D:373:THR:HG22	1.69	0.90
1:C:281:ASN:HD21	1:C:440:ASN:ND2	1.70	0.90
1:C:372:LYS:HE2	1:C:372:LYS:HA	1.52	0.90
1:C:328:ARG:HH11	1:C:328:ARG:HG2	1.37	0.88
1:A:607:ASN:HD22	1:C:621:ARG:HH22	1.19	0.88
1:D:577:LEU:HD11	1:D:582:MET:HE1	1.55	0.88
1:A:460:PRO:O	1:A:463:ILE:HG22	1.74	0.87
1:A:607:ASN:ND2	1:C:621:ARG:HH22	1.74	0.86
1:C:551:ILE:HD11	1:C:577:LEU:HD23	1.55	0.86
1:C:373:THR:CG2	1:C:377:ARG:HH12	1.90	0.85
1:B:466:ALA:CA	1:B:482:MET:CE	2.53	0.84
1:D:577:LEU:HD11	1:D:582:MET:CE	2.09	0.83
1:A:372:LYS:HA	1:A:372:LYS:CE	2.09	0.82
1:D:555:ASN:HD21	1:D:574:THR:HG21	1.43	0.82
1:C:494:ARG:O	1:C:498:GLU:HG2	1.81	0.80
1:D:460:PRO:O	1:D:463:ILE:HG22	1.81	0.80
1:A:373:THR:HG23	1:A:377:ARG:NH1	1.96	0.80
1:D:577:LEU:CD1	1:D:582:MET:HE3	2.11	0.79
1:A:514:GLU:OE1	1:A:624:ARG:HD3	1.82	0.79
1:C:328:ARG:HG2	1:C:328:ARG:NH1	1.94	0.79
1:B:514:GLU:OE1	1:B:624:ARG:HD3	1.83	0.78
1:D:510:LEU:HG	1:D:624:ARG:HB2	1.66	0.78
1:C:396:ASN:O	1:C:400:LEU:HD13	1.84	0.77
1:B:284:ASP:OD1	1:B:373:THR:HG21	1.84	0.77
1:C:466:ALA:N	1:C:482:MET:HE1	1.98	0.77
1:D:370:THR:HG23	1:D:374:ARG:HH22	1.50	0.77
1:C:373:THR:CG2	1:C:377:ARG:NH1	2.48	0.77
1:D:577:LEU:CD1	1:D:582:MET:CE	2.64	0.76
1:A:607:ASN:HD22	1:C:621:ARG:NH2	1.84	0.76
1:D:617:TYR:CE2	1:D:621:ARG:HD2	2.21	0.76
1:A:373:THR:CG2	1:A:377:ARG:HH12	1.99	0.76
1:A:390:ASP:OD1	1:A:546:ARG:NH2	2.19	0.76
1:B:372:LYS:HE3	1:B:372:LYS:HA	1.68	0.76
1:B:551:ILE:HD13	1:B:578:SER:HB2	1.68	0.76
1:A:607:ASN:ND2	1:C:621:ARG:NH2	2.33	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:373:THR:HG22	1:C:377:ARG:NH1	2.01	0.76
1:C:555:ASN:ND2	1:C:574:THR:HG21	2.01	0.75
1:B:330:ARG:HG2	1:B:330:ARG:NH1	1.94	0.75
1:C:396:ASN:O	1:C:400:LEU:CD1	2.35	0.75
1:C:581:VAL:HG12	1:C:582:MET:HE1	1.66	0.75
1:A:571:LEU:HA	1:A:574:THR:HG22	1.68	0.74
1:B:630:ILE:C	2:B:705:HOH:O	2.25	0.74
1:B:373:THR:HG23	1:B:377:ARG:HH12	1.52	0.74
1:B:382:LYS:CE	2:B:722:HOH:O	2.36	0.74
1:D:514:GLU:OE1	1:D:624:ARG:HD3	1.88	0.74
1:B:373:THR:HG23	1:B:377:ARG:NH1	2.02	0.74
1:B:617:TYR:CZ	1:B:621:ARG:HD2	2.24	0.73
1:A:291:PRO:HG3	1:A:358:LYS:HG2	1.71	0.73
1:C:328:ARG:HG3	1:C:329:GLU:N	2.04	0.73
1:B:555:ASN:ND2	1:B:574:THR:HG21	2.04	0.73
1:C:327:ARG:HD2	1:C:386:ASP:OD2	1.89	0.73
1:C:514:GLU:OE1	1:C:624:ARG:HD3	1.89	0.72
1:A:583:PRO:HG3	1:B:472:ARG:NH1	2.04	0.72
1:D:367:ASP:O	1:D:370:THR:HG22	1.90	0.71
1:D:466:ALA:HA	1:D:482:MET:CE	2.21	0.71
1:D:584:ARG:HH11	1:D:615:LEU:HD12	1.56	0.70
1:B:382:LYS:HE3	2:B:722:HOH:O	1.90	0.70
1:B:617:TYR:CE2	1:B:621:ARG:HD2	2.27	0.70
1:C:396:ASN:HB3	1:C:399:LEU:HB2	1.72	0.70
1:A:510:LEU:HG	1:A:624:ARG:HB2	1.73	0.70
1:D:444:GLY:O	1:D:448:VAL:HG12	1.91	0.70
1:C:373:THR:HG22	1:C:377:ARG:HH12	1.57	0.69
1:B:621:ARG:NH2	1:D:607:ASN:HD22	1.90	0.69
1:B:466:ALA:HA	1:B:482:MET:HE3	1.70	0.69
1:B:367:ASP:O	1:B:370:THR:HG22	1.93	0.69
1:C:555:ASN:HD21	1:C:574:THR:CG2	2.02	0.69
1:C:460:PRO:O	1:C:463:ILE:HG22	1.92	0.68
1:D:334:GLU:HA	1:D:334:GLU:OE2	1.93	0.68
1:D:617:TYR:CZ	1:D:621:ARG:HD2	2.28	0.68
1:A:583:PRO:HG3	1:B:472:ARG:HH12	1.59	0.67
1:D:370:THR:HG23	1:D:374:ARG:NH2	2.08	0.67
1:D:358:LYS:HD2	1:D:363:ASN:ND2	2.07	0.67
1:A:617:TYR:CE2	1:A:621:ARG:HD2	2.29	0.67
1:B:330:ARG:CG	1:B:330:ARG:NH1	2.44	0.66
1:C:390:ASP:OD1	1:C:546:ARG:NH2	2.26	0.66
1:D:382:LYS:HE3	1:D:508:ASP:OD1	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:607:ASN:OD1	1:D:621:ARG:NH2	2.28	0.66
1:C:571:LEU:O	1:C:574:THR:HG22	1.96	0.66
1:D:466:ALA:HA	1:D:482:MET:HE2	1.78	0.66
1:D:327:ARG:HH22	1:D:515:ASN:HD22	1.44	0.66
1:A:624:ARG:NH2	1:C:606:GLU:OE1	2.28	0.65
1:D:373:THR:HG23	1:D:377:ARG:HH12	1.61	0.65
1:A:552:HIS:HD2	1:B:480:ASP:OD1	1.80	0.65
1:B:327:ARG:HH22	1:B:515:ASN:HD22	1.41	0.65
1:B:372:LYS:HA	1:B:372:LYS:CE	2.27	0.65
1:B:621:ARG:HH22	1:D:607:ASN:HD22	1.45	0.65
1:C:571:LEU:HA	1:C:574:THR:HG22	1.79	0.65
1:C:551:ILE:CD1	1:C:577:LEU:HD23	2.27	0.64
1:C:617:TYR:CZ	1:C:621:ARG:HD2	2.31	0.64
1:C:326:ASP:OD2	1:C:330:ARG:HD2	1.96	0.64
1:A:284:ASP:OD1	1:A:373:THR:HG21	1.98	0.64
1:B:472:ARG:HG2	1:B:475:SER:HB2	1.78	0.64
1:D:577:LEU:HD12	1:D:582:MET:HE3	1.79	0.64
1:D:582:MET:CE	1:D:616:VAL:HG22	2.28	0.64
1:B:371:LYS:NZ	1:B:375:ASP:OD2	2.27	0.64
1:C:327:ARG:HD3	1:C:382:LYS:HB3	1.80	0.64
1:B:327:ARG:HD2	1:B:386:ASP:OD2	1.98	0.64
1:A:461:GLN:HA	1:A:464:ASN:HD22	1.63	0.63
1:B:621:ARG:NH2	1:D:607:ASN:ND2	2.46	0.63
1:C:463:ILE:CG2	1:C:464:ASN:N	2.61	0.63
1:D:373:THR:HG23	1:D:377:ARG:NH1	2.14	0.63
1:A:577:LEU:HG	1:A:582:MET:CE	2.29	0.62
1:C:396:ASN:HA	1:C:398:PRO:HD2	1.82	0.62
1:D:454:GLN:O	1:D:458:LEU:HB2	2.00	0.62
1:D:631:ARG:HG2	1:D:631:ARG:HH11	1.65	0.62
1:D:327:ARG:HH22	1:D:515:ASN:ND2	1.97	0.62
1:A:617:TYR:CZ	1:A:621:ARG:HD2	2.35	0.61
1:D:300:ARG:CB	1:D:301:PRO:HD3	2.23	0.61
1:A:330:ARG:NH2	1:C:532:ASP:OD2	2.33	0.61
1:D:300:ARG:HB3	1:D:301:PRO:CD	2.20	0.61
1:D:582:MET:HE2	1:D:616:VAL:HG22	1.82	0.61
1:D:584:ARG:NH1	1:D:615:LEU:HD12	2.15	0.61
1:A:271:THR:HA	1:A:442:GLU:OE2	2.01	0.60
1:D:390:ASP:OD1	1:D:546:ARG:NH2	2.34	0.60
1:A:280:LEU:HD13	1:A:310:ILE:HD13	1.83	0.60
1:A:327:ARG:HD2	1:A:386:ASP:OD2	2.01	0.60
1:A:370:THR:CG2	1:A:374:ARG:NH1	2.64	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:300:ARG:HB3	1:C:301:PRO:HD3	1.84	0.60
1:D:293:THR:O	1:D:293:THR:HG23	2.02	0.60
1:A:382:LYS:HE3	1:A:508:ASP:OD1	2.02	0.60
1:B:571:LEU:HA	1:B:574:THR:HG22	1.82	0.60
1:D:322:CYS:HB2	1:D:387:HIS:CD2	2.37	0.60
1:D:292:MET:HB2	1:D:356:LYS:HA	1.82	0.59
1:C:617:TYR:CE2	1:C:621:ARG:HD2	2.38	0.59
1:A:571:LEU:O	1:A:574:THR:HG22	2.02	0.59
1:B:327:ARG:HH22	1:B:515:ASN:HD21	1.49	0.58
1:A:624:ARG:HH22	1:C:606:GLU:CD	2.06	0.58
1:A:278:ALA:O	1:A:282:GLU:HG2	2.04	0.58
1:C:419:GLN:HE21	1:C:419:GLN:CA	1.93	0.58
1:C:419:GLN:NE2	1:C:419:GLN:CA	2.49	0.58
1:B:621:ARG:HH22	1:D:607:ASN:ND2	2.01	0.57
1:D:295:SER:O	1:D:299:PHE:HB2	2.04	0.57
1:A:571:LEU:HD12	1:A:574:THR:HG21	1.86	0.57
1:B:419:GLN:HA	1:B:419:GLN:NE2	2.17	0.57
1:D:448:VAL:HG11	1:D:503:ILE:HD13	1.86	0.57
1:D:575:LYS:O	1:D:579:GLU:HB2	2.04	0.57
1:B:443:GLU:CD	1:B:443:GLU:H	2.08	0.57
1:B:544:ARG:HD2	2:B:768:HOH:O	2.05	0.57
1:B:300:ARG:HB3	1:B:301:PRO:HD3	1.88	0.56
1:B:382:LYS:HE2	2:B:722:HOH:O	1.98	0.56
1:B:390:ASP:OD1	1:B:546:ARG:NH2	2.36	0.56
1:C:463:ILE:HG22	1:C:464:ASN:N	2.22	0.55
1:D:399:LEU:O	1:D:399:LEU:HD22	2.06	0.55
1:D:498:GLU:OE2	1:D:552:HIS:HE1	1.89	0.55
1:A:364:ILE:HG13	1:A:365:ALA:N	2.22	0.55
1:A:571:LEU:HA	1:A:574:THR:CG2	2.35	0.54
1:B:575:LYS:HG3	1:B:579:GLU:OE2	2.07	0.54
1:A:577:LEU:HG	1:A:582:MET:HE1	1.88	0.53
1:C:397:VAL:N	1:C:398:PRO:CD	2.71	0.53
1:A:372:LYS:HE2	1:A:372:LYS:CA	2.15	0.53
1:D:466:ALA:HA	1:D:482:MET:HE1	1.91	0.53
1:D:469:LEU:HD22	1:D:473:PRO:HA	1.89	0.53
1:B:608:GLU:HG2	2:B:726:HOH:O	2.08	0.53
1:A:454:GLN:O	1:A:458:LEU:HB2	2.07	0.53
1:A:360:ASP:O	1:A:364:ILE:HG23	2.09	0.53
1:B:466:ALA:CB	1:B:482:MET:HE3	2.39	0.53
1:B:290:ASP:O	1:B:293:THR:HG22	2.09	0.53
1:B:555:ASN:HD21	1:B:574:THR:CG2	2.13	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:358:LYS:O	1:D:358:LYS:HG3	2.09	0.52
1:B:625:LYS:NZ	1:D:607:ASN:HD21	2.08	0.52
1:C:281:ASN:ND2	1:C:440:ASN:HD21	1.90	0.52
1:D:327:ARG:HD3	1:D:382:LYS:HB3	1.91	0.52
1:B:520:ASP:OD2	1:B:546:ARG:NH1	2.41	0.51
1:A:372:LYS:CE	1:A:372:LYS:CA	2.82	0.51
1:C:605:GLU:OE1	1:C:605:GLU:N	2.42	0.51
1:D:383:ALA:O	1:D:386:ASP:HB2	2.10	0.51
1:A:327:ARG:HD3	1:A:382:LYS:HB3	1.93	0.51
1:B:501:ASP:CG	1:B:553:ILE:HD11	2.31	0.50
1:B:568:GLU:OE2	1:B:568:GLU:HA	2.10	0.50
1:C:330:ARG:NH1	2:C:703:HOH:O	2.44	0.50
1:A:410:ASN:CB	1:A:413:GLU:HG3	2.42	0.50
1:D:371:LYS:O	1:D:374:ARG:HB3	2.11	0.50
1:C:360:ASP:HB2	1:C:361:PRO:HD2	1.94	0.50
1:A:577:LEU:HG	1:A:582:MET:HE2	1.93	0.50
1:D:282:GLU:O	1:D:286:LYS:HB2	2.11	0.50
1:C:328:ARG:HH11	1:C:328:ARG:CG	2.14	0.49
1:D:284:ASP:OD1	1:D:373:THR:HG21	2.12	0.49
1:D:399:LEU:HD22	1:D:403:ILE:HG13	1.93	0.49
1:A:524:CYS:HB3	1:A:609:PHE:CZ	2.47	0.49
1:D:498:GLU:OE2	1:D:552:HIS:CE1	2.66	0.49
1:B:378:ARG:HD3	2:B:716:HOH:O	2.11	0.49
1:C:290:ASP:HB3	1:C:293:THR:HG23	1.95	0.49
1:C:598:ALA:C	1:C:600:VAL:H	2.16	0.49
1:B:624:ARG:NH2	1:D:606:GLU:OE1	2.45	0.49
1:D:326:ASP:OD2	1:D:330:ARG:NH1	2.45	0.49
1:A:571:LEU:CA	1:A:574:THR:HG22	2.39	0.49
1:D:327:ARG:HD2	1:D:386:ASP:OD2	2.12	0.49
1:B:278:ALA:O	1:B:282:GLU:HG2	2.12	0.48
1:C:571:LEU:HA	1:C:574:THR:CG2	2.43	0.48
1:A:300:ARG:HB3	1:A:301:PRO:HD3	1.96	0.48
1:C:396:ASN:HD21	1:C:424:HIS:CD2	2.32	0.48
1:D:355:ARG:HG2	1:D:355:ARG:HH11	1.79	0.48
1:B:510:LEU:HG	1:B:624:ARG:HB2	1.96	0.48
1:A:282:GLU:OE2	2:A:701:HOH:O	2.20	0.48
1:B:271:THR:HA	1:B:442:GLU:OE2	2.14	0.47
1:B:382:LYS:NZ	1:D:530:GLU:OE1	2.48	0.47
1:A:294:PHE:CD1	1:A:294:PHE:C	2.87	0.47
1:A:552:HIS:CD2	1:B:480:ASP:OD1	2.66	0.47
1:B:557:GLU:O	1:B:557:GLU:HG3	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:520:ASP:HB3	1:C:543:ILE:HG13	1.97	0.47
1:C:548:ALA:O	1:C:552:HIS:CD2	2.68	0.47
1:D:352:ASN:CG	1:D:352:ASN:O	2.53	0.47
1:B:528:LEU:CD1	1:B:536:LEU:HD22	2.45	0.46
1:D:322:CYS:HB2	1:D:387:HIS:CG	2.51	0.46
1:C:381:ARG:HG2	1:C:438:ILE:HD12	1.98	0.46
1:B:372:LYS:HE3	1:B:375:ASP:OD2	2.14	0.46
1:B:571:LEU:O	1:B:574:THR:HG22	2.16	0.46
1:B:422:ARG:HH22	1:B:460:PRO:HB3	1.80	0.46
1:B:629:MET:O	1:B:630:ILE:CB	2.63	0.46
1:A:461:GLN:HA	1:A:464:ASN:ND2	2.30	0.46
1:B:466:ALA:CA	1:B:482:MET:HE3	2.37	0.46
1:B:459:CYS:HB3	1:B:460:PRO:CD	2.46	0.45
1:B:533:VAL:HG22	2:B:738:HOH:O	2.15	0.45
1:A:614:ARG:HD3	1:C:614:ARG:HD3	1.97	0.45
1:D:359:GLY:O	1:D:360:ASP:C	2.55	0.45
1:B:464:ASN:O	1:B:468:THR:HG23	2.16	0.45
1:C:412:LYS:O	1:C:415:LYS:HB2	2.17	0.45
1:D:294:PHE:HD1	1:D:353:THR:HG1	1.63	0.45
1:C:571:LEU:CA	1:C:574:THR:HG22	2.46	0.45
1:D:280:LEU:HB3	1:D:377:ARG:NH2	2.32	0.45
1:B:370:THR:CG2	2:B:753:HOH:O	2.64	0.45
1:C:593:ILE:O	1:C:597:SER:HB2	2.16	0.45
1:B:469:LEU:HD22	1:B:473:PRO:HA	1.98	0.45
1:C:361:PRO:HA	1:C:364:ILE:CG2	2.47	0.45
1:C:459:CYS:N	1:C:460:PRO:HD2	2.32	0.45
1:A:519:GLU:OE1	2:A:702:HOH:O	2.21	0.45
1:B:575:LYS:O	1:B:579:GLU:HB2	2.16	0.45
1:D:577:LEU:HD12	1:D:582:MET:CE	2.43	0.45
1:A:587:GLU:OE1	1:A:587:GLU:HA	2.16	0.45
1:C:533:VAL:HG23	1:C:534:ASP:N	2.32	0.44
1:D:402:LEU:O	1:D:402:LEU:HG	2.16	0.44
1:C:396:ASN:HB3	1:C:399:LEU:H	1.81	0.44
1:C:361:PRO:O	1:C:364:ILE:HG23	2.17	0.44
1:C:393:LEU:HD12	1:C:393:LEU:HA	1.63	0.44
1:C:396:ASN:CB	1:C:399:LEU:HB2	2.42	0.44
1:A:455:ILE:HG23	1:A:489:TRP:CH2	2.53	0.44
1:B:407:LYS:HZ2	1:B:407:LYS:HB3	1.82	0.44
1:C:463:ILE:HG22	1:C:464:ASN:H	1.82	0.44
1:A:498:GLU:OE2	1:A:552:HIS:HE1	2.01	0.44
1:D:290:ASP:HB3	1:D:293:THR:HG22	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:569:LYS:HB2	1:B:569:LYS:HE2	1.63	0.44
1:B:624:ARG:HH22	1:D:606:GLU:CD	2.22	0.44
1:D:290:ASP:CG	1:D:293:THR:HG22	2.38	0.43
1:D:292:MET:HA	1:D:355:ARG:O	2.18	0.43
1:A:293:THR:HG22	2:A:752:HOH:O	2.19	0.43
1:D:395:THR:HG22	2:D:707:HOH:O	2.17	0.43
1:A:506:VAL:O	1:A:510:LEU:HD22	2.18	0.43
1:A:327:ARG:NH1	1:A:386:ASP:OD1	2.51	0.43
1:B:373:THR:HG22	2:B:713:HOH:O	2.19	0.43
1:C:303:LEU:HD12	1:C:303:LEU:HA	1.88	0.43
1:C:510:LEU:HG	1:C:624:ARG:HB2	1.99	0.43
1:C:577:LEU:HD11	1:C:582:MET:HE3	1.99	0.43
1:D:520:ASP:OD2	1:D:546:ARG:NH1	2.46	0.43
1:A:393:LEU:O	1:A:394:GLU:C	2.57	0.43
1:C:494:ARG:O	1:C:498:GLU:CG	2.59	0.43
1:C:452:ALA:O	1:C:455:ILE:HB	2.19	0.43
1:B:373:THR:CG2	1:B:377:ARG:HH12	2.27	0.43
1:D:394:GLU:OE2	1:D:394:GLU:HA	2.19	0.43
1:A:293:THR:CG2	2:A:752:HOH:O	2.66	0.42
1:D:463:ILE:CG2	1:D:464:ASN:N	2.82	0.42
1:D:287:ILE:HD13	1:D:369:MET:HG2	2.01	0.42
1:D:489:TRP:O	1:D:490:GLU:C	2.57	0.42
1:A:474:GLN:H	1:A:474:GLN:CD	2.23	0.42
1:A:552:HIS:HD2	1:B:480:ASP:CG	2.23	0.42
1:B:625:LYS:HZ1	1:D:607:ASN:HD21	1.66	0.42
1:C:520:ASP:OD2	1:C:546:ARG:NH1	2.51	0.42
1:C:327:ARG:CD	1:C:382:LYS:HB3	2.48	0.42
1:C:396:ASN:HB3	1:C:399:LEU:CB	2.47	0.42
1:C:520:ASP:HB3	1:C:543:ILE:CG1	2.49	0.42
1:C:582:MET:HE1	1:C:616:VAL:HG22	2.01	0.42
1:D:370:THR:CG2	1:D:374:ARG:HH22	2.25	0.42
1:D:376:LEU:O	1:D:377:ARG:C	2.58	0.42
1:D:281:ASN:HD21	1:D:440:ASN:ND2	2.18	0.42
1:D:422:ARG:HH22	1:D:460:PRO:HB3	1.85	0.42
1:B:388:ILE:HG23	1:B:392:PHE:CD2	2.54	0.42
1:C:360:ASP:O	1:C:364:ILE:HG22	2.20	0.42
1:D:606:GLU:HG3	1:D:607:ASN:N	2.34	0.42
1:D:276:LEU:HB3	1:D:437:SER:O	2.20	0.41
1:C:371:LYS:O	1:C:374:ARG:HB3	2.20	0.41
1:A:446:LYS:O	1:A:450:MET:HG3	2.20	0.41
1:A:331:ILE:HD11	1:A:383:ALA:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:370:THR:HG23	1:A:374:ARG:NH1	2.36	0.41
1:D:571:LEU:HD12	1:D:571:LEU:HA	1.90	0.41
1:C:577:LEU:HG	1:C:582:MET:HE2	2.01	0.41
1:A:399:LEU:HD23	1:A:399:LEU:HA	1.90	0.41
1:A:406:ALA:HA	1:A:470:ALA:HB2	2.02	0.41
1:C:426:ASN:O	1:C:430:GLU:HG3	2.19	0.41
1:D:631:ARG:NH1	1:D:631:ARG:CG	2.83	0.41
1:B:303:LEU:HD12	1:B:303:LEU:HA	1.87	0.41
1:C:571:LEU:C	1:C:574:THR:HG22	2.41	0.41
1:A:463:ILE:CG2	1:A:464:ASN:N	2.84	0.41
1:A:621:ARG:NH2	1:C:607:ASN:OD1	2.53	0.41
1:B:327:ARG:NH2	1:B:515:ASN:ND2	2.45	0.41
1:C:418:ALA:HB1	1:C:463:ILE:CG1	2.51	0.41
1:C:445:VAL:O	1:C:449:ARG:HG3	2.21	0.41
1:D:307:LEU:O	1:D:310:ILE:HB	2.21	0.41
1:D:496:LEU:HD12	1:D:496:LEU:O	2.21	0.41
1:C:533:VAL:CG2	1:C:534:ASP:N	2.83	0.41
1:D:415:LYS:HA	1:D:415:LYS:HD3	1.82	0.41
1:A:456:ASP:O	1:A:460:PRO:HD2	2.21	0.40
1:D:631:ARG:HH11	1:D:631:ARG:CG	2.29	0.40
1:A:290:ASP:HA	1:A:291:PRO:HD3	1.91	0.40
1:B:410:ASN:O	1:B:414:VAL:HG23	2.21	0.40
1:B:445:VAL:O	1:B:449:ARG:HG3	2.21	0.40
1:B:568:GLU:OE2	1:B:568:GLU:CA	2.69	0.40
1:D:355:ARG:HG2	1:D:355:ARG:NH1	2.37	0.40
1:A:569:LYS:HE3	1:A:569:LYS:HB2	1.95	0.40
1:D:351:ASN:C	1:D:353:THR:H	2.23	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	350/375 (93%)	343 (98%)	6 (2%)	1 (0%)	41	49
1	B	352/375 (94%)	347 (99%)	5 (1%)	0	100	100
1	C	357/375 (95%)	347 (97%)	7 (2%)	3 (1%)	19	22
1	D	352/375 (94%)	333 (95%)	15 (4%)	4 (1%)	14	14
All	All	1411/1500 (94%)	1370 (97%)	33 (2%)	8 (1%)	25	29

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	358	LYS
1	C	396	ASN
1	C	599	ASN
1	D	360	ASP
1	A	410	ASN
1	C	605	GLU
1	D	410	ASN
1	D	273	ILE

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	295/319 (92%)	257 (87%)	38 (13%)	4	3
1	B	295/319 (92%)	257 (87%)	38 (13%)	4	3
1	C	305/319 (96%)	262 (86%)	43 (14%)	3	2
1	D	296/319 (93%)	253 (86%)	43 (14%)	3	2
All	All	1191/1276 (93%)	1029 (86%)	162 (14%)	3	3

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

Mol	Chain	Res	Type
1	A	280	LEU
1	A	293	THR
1	A	298	ARG

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Mol	Chain	Res	Type
1	A	300	ARG
1	A	303	LEU
1	A	305	GLU
1	A	352	ASN
1	A	360	ASP
1	A	364	ILE
1	A	370	THR
1	A	373	THR
1	A	380	LEU
1	A	382	LYS
1	A	393	LEU
1	A	396	ASN
1	A	399	LEU
1	A	400	LEU
1	A	415	LYS
1	A	428	LEU
1	A	447	LEU
1	A	448	VAL
1	A	463	ILE
1	A	469	LEU
1	A	476	LYS
1	A	491	LYS
1	A	510	LEU
1	A	523	LYS
1	A	528	LEU
1	A	530	GLU
1	A	550	VAL
1	A	571	LEU
1	A	579	GLU
1	A	584	ARG
1	A	588	GLN
1	A	590	GLU
1	A	604	PHE
1	A	606	GLU
1	A	629	MET
1	B	280	LEU
1	B	293	THR
1	B	300	ARG
1	B	302	SER
1	B	303	LEU
1	B	305	GLU
1	B	312	SER

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Mol	Chain	Res	Type
1	B	327	ARG
1	B	330	ARG
1	B	352	ASN
1	B	355	ARG
1	B	360	ASP
1	B	370	THR
1	B	372	LYS
1	B	373	THR
1	B	380	LEU
1	B	393	LEU
1	B	399	LEU
1	B	407	LYS
1	B	415	LYS
1	B	428	LEU
1	B	447	LEU
1	B	448	VAL
1	B	468	THR
1	B	469	LEU
1	B	472	ARG
1	B	505	SER
1	B	510	LEU
1	B	528	LEU
1	B	530	GLU
1	B	568	GLU
1	B	569	LYS
1	B	571	LEU
1	B	579	GLU
1	B	584	ARG
1	B	590	GLU
1	B	608	GLU
1	B	615	LEU
1	C	280	LEU
1	C	282	GLU
1	C	286	LYS
1	C	291	PRO
1	C	298	ARG
1	C	303	LEU
1	C	309	SER
1	C	327	ARG
1	C	328	ARG
1	C	330	ARG
1	C	355	ARG

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Mol	Chain	Res	Type
1	C	363	ASN
1	C	364	ILE
1	C	373	THR
1	C	380	LEU
1	C	382	LYS
1	C	393	LEU
1	C	395	THR
1	C	399	LEU
1	C	415	LYS
1	C	419	GLN
1	C	427	LYS
1	C	428	LEU
1	C	447	LEU
1	C	463	ILE
1	C	469	LEU
1	C	476	LYS
1	C	480	ASP
1	C	488	GLN
1	C	498	GLU
1	C	510	LEU
1	C	523	LYS
1	C	528	LEU
1	C	550	VAL
1	C	553	ILE
1	C	568	GLU
1	C	571	LEU
1	C	575	LYS
1	C	578	SER
1	C	597	SER
1	C	599	ASN
1	C	605	GLU
1	C	608	GLU
1	D	271	THR
1	D	286	LYS
1	D	292	MET
1	D	298	ARG
1	D	312	SER
1	D	316	LEU
1	D	327	ARG
1	D	350	MET
1	D	352	ASN
1	D	353	THR

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Mol	Chain	Res	Type
1	D	358	LYS
1	D	360	ASP
1	D	364	ILE
1	D	368	LYS
1	D	370	THR
1	D	373	THR
1	D	380	LEU
1	D	382	LYS
1	D	393	LEU
1	D	399	LEU
1	D	408	SER
1	D	413	GLU
1	D	415	LYS
1	D	427	LYS
1	D	447	LEU
1	D	448	VAL
1	D	454	GLN
1	D	463	ILE
1	D	469	LEU
1	D	475	SER
1	D	486	LYS
1	D	510	LEU
1	D	523	LYS
1	D	528	LEU
1	D	550	VAL
1	D	560	ASN
1	D	569	LYS
1	D	571	LEU
1	D	572	GLU
1	D	574	THR
1	D	583	PRO
1	D	584	ARG
1	D	628	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	351	ASN
1	A	396	ASN
1	A	419	GLN
1	A	464	ASN
1	A	481	ASN

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Mol	Chain	Res	Type
1	A	522	ASN
1	A	552	HIS
1	A	555	ASN
1	A	588	GLN
1	A	607	ASN
1	B	351	ASN
1	B	363	ASN
1	B	419	GLN
1	B	479	GLN
1	B	515	ASN
1	B	522	ASN
1	B	555	ASN
1	B	560	ASN
1	C	419	GLN
1	C	440	ASN
1	C	479	GLN
1	C	488	GLN
1	C	492	GLN
1	C	522	ASN
1	C	552	HIS
1	C	555	ASN
1	C	588	GLN
1	D	363	ASN
1	D	440	ASN
1	D	454	GLN
1	D	479	GLN
1	D	515	ASN
1	D	522	ASN
1	D	552	HIS
1	D	555	ASN
1	D	560	ASN
1	D	607	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no 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 (i)

6.1 Protein, DNA and RNA chains (i)

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	354/375 (94%)	-0.25	0	100	100	18, 36, 70, 89	0
1	B	356/375 (94%)	-0.32	1 (0%)	94	94	19, 35, 62, 88	0
1	C	359/375 (95%)	-0.13	7 (1%)	66	64	24, 43, 78, 102	0
1	D	356/375 (94%)	0.20	23 (6%)	18	15	26, 59, 128, 159	0
All	All	1425/1500 (95%)	-0.13	31 (2%)	62	58	18, 41, 88, 159	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	349	TYR	4.7
1	D	289	LEU	4.6
1	D	292	MET	4.5
1	D	299	PHE	4.3
1	D	471	ALA	4.0
1	C	600	VAL	3.8
1	D	363	ASN	3.7
1	D	358	LYS	3.2
1	D	345	LEU	3.2
1	C	602	GLN	3.1
1	D	288	ILE	3.0
1	D	293	THR	2.9
1	D	347	SER	2.7
1	C	271	THR	2.7
1	D	301	PRO	2.6
1	C	394	GLU	2.5
1	C	603	PRO	2.5
1	D	362	LEU	2.5
1	C	604	PHE	2.5
1	D	355	ARG	2.5
1	D	357	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	297	ALA	2.5
1	D	350	MET	2.3
1	D	470	ALA	2.3
1	C	298	ARG	2.2
1	D	346	LEU	2.2
1	D	276	LEU	2.1
1	D	354	GLY	2.1
1	D	353	THR	2.0
1	D	359	GLY	2.0
1	B	630	ILE	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.