Full wwPDB X-ray Structure Validation Report ⓘ

Oct 31, 2023 – 06:40 PM JST

PDB ID	:	5GT6
Title	:	Apo structure of Aldehyde Dehydrogenase from Bacillus cereus
Authors	:	Ngo, H.P.T.; Hong, S.H.; Ho, T.H.; Oh, D.K.; Kang, L.W.
Deposited on	:	2016-08-18
Resolution	:	2.60 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:

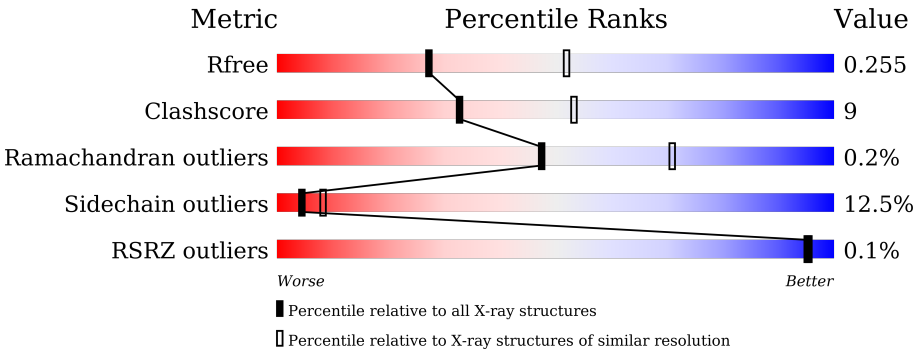
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	508	76% 18%
1	B	508	71% 19% 5%
1	C	508	77% 16%
1	D	508	75% 17%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 15213 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 Betaine-aldehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	491	Total	C	N	O	S	0	0	0
			3791	2414	630	733	14			
1	B	489	Total	C	N	O	S	0	0	0
			3776	2406	627	729	14			
1	C	487	Total	C	N	O	S	0	0	0
			3759	2395	625	725	14			
1	D	491	Total	C	N	O	S	0	0	0
			3791	2414	630	733	14			

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	expression tag	UNP A0A150BLG9
A	-12	GLY	-	expression tag	UNP A0A150BLG9
A	-11	SER	-	expression tag	UNP A0A150BLG9
A	-10	SER	-	expression tag	UNP A0A150BLG9
A	-9	HIS	-	expression tag	UNP A0A150BLG9
A	-8	HIS	-	expression tag	UNP A0A150BLG9
A	-7	HIS	-	expression tag	UNP A0A150BLG9
A	-6	HIS	-	expression tag	UNP A0A150BLG9
A	-5	HIS	-	expression tag	UNP A0A150BLG9
A	-4	HIS	-	expression tag	UNP A0A150BLG9
A	-3	SER	-	expression tag	UNP A0A150BLG9
A	-2	GLN	-	expression tag	UNP A0A150BLG9
A	-1	ASP	-	expression tag	UNP A0A150BLG9
A	0	PRO	-	expression tag	UNP A0A150BLG9
B	-13	MET	-	expression tag	UNP A0A150BLG9
B	-12	GLY	-	expression tag	UNP A0A150BLG9
B	-11	SER	-	expression tag	UNP A0A150BLG9
B	-10	SER	-	expression tag	UNP A0A150BLG9
B	-9	HIS	-	expression tag	UNP A0A150BLG9
B	-8	HIS	-	expression tag	UNP A0A150BLG9
B	-7	HIS	-	expression tag	UNP A0A150BLG9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	HIS	-	expression tag	UNP A0A150BLG9
B	-5	HIS	-	expression tag	UNP A0A150BLG9
B	-4	HIS	-	expression tag	UNP A0A150BLG9
B	-3	SER	-	expression tag	UNP A0A150BLG9
B	-2	GLN	-	expression tag	UNP A0A150BLG9
B	-1	ASP	-	expression tag	UNP A0A150BLG9
B	0	PRO	-	expression tag	UNP A0A150BLG9
C	-13	MET	-	expression tag	UNP A0A150BLG9
C	-12	GLY	-	expression tag	UNP A0A150BLG9
C	-11	SER	-	expression tag	UNP A0A150BLG9
C	-10	SER	-	expression tag	UNP A0A150BLG9
C	-9	HIS	-	expression tag	UNP A0A150BLG9
C	-8	HIS	-	expression tag	UNP A0A150BLG9
C	-7	HIS	-	expression tag	UNP A0A150BLG9
C	-6	HIS	-	expression tag	UNP A0A150BLG9
C	-5	HIS	-	expression tag	UNP A0A150BLG9
C	-4	HIS	-	expression tag	UNP A0A150BLG9
C	-3	SER	-	expression tag	UNP A0A150BLG9
C	-2	GLN	-	expression tag	UNP A0A150BLG9
C	-1	ASP	-	expression tag	UNP A0A150BLG9
C	0	PRO	-	expression tag	UNP A0A150BLG9
D	-13	MET	-	expression tag	UNP A0A150BLG9
D	-12	GLY	-	expression tag	UNP A0A150BLG9
D	-11	SER	-	expression tag	UNP A0A150BLG9
D	-10	SER	-	expression tag	UNP A0A150BLG9
D	-9	HIS	-	expression tag	UNP A0A150BLG9
D	-8	HIS	-	expression tag	UNP A0A150BLG9
D	-7	HIS	-	expression tag	UNP A0A150BLG9
D	-6	HIS	-	expression tag	UNP A0A150BLG9
D	-5	HIS	-	expression tag	UNP A0A150BLG9
D	-4	HIS	-	expression tag	UNP A0A150BLG9
D	-3	SER	-	expression tag	UNP A0A150BLG9
D	-2	GLN	-	expression tag	UNP A0A150BLG9
D	-1	ASP	-	expression tag	UNP A0A150BLG9
D	0	PRO	-	expression tag	UNP A0A150BLG9

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Na 1 1	0	0
2	B	1	Total Na 1 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	C	1	Total Na 1 1	0	0
2	D	1	Total Na 1 1	0	0

- Molecule 3 is water.

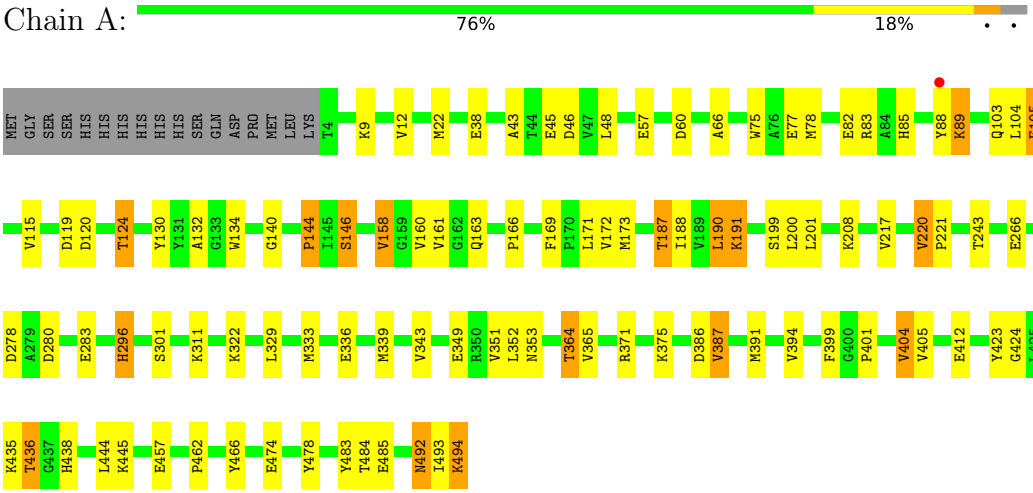
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	26	Total O 26 26	0	0
3	B	23	Total O 23 23	0	0
3	C	26	Total O 26 26	0	0
3	D	17	Total O 17 17	0	0

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.

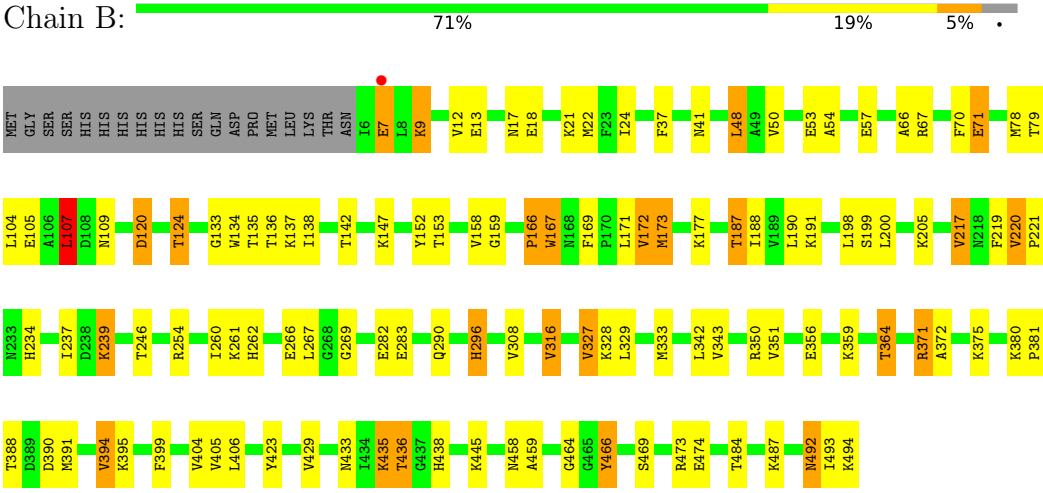
• Molecule 1: Betaine-aldehyde dehydrogenase

Chain A:



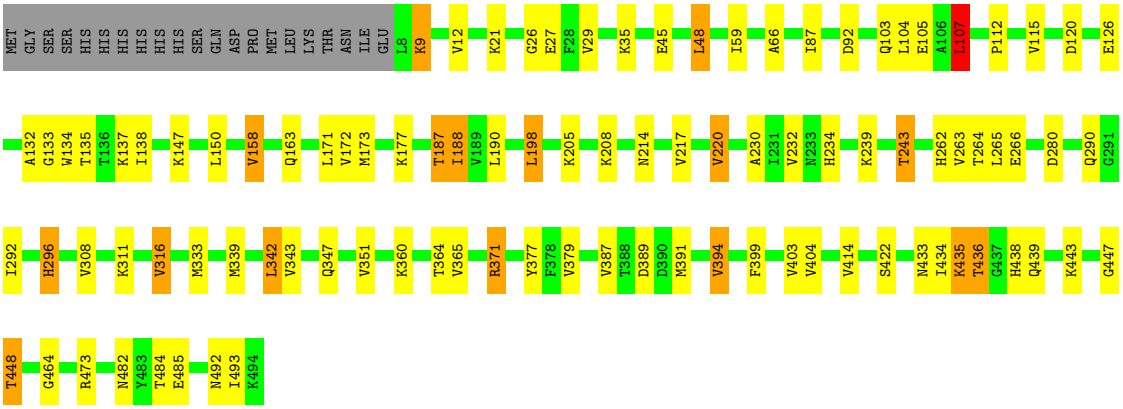
• Molecule 1: Betaine-aldehyde dehydrogenase

Chain B:



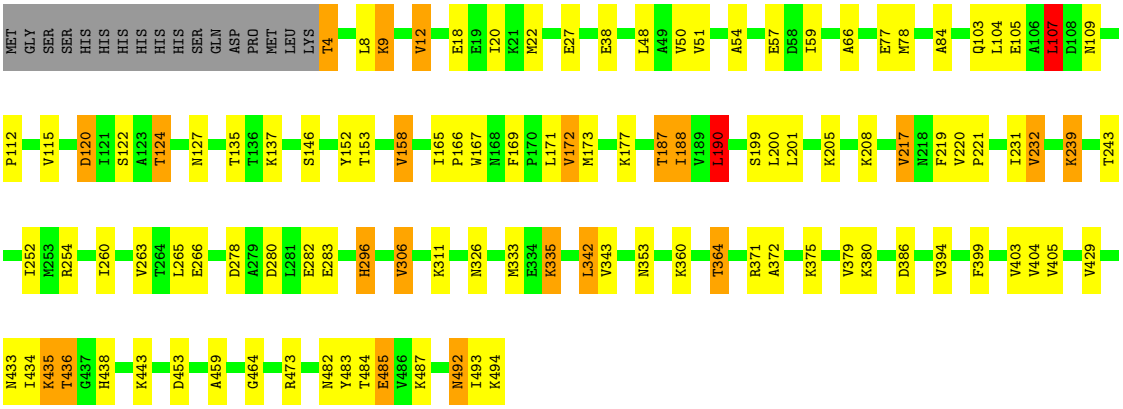
• Molecule 1: Betaine-aldehyde dehydrogenase

Chain C:



• Molecule 1: Betaine-aldehyde dehydrogenase

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	83.53Å 93.28Å 145.47Å 90.00° 98.05° 90.00°	Depositor
Resolution (Å)	49.57 – 2.60 49.57 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.1 (49.57-2.60) 99.1 (49.57-2.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	38.30 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.182 , 0.251 0.194 , 0.255	Depositor DCC
R_{free} test set	3423 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	22.8	Xtriage
Anisotropy	0.127	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 28.3	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	0.92	EDS
Total number of atoms	15213	wwPDB-VP
Average B, all atoms (Å^2)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.56% 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

5.1 Standard geometry

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

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	1.28	5/3870 (0.1%)	1.02	8/5250 (0.2%)
1	B	1.20	2/3855 (0.1%)	1.00	6/5229 (0.1%)
1	C	1.19	2/3838 (0.1%)	1.00	7/5206 (0.1%)
1	D	1.18	4/3870 (0.1%)	0.96	7/5250 (0.1%)
All	All	1.22	13/15433 (0.1%)	0.99	28/20935 (0.1%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	485	GLU	CD-OE2	-6.54	1.18	1.25
1	C	45	GLU	CD-OE2	-6.12	1.19	1.25
1	A	45	GLU	CD-OE1	-5.94	1.19	1.25
1	D	167	TRP	CE3-CZ3	-5.71	1.28	1.38
1	A	483	TYR	CE1-CZ	-5.66	1.31	1.38
1	D	485	GLU	CD-OE2	-5.36	1.19	1.25
1	B	167	TRP	CE3-CZ3	-5.31	1.29	1.38
1	A	474	GLU	CD-OE1	-5.22	1.20	1.25
1	A	485	GLU	CD-OE1	-5.22	1.20	1.25
1	B	466	TYR	CE1-CZ	-5.21	1.31	1.38
1	D	483	TYR	CE1-CZ	-5.19	1.31	1.38
1	C	126	GLU	CG-CD	5.04	1.59	1.51
1	D	38	GLU	CD-OE2	-5.03	1.20	1.25

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	278	ASP	CB-CG-OD1	9.12	126.51	118.30
1	B	350	ARG	NE-CZ-NH1	6.89	123.74	120.30
1	B	107	LEU	CA-CB-CG	6.79	130.92	115.30
1	B	350	ARG	NE-CZ-NH2	-6.74	116.93	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	280	ASP	N-CA-C	-6.68	92.96	111.00
1	C	107	LEU	CA-CB-CG	6.38	129.98	115.30
1	A	278	ASP	CB-CG-OD1	6.31	123.98	118.30
1	C	371	ARG	NE-CZ-NH1	6.23	123.42	120.30
1	D	190	LEU	CA-CB-CG	6.12	129.39	115.30
1	A	190	LEU	CA-CB-CG	6.08	129.28	115.30
1	A	46	ASP	CB-CG-OD1	6.02	123.72	118.30
1	A	352	LEU	CB-CG-CD1	6.00	121.21	111.00
1	B	371	ARG	NE-CZ-NH1	5.67	123.13	120.30
1	D	254	ARG	NE-CZ-NH1	5.61	123.10	120.30
1	D	158	VAL	CB-CA-C	-5.58	100.79	111.40
1	A	107	LEU	CA-CB-CG	5.58	128.12	115.30
1	C	280	ASP	CB-CG-OD1	5.51	123.26	118.30
1	C	92	ASP	CB-CG-OD2	-5.42	113.42	118.30
1	A	158	VAL	CB-CA-C	-5.27	101.39	111.40
1	B	254	ARG	NE-CZ-NH1	5.27	122.93	120.30
1	D	280	ASP	N-CA-C	-5.24	96.85	111.00
1	C	198	LEU	N-CA-C	5.11	124.79	111.00
1	D	107	LEU	CA-CB-CG	5.11	127.05	115.30
1	D	165	ILE	C-N-CD	5.09	139.09	128.40
1	B	41	ASN	N-CA-C	-5.06	97.34	111.00
1	C	150	LEU	N-CA-C	-5.03	97.43	111.00
1	C	158	VAL	CB-CA-C	-5.01	101.88	111.40
1	A	188	ILE	N-CA-C	5.00	124.51	111.00

There are no chirality outliers.

There are no planarity outliers.

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	3791	0	3740	70	1
1	B	3776	0	3727	84	0
1	C	3759	0	3710	68	1
1	D	3791	0	3739	66	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	26	0	0	1	0
3	B	23	0	0	0	0
3	C	26	0	0	1	0
3	D	17	0	0	0	0
All	All	15213	0	14916	261	1

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 9.

All (261) 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:54:ALA:CB	1:B:220:VAL:HG12	1.73	1.18
1:A:494:LYS:HD3	1:C:439:GLN:HE22	1.10	1.11
1:B:54:ALA:HB2	1:B:220:VAL:HG12	1.42	1.00
1:B:7:GLU:N	1:B:7:GLU:OE2	1.94	0.99
1:B:54:ALA:CB	1:B:220:VAL:CG1	2.40	0.98
1:A:494:LYS:HD3	1:C:439:GLN:NE2	1.81	0.95
1:B:237:ILE:O	1:B:261:LYS:NZ	2.02	0.93
1:B:54:ALA:HB2	1:B:220:VAL:CG1	2.01	0.90
1:A:433:ASN:HB3	1:A:436:THR:HG23	1.57	0.87
1:A:492:ASN:HD22	1:A:494:LYS:H	1.22	0.87
1:C:433:ASN:HB3	1:C:436:THR:HG23	1.56	0.86
1:D:266:GLU:OE1	1:D:464:GLY:HA2	1.78	0.83
1:A:492:ASN:ND2	1:A:494:LYS:H	1.77	0.82
1:B:492:ASN:ND2	1:B:494:LYS:H	1.77	0.82
1:A:9:LYS:HB3	1:A:12:VAL:HG13	1.62	0.81
1:A:433:ASN:HD22	1:A:436:THR:H	1.29	0.79
1:A:105:GLU:OE1	1:A:199:SER:OG	2.03	0.77
1:B:66:ALA:HB1	1:B:187:THR:CG2	2.15	0.76
1:B:66:ALA:O	1:B:187:THR:HG21	1.85	0.76
1:D:9:LYS:H	1:D:103:GLN:HE22	1.36	0.74
1:C:188:ILE:HD11	1:C:190:LEU:HB2	1.70	0.73
1:D:66:ALA:HB1	1:D:187:THR:HG23	1.68	0.73
1:D:342:LEU:HD22	1:D:379:VAL:HG13	1.71	0.72
1:D:492:ASN:ND2	1:D:494:LYS:H	1.87	0.72
1:A:66:ALA:O	1:A:187:THR:HG21	1.89	0.71
1:B:364:THR:HG21	1:B:386:ASP:OD2	1.89	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:177:LYS:NZ	1:C:243:THR:HG22	2.06	0.71
1:D:120:ASP:O	1:D:124:THR:HG23	1.89	0.71
1:D:54:ALA:CB	1:D:220:VAL:CG1	2.69	0.70
1:C:66:ALA:O	1:C:187:THR:HG21	1.91	0.70
1:B:364:THR:CG2	1:B:386:ASP:OD2	2.39	0.69
1:C:243:THR:HB	1:C:266:GLU:HB2	1.74	0.69
1:A:492:ASN:HD21	1:A:494:LYS:HA	1.58	0.69
1:B:54:ALA:HB3	1:B:220:VAL:HG12	1.68	0.69
1:B:435:LYS:HG3	1:D:493:ILE:HA	1.75	0.68
1:C:387:VAL:HG11	1:C:404:VAL:HG13	1.75	0.68
1:B:433:ASN:HB3	1:B:436:THR:HG23	1.77	0.67
1:D:124:THR:HG22	1:D:172:VAL:HB	1.76	0.67
1:C:177:LYS:HZ1	1:C:243:THR:HG22	1.60	0.67
1:A:387:VAL:HG13	1:A:394:VAL:HG11	1.77	0.67
1:B:80:THR:CG2	1:B:136:THR:HG22	2.24	0.67
1:B:246:THR:HA	1:B:267:LEU:HD13	1.77	0.67
1:A:462:PRO:HG3	1:A:478:TYR:CD2	2.30	0.66
1:C:448:THR:HG23	3:C:617:HOH:O	1.96	0.66
1:A:9:LYS:H	1:A:103:GLN:HE22	1.45	0.64
1:C:59:ILE:CD1	1:C:220:VAL:HG11	2.27	0.64
1:D:188:ILE:CD1	1:D:190:LEU:HB2	2.27	0.64
1:C:163:GLN:OE1	1:C:188:ILE:HD13	1.98	0.64
1:C:188:ILE:CD1	1:C:190:LEU:HB2	2.27	0.64
1:B:80:THR:HG23	1:B:136:THR:HG22	1.80	0.63
1:B:387:VAL:HG11	1:B:404:VAL:CG1	2.29	0.63
1:A:120:ASP:O	1:A:124:THR:HG23	1.99	0.63
1:B:37:PHE:HD2	1:B:53:GLU:HG3	1.64	0.63
1:B:107:LEU:HD13	1:B:333:MET:HG3	1.79	0.63
1:B:54:ALA:HB1	1:B:220:VAL:CG1	2.27	0.62
1:B:493:ILE:O	1:B:494:LYS:HB2	1.99	0.62
1:D:107:LEU:HD13	1:D:333:MET:HG3	1.80	0.62
1:D:54:ALA:CB	1:D:220:VAL:HG12	2.30	0.62
1:D:66:ALA:O	1:D:187:THR:HG21	2.00	0.62
1:B:433:ASN:HD22	1:B:436:THR:H	1.48	0.61
1:D:243:THR:HG23	1:D:266:GLU:HB3	1.83	0.60
1:A:435:LYS:HG2	1:C:493:ILE:HA	1.84	0.60
1:C:59:ILE:HD13	1:C:220:VAL:HG11	1.83	0.60
1:A:112:PRO:HB2	1:A:115:VAL:HG13	1.84	0.60
1:D:342:LEU:HD22	1:D:379:VAL:CG1	2.31	0.59
1:B:493:ILE:HA	1:D:435:LYS:HG3	1.84	0.59
1:D:177:LYS:NZ	1:D:243:THR:OG1	2.34	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:433:ASN:HD22	1:D:436:THR:H	1.50	0.59
1:D:127:ASN:OD1	1:D:459:ALA:HB2	2.03	0.59
1:C:387:VAL:HG23	1:C:391:MET:SD	2.43	0.58
1:D:66:ALA:HB1	1:D:187:THR:CG2	2.33	0.58
1:B:177:LYS:NZ	1:B:474:GLU:OE1	2.35	0.58
1:C:66:ALA:HB1	1:C:187:THR:CG2	2.34	0.58
1:A:492:ASN:ND2	1:A:494:LYS:N	2.50	0.58
1:C:308:VAL:HG11	1:C:316:VAL:CG1	2.34	0.58
1:C:27:GLU:OE2	1:C:27:GLU:HA	2.04	0.57
1:A:75:TRP:CH2	1:A:83:ARG:HD2	2.39	0.57
1:B:492:ASN:C	1:B:492:ASN:HD22	2.07	0.57
1:C:243:THR:HA	1:C:266:GLU:O	2.04	0.57
1:B:37:PHE:CD2	1:B:53:GLU:HG3	2.40	0.57
1:A:22:MET:HG2	1:A:221:PRO:HD2	1.87	0.57
1:C:264:THR:HG21	1:C:482:ASN:ND2	2.20	0.56
1:C:387:VAL:HG22	1:C:394:VAL:HG11	1.87	0.56
1:B:22:MET:HE3	1:B:24:ILE:HD11	1.88	0.56
1:B:435:LYS:NZ	1:D:492:ASN:ND2	2.54	0.56
1:D:492:ASN:C	1:D:492:ASN:HD22	2.08	0.56
1:A:190:LEU:HD12	1:A:200:LEU:HD21	1.87	0.55
1:B:70:PHE:CE1	1:B:159:GLY:HA2	2.42	0.55
1:D:394:VAL:HB	1:D:404:VAL:HG11	1.89	0.55
1:C:26:GLY:HA3	1:C:214:ASN:ND2	2.22	0.55
1:C:447:GLY:HA3	1:C:464:GLY:O	2.07	0.55
1:A:466:TYR:CE1	1:D:485:GLU:HG3	2.42	0.54
1:A:243:THR:HG23	1:A:266:GLU:HB2	1.90	0.54
1:A:433:ASN:ND2	1:A:436:THR:H	2.03	0.54
1:C:188:ILE:HD11	1:C:190:LEU:CB	2.36	0.54
1:A:66:ALA:HB1	1:A:187:THR:CG2	2.38	0.54
1:B:239:LYS:NZ	1:B:262:HIS:HD2	2.06	0.54
1:B:80:THR:HG23	1:B:136:THR:CG2	2.37	0.53
1:B:67:ARG:O	1:B:71:GLU:HG3	2.07	0.53
1:C:264:THR:HG21	1:C:482:ASN:CG	2.28	0.53
1:A:107:LEU:HD13	1:A:333:MET:HG3	1.88	0.53
1:B:22:MET:HG2	1:B:221:PRO:HD2	1.90	0.53
1:D:217:VAL:HG13	1:D:219:PHE:CZ	2.44	0.53
1:B:266:GLU:OE2	1:B:474:GLU:HG3	2.09	0.53
1:B:66:ALA:HB1	1:B:187:THR:HG23	1.90	0.53
1:D:4:THR:HB	1:D:335:LYS:HD3	1.91	0.52
1:B:387:VAL:HG13	1:B:394:VAL:HG11	1.90	0.52
1:C:387:VAL:HG22	1:C:394:VAL:CG1	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:LYS:HG2	1:A:220:VAL:O	2.10	0.52
1:C:112:PRO:HB2	1:C:115:VAL:HG13	1.92	0.52
1:C:66:ALA:HB1	1:C:187:THR:HG23	1.92	0.51
1:A:493:ILE:HA	1:C:435:LYS:HG3	1.92	0.51
1:B:387:VAL:HG21	1:B:404:VAL:HG13	1.91	0.51
1:C:389:ASP:HA	1:C:394:VAL:HG21	1.91	0.51
1:B:166:PRO:HD2	1:B:173:MET:HG2	1.93	0.51
1:B:356:GLU:OE1	1:B:359:LYS:NZ	2.44	0.51
1:B:492:ASN:HD22	1:B:494:LYS:H	1.56	0.51
1:D:9:LYS:HB3	1:D:12:VAL:HG13	1.91	0.51
1:A:462:PRO:HG3	1:A:478:TYR:CE2	2.45	0.51
1:D:200:LEU:O	1:D:200:LEU:HG	2.10	0.50
1:D:364:THR:CG2	1:D:386:ASP:OD2	2.60	0.50
1:B:260:ILE:O	1:B:260:ILE:HG22	2.12	0.50
1:D:54:ALA:CB	1:D:220:VAL:HG13	2.42	0.50
1:A:339:MET:CE	1:A:401:PRO:HG3	2.42	0.49
1:A:103:GLN:O	1:A:107:LEU:HB2	2.11	0.49
1:C:308:VAL:HG11	1:C:316:VAL:HG13	1.94	0.49
1:B:152:TYR:OH	1:C:438:HIS:HD2	1.96	0.49
1:C:107:LEU:HD13	1:C:333:MET:HG3	1.95	0.49
1:D:54:ALA:HB1	1:D:220:VAL:CG1	2.41	0.49
1:D:464:GLY:HA3	1:D:473:ARG:HD3	1.94	0.49
1:D:492:ASN:HD22	1:D:494:LYS:H	1.59	0.49
1:B:66:ALA:HB1	1:B:187:THR:HG22	1.95	0.49
1:D:433:ASN:ND2	1:D:436:THR:HG23	2.28	0.49
1:A:339:MET:HE3	1:A:401:PRO:HG3	1.94	0.49
1:B:9:LYS:CB	1:B:12:VAL:HG13	2.43	0.48
1:B:22:MET:CE	1:B:24:ILE:HD11	2.42	0.48
1:C:433:ASN:HD22	1:C:436:THR:H	1.61	0.48
1:D:59:ILE:HD12	1:D:231:ILE:HG13	1.95	0.48
1:B:54:ALA:HB1	1:B:220:VAL:HG11	1.96	0.48
1:D:20:ILE:HB	1:D:51:VAL:HG13	1.96	0.48
1:C:21:LYS:HG2	1:C:29:VAL:O	2.14	0.47
1:B:9:LYS:HB2	1:B:12:VAL:HG13	1.96	0.47
1:B:169:PHE:O	1:B:173:MET:HB2	2.15	0.47
1:C:296:HIS:N	1:C:296:HIS:CD2	2.82	0.47
1:A:349:GLU:O	1:A:353:ASN:HB2	2.14	0.47
1:A:438:HIS:HD2	1:D:152:TYR:OH	1.98	0.47
1:B:48:LEU:HD13	1:B:198:LEU:CD1	2.45	0.47
1:C:220:VAL:O	1:C:220:VAL:HG13	2.14	0.47
1:C:9:LYS:H	1:C:103:GLN:HE22	1.62	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:TYR:CD2	1:A:132:ALA:HB1	2.50	0.47
1:A:296:HIS:N	1:A:296:HIS:CD2	2.83	0.47
1:B:464:GLY:HA3	1:B:473:ARG:HD3	1.97	0.47
1:C:292:ILE:HG21	1:C:403:VAL:HB	1.96	0.47
1:B:120:ASP:O	1:B:124:THR:HG23	2.14	0.46
1:B:124:THR:HG22	1:B:172:VAL:HA	1.97	0.46
1:B:134:TRP:CD1	1:D:137:LYS:HE3	2.51	0.46
1:A:438:HIS:CE1	1:B:438:HIS:HE1	2.33	0.46
1:D:201:LEU:HD11	1:D:221:PRO:HG3	1.97	0.46
1:D:22:MET:CE	1:D:220:VAL:HG13	2.44	0.46
1:A:364:THR:HG23	1:A:386:ASP:HB2	1.96	0.46
1:A:120:ASP:O	1:A:124:THR:CG2	2.64	0.46
1:D:364:THR:HG21	1:D:386:ASP:OD2	2.16	0.46
1:C:308:VAL:HG11	1:C:316:VAL:HG11	1.98	0.46
1:A:444:LEU:O	1:D:487:LYS:NZ	2.44	0.46
1:C:239:LYS:HZ3	1:C:262:HIS:HD2	1.64	0.46
1:C:48:LEU:HD11	1:C:107:LEU:HB3	1.97	0.45
1:B:133:GLY:O	1:B:137:LYS:HD2	2.16	0.45
1:A:394:VAL:HB	1:A:404:VAL:HG11	1.98	0.45
1:D:54:ALA:HB3	1:D:220:VAL:HG12	1.99	0.45
1:D:306:VAL:HG12	1:D:403:VAL:CG2	2.46	0.45
1:A:85:HIS:CE1	1:A:89:LYS:HD2	2.52	0.45
1:A:492:ASN:HD22	1:A:492:ASN:C	2.19	0.45
1:A:492:ASN:ND2	1:A:494:LYS:HA	2.30	0.45
1:D:112:PRO:HB2	1:D:115:VAL:HG13	1.99	0.45
1:A:492:ASN:O	1:C:435:LYS:HD2	2.16	0.45
1:D:239:LYS:HD2	1:D:482:ASN:O	2.16	0.45
1:A:160:VAL:HG22	1:A:187:THR:HG22	1.99	0.45
1:A:190:LEU:HD12	1:A:200:LEU:CD2	2.46	0.45
1:C:387:VAL:CG2	1:C:394:VAL:CG1	2.95	0.45
1:B:190:LEU:HD13	1:B:191:LYS:N	2.33	0.44
1:D:120:ASP:O	1:D:124:THR:CG2	2.64	0.44
1:D:84:ALA:HB2	1:D:135:THR:OG1	2.16	0.44
1:B:138:ILE:HG22	1:B:138:ILE:O	2.16	0.44
1:B:372:ALA:HB2	1:B:380:LYS:HG3	1.98	0.44
1:C:133:GLY:O	1:C:137:LYS:HD2	2.17	0.44
1:C:230:ALA:O	1:C:234:HIS:HB2	2.18	0.44
1:A:269:GLY:HA2	1:A:423:TYR:CG	2.53	0.44
1:B:391:MET:HB2	1:B:394:VAL:HG13	1.99	0.44
1:B:296:HIS:CD2	1:B:296:HIS:N	2.85	0.44
1:A:140:GLY:HA3	1:B:142:THR:OG1	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:433:ASN:HD22	1:C:436:THR:CG2	2.31	0.44
1:D:22:MET:HE3	1:D:220:VAL:HG13	1.99	0.44
1:A:438:HIS:HE1	1:B:438:HIS:HE1	1.64	0.43
1:C:21:LYS:HG2	1:C:29:VAL:C	2.38	0.43
1:D:296:HIS:N	1:D:296:HIS:CD2	2.87	0.43
1:B:435:LYS:HZ3	1:D:492:ASN:ND2	2.17	0.43
1:C:134:TRP:O	1:C:138:ILE:HG13	2.18	0.43
1:C:296:HIS:CD2	1:C:339:MET:HG3	2.54	0.43
1:C:434:ILE:HA	1:D:434:ILE:HB	2.00	0.43
1:A:391:MET:HB2	1:A:394:VAL:HG13	2.01	0.43
1:A:492:ASN:HD21	1:A:494:LYS:CA	2.28	0.43
1:B:492:ASN:ND2	1:B:492:ASN:C	2.72	0.43
1:D:153:THR:HA	1:D:487:LYS:O	2.19	0.43
1:A:466:TYR:CZ	1:D:485:GLU:HG3	2.54	0.43
1:A:43:ALA:HB2	3:A:609:HOH:O	2.18	0.43
1:B:308:VAL:HG21	1:B:316:VAL:CG2	2.48	0.43
1:C:433:ASN:ND2	1:C:436:THR:CG2	2.82	0.43
1:B:66:ALA:C	1:B:187:THR:HG21	2.38	0.43
1:B:269:GLY:HA2	1:B:423:TYR:HB3	2.01	0.43
1:B:466:TYR:CE1	1:C:485:GLU:HG3	2.53	0.43
1:C:438:HIS:HE1	1:D:438:HIS:HE1	1.67	0.43
1:A:269:GLY:HA2	1:A:423:TYR:CD1	2.54	0.42
1:C:342:LEU:HB2	1:C:377:TYR:O	2.19	0.42
1:A:364:THR:HG23	1:A:386:ASP:OD2	2.18	0.42
1:B:217:VAL:HG13	1:B:219:PHE:CZ	2.54	0.42
1:A:130:TYR:O	1:C:137:LYS:NZ	2.38	0.42
1:A:134:TRP:CD1	1:C:137:LYS:HE2	2.54	0.42
1:A:201:LEU:HD11	1:A:221:PRO:HG3	2.02	0.42
1:A:364:THR:CG2	1:A:386:ASP:OD2	2.68	0.42
1:B:327:VAL:CG1	1:B:381:PRO:HG2	2.49	0.42
1:B:153:THR:HA	1:B:487:LYS:O	2.20	0.42
1:C:48:LEU:HD13	1:C:198:LEU:CD1	2.50	0.42
1:C:263:VAL:HG12	1:C:265:LEU:HG	2.02	0.42
1:D:263:VAL:HG12	1:D:265:LEU:HG	2.01	0.42
1:D:372:ALA:HB2	1:D:380:LYS:HG3	2.01	0.42
1:B:134:TRP:NE1	1:D:137:LYS:HE3	2.34	0.41
1:B:239:LYS:HZ3	1:B:262:HIS:HD2	1.66	0.41
1:A:435:LYS:CG	1:C:493:ILE:HA	2.49	0.41
1:D:166:PRO:HD2	1:D:173:MET:SD	2.61	0.41
1:D:232:VAL:HG21	1:D:252:ILE:HG12	2.02	0.41
1:D:492:ASN:ND2	1:D:492:ASN:C	2.72	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:492:ASN:ND2	1:D:494:LYS:N	2.64	0.41
1:A:163:GLN:HB2	1:A:190:LEU:HD23	2.03	0.41
1:A:22:MET:HB3	1:A:22:MET:HE2	1.94	0.41
1:C:342:LEU:HD22	1:C:379:VAL:HG23	2.02	0.41
1:D:54:ALA:HB2	1:D:220:VAL:HG13	2.01	0.41
1:A:166:PRO:HD3	1:A:243:THR:O	2.20	0.41
1:A:107:LEU:HD13	1:A:333:MET:CG	2.51	0.41
1:B:458:ASN:O	1:B:459:ALA:C	2.58	0.41
1:C:87:ILE:HG22	1:C:132:ALA:HB2	2.03	0.41
1:C:464:GLY:HA3	1:C:473:ARG:HD3	2.03	0.41
1:A:144:PRO:HG2	1:A:144:PRO:O	2.21	0.41
1:A:146:SER:HB2	1:B:79:THR:HG22	2.03	0.41
1:B:387:VAL:HG11	1:B:404:VAL:HG13	2.03	0.41
1:D:260:ILE:O	1:D:260:ILE:HG22	2.21	0.41
1:A:78:MET:HG2	1:A:82:GLU:CB	2.52	0.40
1:A:271:SER:HB3	1:A:426:ALA:O	2.22	0.40
1:B:167:TRP:CE3	1:B:343:VAL:HG11	2.56	0.40
1:B:364:THR:HG23	1:B:385:THR:HG22	2.02	0.40
1:C:220:VAL:O	1:C:220:VAL:CG1	2.68	0.40
1:A:169:PHE:O	1:A:173:MET:HB2	2.20	0.40
1:B:492:ASN:ND2	1:D:435:LYS:NZ	2.70	0.40
1:C:48:LEU:HD13	1:C:198:LEU:HD12	2.03	0.40
1:A:115:VAL:O	1:A:119:ASP:HB2	2.21	0.40
1:B:78:MET:HG3	1:B:82:GLU:HB2	2.03	0.40
1:B:167:TRP:CD2	1:B:343:VAL:HG11	2.57	0.40
1:B:230:ALA:O	1:B:234:HIS:HB2	2.21	0.40
1:C:347:GLN:O	1:C:351:VAL:HG13	2.21	0.40
1:D:169:PHE:HB3	1:D:172:VAL:HG13	2.02	0.40
1:B:107:LEU:HD13	1:B:333:MET:CG	2.50	0.40
1:C:433:ASN:ND2	1:C:436:THR:HG22	2.37	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:GLU:OE1	1:C:35:LYS:NZ[1_655]	2.13	0.07

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	489/508 (96%)	473 (97%)	15 (3%)	1 (0%)	47 71
1	B	487/508 (96%)	466 (96%)	20 (4%)	1 (0%)	47 71
1	C	485/508 (96%)	473 (98%)	12 (2%)	0	100 100
1	D	489/508 (96%)	467 (96%)	21 (4%)	1 (0%)	47 71
All	All	1950/2032 (96%)	1879 (96%)	68 (4%)	3 (0%)	47 71

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	78	MET
1	A	424	GLY
1	B	166	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	398/414 (96%)	354 (89%)	44 (11%)	6 11
1	B	396/414 (96%)	335 (85%)	61 (15%)	2 4
1	C	394/414 (95%)	353 (90%)	41 (10%)	7 13
1	D	398/414 (96%)	345 (87%)	53 (13%)	4 7
All	All	1586/1656 (96%)	1387 (88%)	199 (12%)	4 8

All (199) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	LEU
1	A	57	GLU
1	A	60	ASP
1	A	77	GLU
1	A	89	LYS
1	A	104	LEU
1	A	105	GLU
1	A	107	LEU
1	A	124	THR
1	A	144	PRO
1	A	146	SER
1	A	158	VAL
1	A	161	VAL
1	A	171	LEU
1	A	172	VAL
1	A	187	THR
1	A	191	LYS
1	A	208	LYS
1	A	217	VAL
1	A	220	VAL
1	A	283	GLU
1	A	296	HIS
1	A	301	SER
1	A	311	LYS
1	A	322	LYS
1	A	329	LEU
1	A	336	GLU
1	A	343	VAL
1	A	351	VAL
1	A	364	THR
1	A	365	VAL
1	A	371	ARG
1	A	375	LYS
1	A	387	VAL
1	A	399	PHE
1	A	404	VAL
1	A	405	VAL
1	A	412	GLU
1	A	436	THR
1	A	445	LYS
1	A	457	GLU
1	A	484	THR
1	A	492	ASN

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Mol	Chain	Res	Type
1	A	494	LYS
1	B	7	GLU
1	B	9	LYS
1	B	13	GLU
1	B	17	ASN
1	B	18	GLU
1	B	21	LYS
1	B	48	LEU
1	B	50	VAL
1	B	57	GLU
1	B	71	GLU
1	B	80	THR
1	B	104	LEU
1	B	105	GLU
1	B	107	LEU
1	B	109	ASN
1	B	120	ASP
1	B	124	THR
1	B	135	THR
1	B	147	LYS
1	B	158	VAL
1	B	171	LEU
1	B	172	VAL
1	B	173	MET
1	B	187	THR
1	B	188	ILE
1	B	199	SER
1	B	200	LEU
1	B	205	LYS
1	B	217	VAL
1	B	220	VAL
1	B	232	VAL
1	B	239	LYS
1	B	282	GLU
1	B	283	GLU
1	B	290	GLN
1	B	296	HIS
1	B	316	VAL
1	B	327	VAL
1	B	328	LYS
1	B	329	LEU
1	B	342	LEU

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Mol	Chain	Res	Type
1	B	351	VAL
1	B	364	THR
1	B	371	ARG
1	B	375	LYS
1	B	385	THR
1	B	387	VAL
1	B	388	THR
1	B	390	ASP
1	B	394	VAL
1	B	395	LYS
1	B	399	PHE
1	B	405	VAL
1	B	406	LEU
1	B	429	VAL
1	B	435	LYS
1	B	436	THR
1	B	445	LYS
1	B	469	SER
1	B	484	THR
1	B	492	ASN
1	C	9	LYS
1	C	12	VAL
1	C	48	LEU
1	C	104	LEU
1	C	105	GLU
1	C	107	LEU
1	C	120	ASP
1	C	135	THR
1	C	147	LYS
1	C	158	VAL
1	C	171	LEU
1	C	172	VAL
1	C	173	MET
1	C	187	THR
1	C	188	ILE
1	C	205	LYS
1	C	208	LYS
1	C	217	VAL
1	C	220	VAL
1	C	232	VAL
1	C	243	THR
1	C	290	GLN

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Mol	Chain	Res	Type
1	C	296	HIS
1	C	311	LYS
1	C	316	VAL
1	C	342	LEU
1	C	343	VAL
1	C	360	LYS
1	C	364	THR
1	C	365	VAL
1	C	371	ARG
1	C	394	VAL
1	C	399	PHE
1	C	414	VAL
1	C	422	SER
1	C	435	LYS
1	C	436	THR
1	C	443	LYS
1	C	448	THR
1	C	484	THR
1	C	492	ASN
1	D	4	THR
1	D	8	LEU
1	D	9	LYS
1	D	12	VAL
1	D	18	GLU
1	D	27	GLU
1	D	48	LEU
1	D	50	VAL
1	D	57	GLU
1	D	77	GLU
1	D	104	LEU
1	D	105	GLU
1	D	107	LEU
1	D	109	ASN
1	D	120	ASP
1	D	122	SER
1	D	124	THR
1	D	146	SER
1	D	158	VAL
1	D	171	LEU
1	D	172	VAL
1	D	187	THR
1	D	188	ILE

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Mol	Chain	Res	Type
1	D	190	LEU
1	D	199	SER
1	D	205	LYS
1	D	208	LYS
1	D	217	VAL
1	D	232	VAL
1	D	239	LYS
1	D	282	GLU
1	D	283	GLU
1	D	296	HIS
1	D	306	VAL
1	D	311	LYS
1	D	326	ASN
1	D	335	LYS
1	D	342	LEU
1	D	343	VAL
1	D	353	ASN
1	D	360	LYS
1	D	364	THR
1	D	371	ARG
1	D	375	LYS
1	D	399	PHE
1	D	405	VAL
1	D	429	VAL
1	D	435	LYS
1	D	436	THR
1	D	443	LYS
1	D	453	ASP
1	D	484	THR
1	D	492	ASN

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

Mol	Chain	Res	Type
1	A	85	HIS
1	A	103	GLN
1	A	214	ASN
1	A	262	HIS
1	A	325	ASN
1	A	326	ASN
1	A	432	GLN
1	A	433	ASN

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Mol	Chain	Res	Type
1	A	438	HIS
1	A	442	ASN
1	A	482	ASN
1	A	492	ASN
1	B	114	GLN
1	B	214	ASN
1	B	262	HIS
1	B	312	HIS
1	B	325	ASN
1	B	326	ASN
1	B	432	GLN
1	B	433	ASN
1	B	438	HIS
1	B	442	ASN
1	B	492	ASN
1	C	85	HIS
1	C	103	GLN
1	C	214	ASN
1	C	262	HIS
1	C	290	GLN
1	C	326	ASN
1	C	357	GLN
1	C	432	GLN
1	C	433	ASN
1	C	438	HIS
1	C	439	GLN
1	C	442	ASN
1	C	482	ASN
1	C	492	ASN
1	D	103	GLN
1	D	214	ASN
1	D	262	HIS
1	D	312	HIS
1	D	325	ASN
1	D	432	GLN
1	D	433	ASN
1	D	438	HIS
1	D	442	ASN
1	D	492	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

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	491/508 (96%)	-0.52	1 (0%)	95	95	11, 18, 34, 46	0
1	B	489/508 (96%)	-0.57	1 (0%)	95	95	10, 19, 33, 63	0
1	C	487/508 (95%)	-0.56	0	100	100	10, 19, 33, 46	0
1	D	491/508 (96%)	-0.44	0	100	100	11, 21, 36, 57	0
All	All	1958/2032 (96%)	-0.52	2 (0%)	95	95	10, 19, 34, 63	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	7	GLU	3.1
1	A	88	TYR	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 monosaccharides in this entry.

6.4 Ligands

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, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	NA	B	501	1/1	0.73	0.24	19,19,19,19	0
2	NA	A	501	1/1	0.81	0.15	16,16,16,16	0
2	NA	D	501	1/1	0.85	0.18	20,20,20,20	0
2	NA	C	501	1/1	0.93	0.22	18,18,18,18	0

6.5 Other polymers (i)

There are no such residues in this entry.