



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 09:08 AM UTC

PDB ID : 3EUW / pdb_00003euw
Title : Crystal Structure of a Myo-inositol dehydrogenase from *Corynebacterium glutamicum* ATCC 13032
Authors : Kumaran, D.; Mahmood, A.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2008-10-11
Resolution : 2.30 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
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

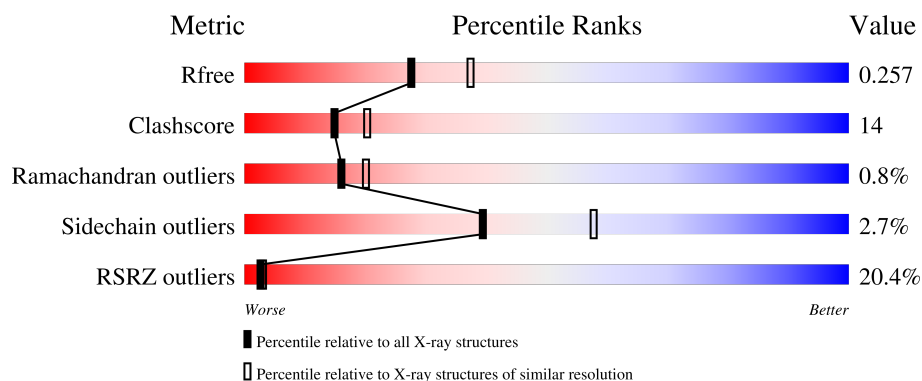
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	344	3% 76% 19% ..
1	B	344	3% 78% 17% ..
1	C	344	35% 64% 31% ..
1	D	344	37% 63% 32% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10763 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 Myo-inositol dehydrogenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	333	Total	C	N	O	S	Se	0	0	0
			2533	1588	445	491	4	5			
1	B	333	Total	C	N	O	S	Se	0	0	0
			2533	1588	445	491	4	5			
1	C	333	Total	C	N	O	S	Se	0	0	0
			2533	1588	445	491	4	5			
1	D	333	Total	C	N	O	S	Se	0	0	0
			2533	1588	445	491	4	5			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	expression tag	UNP Q8NL86
A	0	SER	-	expression tag	UNP Q8NL86
A	1	LEU	-	expression tag	UNP Q8NL86
A	335	GLU	-	expression tag	UNP Q8NL86
A	336	GLY	-	expression tag	UNP Q8NL86
A	337	HIS	-	expression tag	UNP Q8NL86
A	338	HIS	-	expression tag	UNP Q8NL86
A	339	HIS	-	expression tag	UNP Q8NL86
A	340	HIS	-	expression tag	UNP Q8NL86
A	341	HIS	-	expression tag	UNP Q8NL86
A	342	HIS	-	expression tag	UNP Q8NL86
B	-1	MSE	-	expression tag	UNP Q8NL86
B	0	SER	-	expression tag	UNP Q8NL86
B	1	LEU	-	expression tag	UNP Q8NL86
B	335	GLU	-	expression tag	UNP Q8NL86
B	336	GLY	-	expression tag	UNP Q8NL86
B	337	HIS	-	expression tag	UNP Q8NL86
B	338	HIS	-	expression tag	UNP Q8NL86
B	339	HIS	-	expression tag	UNP Q8NL86
B	340	HIS	-	expression tag	UNP Q8NL86
B	341	HIS	-	expression tag	UNP Q8NL86

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Chain	Residue	Modelled	Actual	Comment	Reference
B	342	HIS	-	expression tag	UNP Q8NL86
C	-1	MSE	-	expression tag	UNP Q8NL86
C	0	SER	-	expression tag	UNP Q8NL86
C	1	LEU	-	expression tag	UNP Q8NL86
C	335	GLU	-	expression tag	UNP Q8NL86
C	336	GLY	-	expression tag	UNP Q8NL86
C	337	HIS	-	expression tag	UNP Q8NL86
C	338	HIS	-	expression tag	UNP Q8NL86
C	339	HIS	-	expression tag	UNP Q8NL86
C	340	HIS	-	expression tag	UNP Q8NL86
C	341	HIS	-	expression tag	UNP Q8NL86
C	342	HIS	-	expression tag	UNP Q8NL86
D	-1	MSE	-	expression tag	UNP Q8NL86
D	0	SER	-	expression tag	UNP Q8NL86
D	1	LEU	-	expression tag	UNP Q8NL86
D	335	GLU	-	expression tag	UNP Q8NL86
D	336	GLY	-	expression tag	UNP Q8NL86
D	337	HIS	-	expression tag	UNP Q8NL86
D	338	HIS	-	expression tag	UNP Q8NL86
D	339	HIS	-	expression tag	UNP Q8NL86
D	340	HIS	-	expression tag	UNP Q8NL86
D	341	HIS	-	expression tag	UNP Q8NL86
D	342	HIS	-	expression tag	UNP Q8NL86

- Molecule 2 is SODIUM ION (CCD ID: 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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	231	Total O 231 231	0	0
3	B	201	Total O 201 201	0	0
3	C	96	Total O 96 96	0	0

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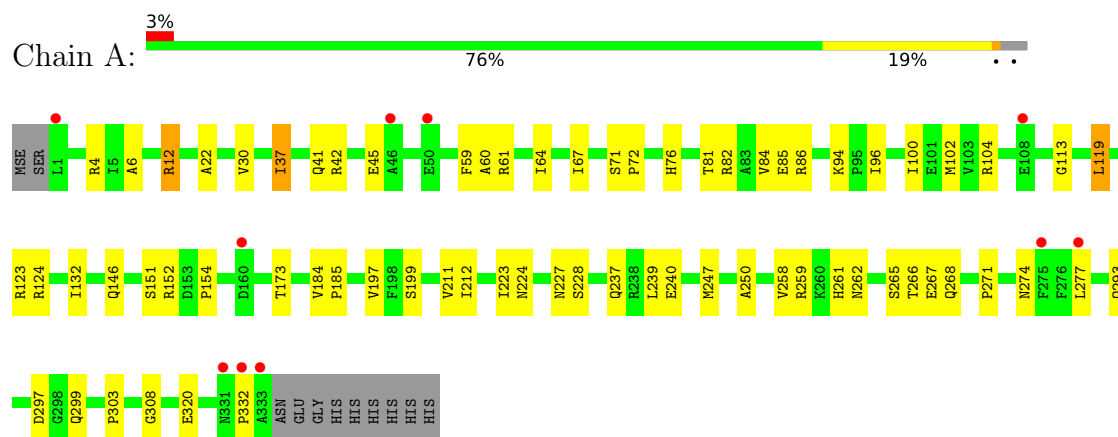
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	101	Total 101	O 101	0	0

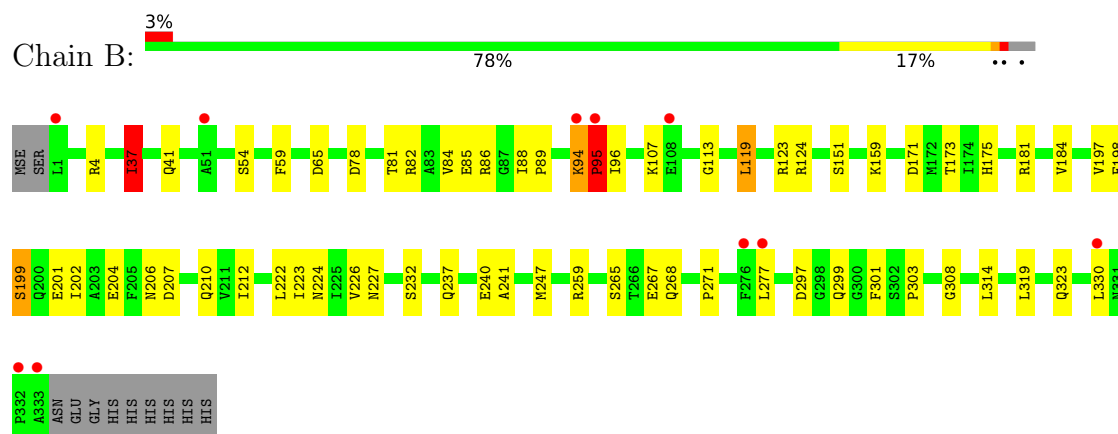
3 Residue-property plots [i](#)

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.

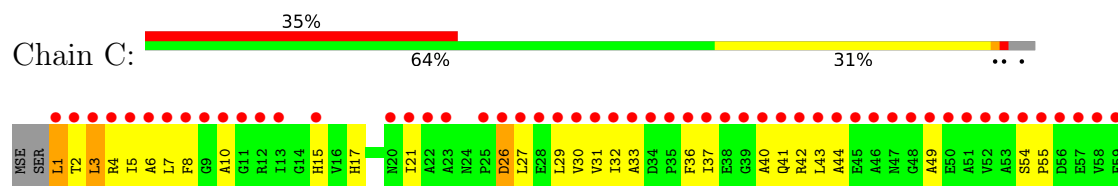
• Molecule 1: Myo-inositol dehydrogenase

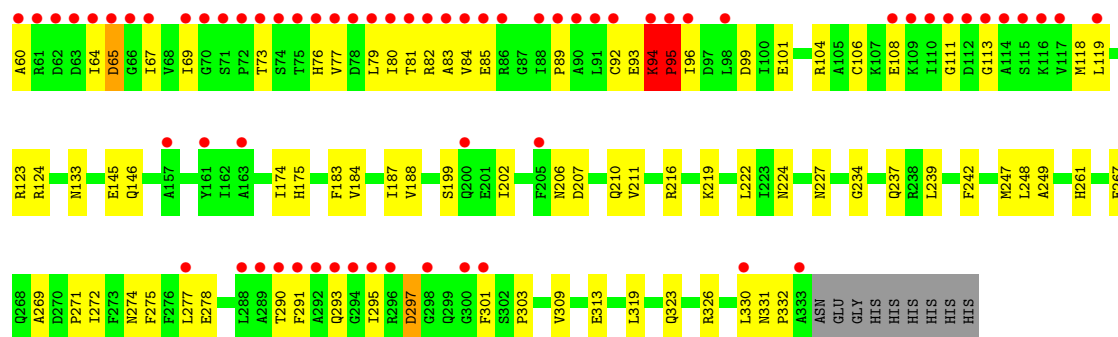


• Molecule 1: Myo-inositol dehydrogenase

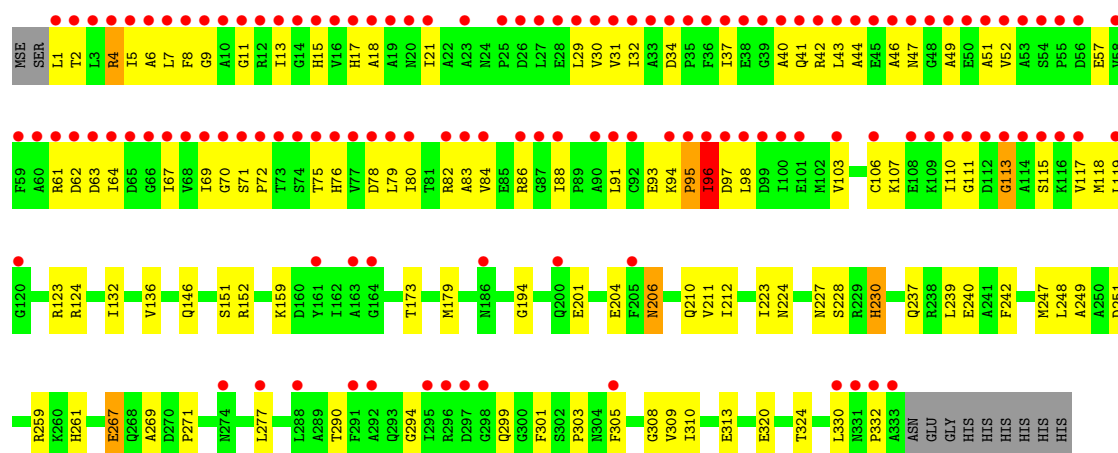


• Molecule 1: Myo-inositol dehydrogenase





• Molecule 1: Myo-inositol dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.04Å 127.59Å 87.16Å 90.00° 92.96° 90.00°	Depositor
Resolution (Å)	38.21 – 2.30 38.21 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.2 (38.21-2.30) 95.1 (38.21-2.30)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.16 (at 2.29Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.223 , 0.258 0.223 , 0.257	Depositor DCC
R_{free} test set	4052 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	21.7	Xtriage
Anisotropy	0.349	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for l,k,-h 0.044 for h,-k,-l 0.017 for l,-k,h	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	10763	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% of the height of the origin peak. No significant pseudotranslation is detected.*

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 ⓘ

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	0.44	0/2572	0.92	10/3481 (0.3%)
1	B	0.45	0/2572	0.98	14/3481 (0.4%)
1	C	0.36	0/2572	0.90	7/3481 (0.2%)
1	D	0.36	0/2572	0.92	5/3481 (0.1%)
All	All	0.40	0/10288	0.93	36/13924 (0.3%)

There are no bond length outliers.

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	96	ILE	N-CA-C	-10.91	99.70	110.72
1	B	94	LYS	CA-C-N	9.85	132.16	119.84
1	B	94	LYS	C-N-CA	9.85	132.16	119.84
1	D	271	PRO	N-CA-C	-7.90	99.56	111.41
1	A	271	PRO	N-CA-C	-7.51	99.44	111.38
1	B	271	PRO	N-CA-C	-7.29	99.78	111.38
1	C	271	PRO	N-CA-C	-7.17	99.97	111.38
1	D	83	ALA	N-CA-C	-6.97	105.32	114.31
1	A	197	VAL	N-CA-C	6.82	120.86	113.43
1	B	95	PRO	CA-N-CD	-6.67	102.66	112.00
1	D	113	GLY	N-CA-C	-6.55	106.75	115.32
1	A	37	ILE	N-CA-C	6.54	117.23	110.36
1	C	94	LYS	CA-C-N	6.41	127.86	119.84
1	C	94	LYS	C-N-CA	6.41	127.86	119.84
1	C	60	ALA	N-CA-C	-6.36	105.47	112.72
1	C	36	PHE	N-CA-C	-6.36	102.64	110.65
1	A	199	SER	N-CA-C	6.19	118.80	108.96
1	B	94	LYS	C-N-CD	-6.08	100.08	125.00
1	A	76	HIS	N-CA-C	5.93	117.55	111.14
1	A	96	ILE	N-CA-C	-5.89	104.77	110.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	42	ARG	N-CA-C	-5.72	106.20	112.72
1	D	117	VAL	N-CA-C	5.72	115.81	106.72
1	B	37	ILE	N-CA-C	5.66	117.07	110.62
1	B	197	VAL	N-CA-C	5.59	119.52	113.43
1	A	60	ALA	N-CA-C	-5.47	106.11	112.89
1	B	184	VAL	CA-C-N	5.46	125.54	119.32
1	B	184	VAL	C-N-CA	5.46	125.54	119.32
1	B	199	SER	N-CA-C	5.42	117.14	108.41
1	B	232	SER	N-CA-C	5.31	117.81	111.71
1	B	88	ILE	N-CA-C	5.30	113.11	107.76
1	C	184	VAL	CA-C-N	5.24	125.30	119.32
1	C	184	VAL	C-N-CA	5.24	125.30	119.32
1	A	184	VAL	CA-C-N	5.14	125.17	119.32
1	A	184	VAL	C-N-CA	5.14	125.17	119.32
1	B	330	LEU	N-CA-C	5.13	116.87	111.28
1	A	258	VAL	N-CA-C	5.10	116.16	109.58

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2533	0	2490	49	0
1	B	2533	0	2490	50	0
1	C	2533	0	2490	92	0
1	D	2533	0	2490	107	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	231	0	0	9	0
3	B	201	0	0	5	0
3	C	96	0	0	5	0
3	D	101	0	0	7	0
All	All	10763	0	9960	277	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 14.

All (277) 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:240:GLU:CD	1:B:247:MSE:HE2	1.92	0.94
3:A:574:HOH:O	1:B:247:MSE:HE3	1.70	0.91
1:C:29:LEU:HD21	1:C:32:ILE:HD11	1.53	0.90
1:B:94:LYS:CB	1:B:95:PRO:HD2	2.04	0.86
1:C:210:GLN:NE2	1:D:146:GLN:HG3	1.91	0.86
1:C:94:LYS:HD2	1:C:95:PRO:HD2	1.57	0.85
1:C:94:LYS:HD3	1:C:175:HIS:NE2	1.99	0.78
1:C:6:ALA:HB2	1:C:64:ILE:HD12	1.67	0.77
1:D:247:MSE:CE	1:D:249:ALA:HB2	2.15	0.77
1:A:224:ASN:HD21	1:B:224:ASN:HD21	1.30	0.77
1:A:12:ARG:HE	1:A:12:ARG:HA	1.51	0.76
1:D:247:MSE:HE1	1:D:249:ALA:HB2	1.69	0.75
1:A:212:ILE:HD13	1:B:222:LEU:HD13	1.67	0.75
1:D:6:ALA:HB2	1:D:64:ILE:HG21	1.69	0.75
1:D:4:ARG:HH11	1:D:4:ARG:HB2	1.53	0.73
1:C:331:ASN:HA	3:C:414:HOH:O	1.90	0.71
1:A:4:ARG:HD2	1:A:30:VAL:HG11	1.74	0.69
1:C:261:HIS:CD2	1:C:267:GLU:H	2.11	0.69
1:A:224:ASN:HD21	1:B:224:ASN:ND2	1.91	0.69
1:A:224:ASN:ND2	1:B:224:ASN:HD21	1.92	0.68
1:C:2:THR:HG22	1:C:3:LEU:H	1.59	0.68
1:D:34:ASP:HB3	1:D:40:ALA:HB2	1.76	0.68
1:D:32:ILE:HG13	3:D:376:HOH:O	1.94	0.67
1:D:107:LYS:HE2	3:D:384:HOH:O	1.93	0.67
1:C:94:LYS:HD2	1:C:95:PRO:CD	2.25	0.67
1:A:81:THR:O	1:A:85:GLU:HG3	1.94	0.66
1:C:123:ARG:CZ	1:C:237:GLN:HE22	2.08	0.66
1:D:247:MSE:HE1	1:D:249:ALA:CB	2.25	0.66
1:D:261:HIS:CD2	1:D:267:GLU:H	2.13	0.66
1:C:332:PRO:HD2	3:C:414:HOH:O	1.95	0.65
1:D:124:ARG:HD2	1:D:303:PRO:HG3	1.78	0.65
1:C:224:ASN:ND2	1:D:224:ASN:HD21	1.94	0.65
1:B:94:LYS:HG3	1:B:95:PRO:HG2	1.79	0.65
1:C:32:ILE:HD13	1:C:44:ALA:HB2	1.79	0.65
1:C:224:ASN:HD21	1:D:224:ASN:HD21	1.45	0.65
1:C:224:ASN:HD21	1:D:224:ASN:ND2	1.95	0.64
1:B:94:LYS:HE3	1:B:171:ASP:O	1.97	0.64
1:A:82:ARG:HD3	3:A:543:HOH:O	1.97	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:210:GLN:HE22	1:D:146:GLN:HG3	1.63	0.63
1:C:31:VAL:C	1:C:32:ILE:HD12	2.23	0.62
1:D:9:GLY:HA3	1:D:70:GLY:O	1.99	0.62
1:C:54:SER:HB2	1:C:55:PRO:HD2	1.80	0.62
1:A:12:ARG:HA	1:A:12:ARG:NE	2.15	0.62
1:B:94:LYS:HB2	1:B:95:PRO:HD2	1.79	0.61
1:C:145:GLU:OE1	1:D:230:HIS:HE1	1.84	0.61
1:D:31:VAL:HG22	1:D:32:ILE:N	2.16	0.61
1:B:94:LYS:HG3	1:B:95:PRO:HD2	1.83	0.60
1:D:61:ARG:HD2	3:D:381:HOH:O	2.00	0.60
1:B:94:LYS:CG	1:B:95:PRO:HD2	2.31	0.60
1:D:67:ILE:HD11	1:D:88:ILE:HD12	1.84	0.60
1:D:30:VAL:C	1:D:49:ALA:HB1	2.27	0.60
1:B:241:ALA:O	1:B:247:MSE:HG2	2.01	0.59
1:D:13:ILE:HG12	1:D:93:GLU:HG3	1.84	0.59
1:C:73:THR:HA	1:C:76:HIS:CE1	2.37	0.59
1:D:118:MSE:HE1	1:D:290:THR:HG22	1.84	0.59
1:C:80:ILE:O	1:C:84:VAL:HG23	2.03	0.59
1:D:4:ARG:O	1:D:5:ILE:HD13	2.03	0.59
1:B:54:SER:HB2	3:B:402:HOH:O	2.03	0.58
1:C:3:LEU:HD12	3:C:404:HOH:O	2.02	0.58
1:C:247:MSE:CE	1:C:249:ALA:HB2	2.33	0.58
1:B:95:PRO:HG2	3:B:347:HOH:O	2.03	0.58
1:C:4:ARG:O	1:C:5:ILE:HD13	2.03	0.58
1:C:84:VAL:HG11	1:C:113:GLY:HA3	1.85	0.58
1:C:124:ARG:HD2	1:C:303:PRO:HG3	1.84	0.58
1:D:123:ARG:HG3	3:D:400:HOH:O	2.04	0.57
1:B:240:GLU:CG	1:B:247:MSE:HE2	2.34	0.57
1:A:42:ARG:HG2	1:A:42:ARG:HH11	1.68	0.57
1:C:26:ASP:O	1:C:27:LEU:HD23	2.05	0.57
1:D:8:PHE:HB2	1:D:69:ILE:HA	1.86	0.56
1:C:37:ILE:O	1:C:41:GLN:HG3	2.05	0.56
1:D:7:LEU:HB2	1:D:29:LEU:HD11	1.86	0.56
1:D:7:LEU:O	1:D:32:ILE:HA	2.06	0.56
1:C:17:HIS:O	1:C:21:ILE:HG13	2.06	0.56
1:D:106:CYS:O	1:D:110:ILE:HG12	2.05	0.56
1:D:247:MSE:HE3	1:D:249:ALA:HB2	1.87	0.55
1:D:63:ASP:O	1:D:64:ILE:HD13	2.07	0.55
1:A:22:ALA:HB2	3:A:481:HOH:O	2.05	0.55
1:A:123:ARG:CZ	1:A:237:GLN:HE22	2.19	0.55
1:A:259:ARG:HA	1:A:268:GLN:O	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:261:HIS:HD2	1:A:266:THR:HB	1.71	0.55
1:C:77:VAL:HG13	1:C:106:CYS:HB2	1.88	0.55
1:B:319:LEU:O	1:B:323:GLN:HG3	2.06	0.54
1:C:37:ILE:HD12	1:C:40:ALA:HB3	1.88	0.54
1:A:261:HIS:CD2	1:A:267:GLU:H	2.25	0.54
1:A:265:SER:HB2	1:C:269:ALA:O	2.07	0.54
1:B:4:ARG:HG3	1:B:4:ARG:HH11	1.73	0.54
1:C:15:HIS:CD2	1:C:43:LEU:HD13	2.42	0.54
1:D:1:LEU:HD23	1:D:2:THR:O	2.08	0.54
1:C:202:ILE:HG23	1:C:207:ASP:HB3	1.90	0.53
1:D:123:ARG:HB3	1:D:179:MSE:CE	2.38	0.53
1:A:262:ASN:HB2	3:D:409:HOH:O	2.07	0.53
1:C:94:LYS:CD	1:C:95:PRO:HD2	2.33	0.53
1:A:119:LEU:HD22	1:A:308:GLY:CA	2.39	0.53
1:C:319:LEU:O	1:C:323:GLN:HG3	2.09	0.53
1:B:202:ILE:HG23	1:B:207:ASP:HB3	1.91	0.52
1:D:123:ARG:NH2	1:D:237:GLN:HE22	2.08	0.52
1:D:94:LYS:CD	1:D:95:PRO:HD2	2.40	0.52
1:B:94:LYS:HG3	1:B:95:PRO:CD	2.39	0.52
1:D:294:GLY:HA3	1:D:301:PHE:CE1	2.45	0.52
1:C:95:PRO:HG3	1:C:174:ILE:HD13	1.92	0.52
1:D:37:ILE:HG13	1:D:41:GLN:OE1	2.09	0.52
1:D:52:VAL:HB	1:D:57:GLU:OE2	2.10	0.52
1:C:76:HIS:O	1:C:80:ILE:HG13	2.10	0.52
1:D:75:THR:O	1:D:79:LEU:HG	2.10	0.52
1:A:4:ARG:HD2	1:A:30:VAL:CG1	2.40	0.52
1:C:30:VAL:C	1:C:49:ALA:HB1	2.35	0.52
1:D:4:ARG:HH11	1:D:4:ARG:CB	2.23	0.52
1:D:94:LYS:HD2	1:D:95:PRO:HD2	1.91	0.52
1:A:61:ARG:HD3	1:A:64:ILE:HD11	1.90	0.51
1:D:44:ALA:CB	1:D:51:ALA:HB2	2.40	0.51
1:C:247:MSE:HE3	1:C:248:LEU:C	2.35	0.51
1:A:151:SER:O	1:A:227:ASN:HA	2.11	0.51
1:B:94:LYS:HG3	1:B:95:PRO:CG	2.40	0.51
1:C:146:GLN:CG	1:D:210:GLN:NE2	2.74	0.51
1:D:6:ALA:CB	1:D:64:ILE:HG21	2.38	0.51
1:A:239:LEU:HD23	1:A:239:LEU:C	2.36	0.51
1:A:293:GLN:HE21	1:A:297:ASP:CG	2.19	0.51
1:B:94:LYS:HD3	1:B:175:HIS:CE1	2.46	0.51
1:C:239:LEU:C	1:C:239:LEU:HD23	2.35	0.51
1:D:31:VAL:HG22	1:D:32:ILE:H	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:113:GLY:C	1:D:115:SER:H	2.19	0.51
1:C:96:ILE:HD12	1:C:96:ILE:H	1.76	0.50
1:D:71:SER:HB2	1:D:72:PRO:HD2	1.93	0.50
1:D:84:VAL:C	1:D:86:ARG:H	2.18	0.50
1:A:100:ILE:O	1:A:104:ARG:HG3	2.11	0.50
1:C:42:ARG:HG2	1:C:42:ARG:HH11	1.76	0.50
1:C:94:LYS:HB3	1:C:95:PRO:HD2	1.93	0.50
1:B:59:PHE:O	1:B:86:ARG:NH1	2.45	0.50
1:B:119:LEU:HD13	1:B:308:GLY:HA3	1.94	0.50
1:C:99:ASP:OD1	1:C:101:GLU:HB2	2.11	0.50
1:C:146:GLN:HB2	1:C:242:PHE:HB3	1.94	0.50
1:A:146:GLN:HE21	1:B:210:GLN:CG	2.24	0.50
1:C:222:LEU:HD12	1:D:210:GLN:HG2	1.93	0.50
1:D:96:ILE:HG21	1:D:103:VAL:HG13	1.94	0.50
1:D:118:MSE:HB2	1:D:301:PHE:CB	2.42	0.49
1:C:3:LEU:HD23	1:C:65:ASP:OD2	2.12	0.49
1:D:118:MSE:HE3	1:D:301:PHE:CD1	2.46	0.49
1:A:102:MSE:HE2	3:A:439:HOH:O	2.11	0.49
1:D:240:GLU:HG3	1:D:247:MSE:SE	2.62	0.49
1:D:247:MSE:CE	1:D:249:ALA:CB	2.86	0.49
1:B:259:ARG:HA	1:B:268:GLN:O	2.12	0.49
1:D:118:MSE:HB2	1:D:301:PHE:HB3	1.95	0.49
1:C:81:THR:C	1:C:83:ALA:H	2.21	0.48
1:D:223:ILE:HD12	1:D:223:ILE:N	2.28	0.48
1:D:18:ALA:HB1	1:D:47:ASN:ND2	2.28	0.48
1:C:10:ALA:HB2	1:C:32:ILE:HG21	1.94	0.48
1:C:291:PHE:O	1:C:295:ILE:HG13	2.14	0.48
1:C:326:ARG:HD2	3:D:356:HOH:O	2.12	0.48
1:D:7:LEU:CB	1:D:29:LEU:HD11	2.43	0.48
3:A:574:HOH:O	1:B:247:MSE:CE	2.45	0.48
1:B:123:ARG:CZ	1:B:237:GLN:HE22	2.27	0.48
1:D:80:ILE:O	1:D:84:VAL:HG23	2.13	0.48
1:C:118:MSE:HE1	1:C:290:THR:HG22	1.95	0.48
1:B:265:SER:HB2	1:D:269:ALA:O	2.14	0.48
1:C:94:LYS:CB	1:C:95:PRO:HD2	2.43	0.48
1:D:78:ASP:O	1:D:82:ARG:HD3	2.13	0.48
1:B:94:LYS:HE2	3:B:347:HOH:O	2.14	0.48
1:D:118:MSE:HE3	1:D:301:PHE:HD1	1.79	0.48
1:B:240:GLU:HG3	1:B:247:MSE:HE2	1.96	0.47
1:B:297:ASP:HB3	1:B:299:GLN:HG2	1.95	0.47
1:C:146:GLN:HE21	1:D:210:GLN:HG3	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:320:GLU:OE2	1:D:324:THR:HG21	2.14	0.47
1:C:309:VAL:O	1:C:313:GLU:HG3	2.14	0.47
1:D:61:ARG:NH2	1:D:63:ASP:OD1	2.46	0.47
1:C:93:GLU:HB3	1:C:94:LYS:H	1.49	0.47
1:C:277:LEU:HD12	1:C:278:GLU:N	2.29	0.47
1:D:17:HIS:O	1:D:21:ILE:HG13	2.14	0.47
1:D:76:HIS:O	1:D:80:ILE:HG13	2.15	0.47
1:D:124:ARG:HD2	1:D:303:PRO:CG	2.43	0.47
1:C:211:VAL:HG22	1:C:227:ASN:HB2	1.97	0.47
1:C:247:MSE:HE3	1:C:249:ALA:HB2	1.96	0.47
1:A:59:PHE:O	1:A:86:ARG:NH1	2.48	0.47
1:C:84:VAL:CG1	1:C:113:GLY:HA3	2.45	0.47
1:C:247:MSE:HE1	1:C:249:ALA:HB2	1.96	0.47
1:C:8:PHE:HD1	1:C:33:ALA:HB3	1.80	0.47
1:D:123:ARG:CZ	1:D:237:GLN:HE22	2.28	0.47
1:C:32:ILE:HD12	1:C:32:ILE:N	2.30	0.46
1:C:133:ASN:HA	1:C:183:PHE:CE1	2.50	0.46
1:A:212:ILE:HD13	1:B:222:LEU:CD1	2.43	0.46
1:C:89:PRO:HG2	1:C:295:ILE:HG23	1.96	0.46
1:D:146:GLN:HB2	1:D:242:PHE:HB3	1.98	0.46
1:D:37:ILE:HA	1:D:40:ALA:HB3	1.97	0.46
1:A:293:GLN:HG3	1:A:297:ASP:OD2	2.16	0.46
1:C:67:ILE:CD1	1:C:83:ALA:HB1	2.45	0.46
1:D:4:ARG:NH1	1:D:63:ASP:O	2.48	0.46
1:D:113:GLY:C	1:D:115:SER:N	2.73	0.46
1:B:301:PHE:HA	3:B:483:HOH:O	2.14	0.46
1:D:15:HIS:ND1	1:D:43:LEU:HD13	2.30	0.46
1:A:240:GLU:HG3	1:A:247:MSE:SE	2.66	0.46
1:C:81:THR:O	1:C:85:GLU:HG2	2.16	0.45
1:D:91:LEU:HD13	1:D:118:MSE:SE	2.66	0.45
1:D:71:SER:HB2	1:D:72:PRO:CD	2.47	0.45
1:B:107:LYS:HE3	3:B:488:HOH:O	2.16	0.45
1:D:41:GLN:HG2	1:D:51:ALA:HB3	1.97	0.45
1:D:152:ARG:HG2	1:D:228:SER:HB3	1.98	0.45
1:A:6:ALA:HB3	1:A:67:ILE:HG12	1.99	0.45
1:A:223:ILE:HD12	1:A:223:ILE:N	2.32	0.45
1:B:159:LYS:HB2	1:B:201:GLU:OE2	2.16	0.45
1:C:219:LYS:HE2	3:C:374:HOH:O	2.16	0.45
1:D:173:THR:HA	1:D:227:ASN:HD21	1.82	0.45
1:C:274:ASN:OD1	1:C:275:PHE:CD2	2.70	0.45
1:D:96:ILE:HB	1:D:97:ASP:H	1.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:PHE:CE1	1:B:199:SER:HB2	2.51	0.45
1:D:37:ILE:HG13	1:D:41:GLN:CD	2.42	0.45
1:D:204:GLU:C	1:D:206:ASN:H	2.24	0.44
3:A:519:HOH:O	1:D:259:ARG:HD2	2.17	0.44
1:D:30:VAL:HG21	1:D:64:ILE:CD1	2.47	0.44
1:D:118:MSE:CE	1:D:290:THR:HG22	2.46	0.44
1:C:187:ILE:O	1:C:330:LEU:HD12	2.18	0.44
1:D:123:ARG:HB3	1:D:179:MSE:HE2	2.00	0.44
1:D:277:LEU:O	1:D:277:LEU:HD12	2.18	0.44
1:C:293:GLN:O	1:C:297:ASP:HB2	2.17	0.44
1:A:124:ARG:HD2	1:A:303:PRO:HG3	1.99	0.44
1:B:201:GLU:O	1:B:204:GLU:HB3	2.17	0.44
1:D:118:MSE:HB2	1:D:301:PHE:CG	2.53	0.44
1:A:12:ARG:HE	1:A:12:ARG:CA	2.18	0.44
1:A:64:ILE:HG12	3:A:391:HOH:O	2.18	0.44
1:A:211:VAL:O	1:A:212:ILE:HD12	2.16	0.44
1:D:151:SER:O	1:D:227:ASN:HA	2.17	0.43
1:C:95:PRO:CG	1:C:174:ILE:HD13	2.47	0.43
1:D:96:ILE:HD12	1:D:305:PHE:CE1	2.53	0.43
1:B:198:PHE:CD1	1:B:198:PHE:C	2.96	0.43
1:C:2:THR:HG22	1:C:3:LEU:N	2.30	0.43
1:D:97:ASP:CG	1:D:98:LEU:H	2.26	0.43
1:D:159:LYS:HB2	1:D:201:GLU:OE2	2.17	0.43
1:A:146:GLN:HE21	1:B:210:GLN:HG3	1.83	0.43
1:C:67:ILE:HD13	1:C:83:ALA:HB1	2.01	0.43
1:C:3:LEU:CD2	1:C:65:ASP:HB2	2.49	0.43
1:A:41:GLN:O	1:A:45:GLU:HG2	2.19	0.43
1:A:211:VAL:C	1:A:212:ILE:HD12	2.44	0.43
1:C:188:VAL:HG22	1:C:216:ARG:O	2.19	0.43
1:D:239:LEU:C	1:D:239:LEU:HD23	2.44	0.43
1:A:152:ARG:HG2	1:A:228:SER:HB3	2.00	0.43
1:D:310:ILE:HA	1:D:313:GLU:OE1	2.19	0.43
1:A:119:LEU:HD22	1:A:308:GLY:HA3	2.00	0.42
1:D:251:ASP:HB3	3:D:367:HOH:O	2.17	0.42
1:C:104:ARG:O	1:C:108:GLU:HG3	2.19	0.42
1:D:31:VAL:CG2	1:D:32:ILE:N	2.82	0.42
1:A:84:VAL:CG1	1:A:113:GLY:HA3	2.49	0.42
1:B:181:ARG:HD2	1:B:314:LEU:HD11	2.02	0.42
1:A:277:LEU:C	1:A:277:LEU:HD12	2.44	0.42
1:C:118:MSE:HB2	1:C:301:PHE:CG	2.54	0.42
1:D:43:LEU:O	1:D:46:ALA:HB3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:194:GLY:HA2	1:D:211:VAL:HG12	2.01	0.42
1:C:7:LEU:HD12	1:C:8:PHE:H	1.84	0.42
1:C:199:SER:OG	1:C:202:ILE:HG13	2.20	0.42
1:D:305:PHE:O	1:D:309:VAL:HG23	2.19	0.42
1:A:132:ILE:HD11	1:A:250:ALA:HB2	2.02	0.42
1:C:69:ILE:HB	1:C:92:CYS:HA	2.00	0.42
1:B:65:ASP:O	1:B:89:PRO:HD2	2.20	0.42
1:D:132:ILE:O	1:D:136:VAL:HG23	2.20	0.42
1:B:37:ILE:O	1:B:41:GLN:HG3	2.19	0.41
1:C:234:GLY:HA3	1:C:272:ILE:HG13	2.02	0.41
1:B:81:THR:O	1:B:85:GLU:HG3	2.19	0.41
1:C:118:MSE:HE3	3:C:391:HOH:O	2.20	0.41
1:C:146:GLN:HG2	1:D:210:GLN:NE2	2.34	0.41
1:A:37:ILE:O	1:A:41:GLN:HG3	2.21	0.41
1:C:31:VAL:HG22	1:C:32:ILE:N	2.35	0.41
1:D:247:MSE:HE3	1:D:248:LEU:C	2.45	0.41
1:A:185:PRO:HD2	3:A:405:HOH:O	2.20	0.41
1:C:224:ASN:HB2	1:D:212:ILE:HD13	2.03	0.41
1:D:119:LEU:CD2	1:D:308:GLY:HA3	2.51	0.41
1:A:71:SER:HB2	1:A:72:PRO:CD	2.50	0.41
1:B:78:ASP:O	1:B:82:ARG:HG3	2.21	0.41
1:B:212:ILE:HD13	1:B:226:VAL:HG22	2.02	0.41
1:C:145:GLU:HB3	1:D:210:GLN:HE22	1.85	0.41
1:B:84:VAL:CG1	1:B:113:GLY:HA3	2.51	0.41
1:B:223:ILE:HD12	1:B:223:ILE:N	2.36	0.41
1:D:52:VAL:HB	1:D:57:GLU:CD	2.46	0.41
1:C:1:LEU:HD13	1:C:2:THR:N	2.36	0.40
1:B:151:SER:O	1:B:227:ASN:HA	2.21	0.40
1:C:8:PHE:CD2	1:C:79:LEU:HD13	2.56	0.40
1:A:299:GLN:HB2	3:A:440:HOH:O	2.21	0.40
1:C:247:MSE:HE1	1:C:249:ALA:CB	2.51	0.40
1:A:154:PRO:HG2	1:A:274:ASN:O	2.22	0.40
1:B:124:ARG:HD2	1:B:303:PRO:HG3	2.03	0.40
1:D:15:HIS:HA	1:D:43:LEU:HD11	2.03	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	331/344 (96%)	321 (97%)	9 (3%)	1 (0%)	36	46
1	B	331/344 (96%)	319 (96%)	11 (3%)	1 (0%)	36	46
1	C	331/344 (96%)	308 (93%)	20 (6%)	3 (1%)	14	17
1	D	331/344 (96%)	294 (89%)	31 (9%)	6 (2%)	6	6
All	All	1324/1376 (96%)	1242 (94%)	71 (5%)	11 (1%)	16	20

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	95	PRO
1	C	95	PRO
1	D	62	ASP
1	D	332	PRO
1	C	82	ARG
1	C	111	GLY
1	D	11	GLY
1	D	96	ILE
1	A	332	PRO
1	D	95	PRO
1	D	111	GLY

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	263/267 (98%)	258 (98%)	5 (2%)	50	69
1	B	263/267 (98%)	256 (97%)	7 (3%)	39	58
1	C	263/267 (98%)	254 (97%)	9 (3%)	32	49
1	D	263/267 (98%)	256 (97%)	7 (3%)	39	58
All	All	1052/1068 (98%)	1024 (97%)	28 (3%)	39	58

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	ARG
1	A	94	LYS
1	A	119	LEU
1	A	173	THR
1	A	320	GLU
1	B	37	ILE
1	B	95	PRO
1	B	119	LEU
1	B	173	THR
1	B	206	ASN
1	B	267	GLU
1	B	277	LEU
1	C	1	LEU
1	C	3	LEU
1	C	26	ASP
1	C	65	ASP
1	C	94	LYS
1	C	95	PRO
1	C	119	LEU
1	C	206	ASN
1	C	297	ASP
1	D	4	ARG
1	D	96	ILE
1	D	206	ASN
1	D	230	HIS
1	D	267	GLU
1	D	299	GLN
1	D	330	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	143	ASN
1	A	210	GLN
1	A	227	ASN
1	A	237	GLN
1	A	261	HIS
1	A	293	GLN
1	A	323	GLN
1	B	143	ASN
1	B	175	HIS
1	B	186	ASN
1	B	196	ASN
1	B	224	ASN
1	B	227	ASN
1	B	237	GLN
1	C	20	ASN
1	C	143	ASN
1	C	186	ASN
1	C	210	GLN
1	C	227	ASN
1	C	237	GLN
1	C	261	HIS
1	C	293	GLN
1	D	17	HIS
1	D	47	ASN
1	D	186	ASN
1	D	210	GLN
1	D	224	ASN
1	D	227	ASN
1	D	230	HIS
1	D	237	GLN
1	D	261	HIS

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](#)

Of 2 ligands modelled in this entry, 2 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 [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	328/344 (95%)	-0.04	10 (3%) 52 54	10, 21, 41, 72	0
1	B	328/344 (95%)	-0.03	10 (3%) 52 54	10, 22, 43, 60	0
1	C	328/344 (95%)	1.61	119 (36%) 1 1	15, 42, 72, 76	0
1	D	328/344 (95%)	1.72	128 (39%) 1 0	14, 41, 76, 80	0
All	All	1312/1376 (95%)	0.81	267 (20%) 3 3	10, 29, 72, 80	0

All (267) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	95	PRO	8.0
1	A	333	ALA	7.9
1	D	43	LEU	7.0
1	D	64	ILE	7.0
1	A	332	PRO	6.9
1	C	60	ALA	6.7
1	D	35	PRO	6.6
1	D	58	VAL	6.5
1	C	64	ILE	6.4
1	D	49	ALA	6.4
1	D	51	ALA	6.2
1	D	8	PHE	6.2
1	C	29	LEU	6.2
1	C	95	PRO	6.2
1	D	40	ALA	6.1
1	D	333	ALA	6.1
1	D	32	ILE	6.0
1	B	95	PRO	6.0
1	C	58	VAL	5.9
1	D	55	PRO	5.9
1	C	30	VAL	5.8

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Mol	Chain	Res	Type	RSRZ
1	D	46	ALA	5.8
1	D	1	LEU	5.7
1	C	52	VAL	5.7
1	D	33	ALA	5.7
1	C	90	ALA	5.6
1	C	1	LEU	5.6
1	C	6	ALA	5.6
1	C	51	ALA	5.6
1	D	29	LEU	5.5
1	D	39	GLY	5.4
1	D	94	LYS	5.3
1	C	333	ALA	5.3
1	D	30	VAL	5.3
1	D	60	ALA	5.3
1	D	31	VAL	5.2
1	D	59	PHE	5.2
1	D	5	ILE	5.2
1	D	7	LEU	5.1
1	C	8	PHE	5.1
1	D	9	GLY	5.1
1	D	113	GLY	5.1
1	C	53	ALA	5.0
1	C	32	ILE	5.0
1	C	67	ILE	5.0
1	C	63	ASP	5.0
1	D	73	THR	5.0
1	D	79	LEU	5.0
1	C	65	ASP	5.0
1	C	114	ALA	5.0
1	C	36	PHE	4.9
1	D	37	ILE	4.9
1	C	3	LEU	4.8
1	C	113	GLY	4.8
1	D	11	GLY	4.8
1	C	39	GLY	4.8
1	D	75	THR	4.8
1	D	10	ALA	4.8
1	D	103	VAL	4.8
1	D	88	ILE	4.7
1	D	53	ALA	4.7
1	D	71	SER	4.6
1	C	55	PRO	4.6

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Mol	Chain	Res	Type	RSRZ
1	D	25	PRO	4.6
1	D	115	SER	4.5
1	C	38	GLU	4.5
1	C	49	ALA	4.5
1	C	56	ASP	4.5
1	D	70	GLY	4.5
1	D	3	LEU	4.4
1	D	298	GLY	4.4
1	C	46	ALA	4.4
1	D	44	ALA	4.4
1	D	332	PRO	4.4
1	C	4	ARG	4.4
1	C	59	PHE	4.4
1	D	52	VAL	4.4
1	D	74	SER	4.4
1	D	330	LEU	4.4
1	C	37	ILE	4.4
1	C	11	GLY	4.3
1	C	83	ALA	4.3
1	C	40	ALA	4.3
1	D	36	PHE	4.2
1	C	50	GLU	4.2
1	D	80	ILE	4.2
1	C	77	VAL	4.2
1	C	31	VAL	4.2
1	C	7	LEU	4.2
1	D	15	HIS	4.1
1	C	5	ILE	4.1
1	C	33	ALA	4.1
1	C	54	SER	4.0
1	D	54	SER	4.0
1	C	27	LEU	4.0
1	D	56	ASP	4.0
1	D	13	ILE	4.0
1	D	34	ASP	3.9
1	D	91	LEU	3.9
1	D	48	GLY	3.9
1	D	65	ASP	3.9
1	C	88	ILE	3.9
1	D	92	CYS	3.9
1	C	96	ILE	3.9
1	C	94	LYS	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	4	ARG	3.9
1	C	79	LEU	3.8
1	C	277	LEU	3.8
1	B	333	ALA	3.8
1	C	44	ALA	3.8
1	D	90	ALA	3.8
1	C	86	ARG	3.8
1	D	67	ILE	3.7
1	D	77	VAL	3.7
1	D	114	ALA	3.7
1	C	25	PRO	3.7
1	C	13	ILE	3.7
1	D	68	VAL	3.7
1	D	110	ILE	3.6
1	D	96	ILE	3.6
1	C	10	ALA	3.6
1	D	38	GLU	3.5
1	D	98	LEU	3.5
1	C	57	GLU	3.5
1	D	69	ILE	3.5
1	C	47	ASN	3.5
1	D	28	GLU	3.5
1	C	35	PRO	3.4
1	C	84	VAL	3.4
1	C	43	LEU	3.4
1	C	9	GLY	3.4
1	C	111	GLY	3.4
1	D	12	ARG	3.4
1	D	18	ALA	3.4
1	C	48	GLY	3.4
1	C	66	GLY	3.4
1	C	291	PHE	3.3
1	D	6	ALA	3.3
1	C	298	GLY	3.3
1	C	69	ILE	3.3
1	D	42	ARG	3.3
1	C	110	ILE	3.3
1	C	12	ARG	3.3
1	C	34	ASP	3.3
1	B	332	PRO	3.3
1	C	71	SER	3.3
1	D	47	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	331	ASN	3.2
1	B	1	LEU	3.2
1	C	295	ILE	3.2
1	C	80	ILE	3.2
1	D	78	ASP	3.2
1	C	116	LYS	3.2
1	C	163	ALA	3.2
1	D	83	ALA	3.2
1	D	292	ALA	3.2
1	C	15	HIS	3.2
1	B	277	LEU	3.2
1	D	72	PRO	3.1
1	C	61	ARG	3.1
1	C	28	GLU	3.1
1	D	27	LEU	3.1
1	D	45	GLU	3.1
1	D	2	THR	3.1
1	D	76	HIS	3.0
1	C	76	HIS	3.0
1	D	17	HIS	3.0
1	D	200	GLN	3.0
1	D	63	ASP	3.0
1	D	97	ASP	3.0
1	C	289	ALA	3.0
1	D	50	GLU	3.0
1	C	92	CYS	3.0
1	D	41	GLN	3.0
1	C	2	THR	2.9
1	C	78	ASP	2.9
1	C	108	GLU	2.9
1	D	14	GLY	2.9
1	D	66	GLY	2.9
1	C	157	ALA	2.9
1	C	26	ASP	2.9
1	C	23	ALA	2.9
1	D	16	VAL	2.9
1	D	117	VAL	2.9
1	C	292	ALA	2.9
1	D	291	PHE	2.9
1	C	73	THR	2.9
1	D	100	ILE	2.9
1	C	300	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	109	LYS	2.8
1	C	91	LEU	2.8
1	B	276	PHE	2.8
1	C	42	ARG	2.8
1	D	108	GLU	2.8
1	D	86	ARG	2.8
1	A	331	ASN	2.8
1	A	277	LEU	2.8
1	C	117	VAL	2.8
1	D	62	ASP	2.7
1	D	119	LEU	2.7
1	C	161	TYR	2.7
1	C	41	GLN	2.7
1	C	75	THR	2.7
1	C	98	LEU	2.7
1	C	21	ILE	2.7
1	D	19	ALA	2.7
1	D	163	ALA	2.7
1	D	161	TYR	2.7
1	D	84	VAL	2.7
1	C	89	PRO	2.7
1	C	62	ASP	2.6
1	C	294	GLY	2.6
1	C	81	THR	2.6
1	C	72	PRO	2.6
1	C	205	PHE	2.6
1	C	200	GLN	2.5
1	D	120	GLY	2.5
1	C	85	GLU	2.5
1	A	1	LEU	2.5
1	D	82	ARG	2.5
1	D	274	ASN	2.5
1	A	50	GLU	2.5
1	C	296	ARG	2.5
1	C	112	ASP	2.5
1	D	106	CYS	2.4
1	A	275	PHE	2.4
1	A	160	ASP	2.4
1	D	61	ARG	2.4
1	D	111	GLY	2.4
1	D	305	PHE	2.4
1	B	108	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	301	PHE	2.4
1	D	295	ILE	2.4
1	C	288	LEU	2.4
1	D	23	ALA	2.4
1	C	109	LYS	2.4
1	C	293	GLN	2.4
1	D	101	GLU	2.4
1	D	288	LEU	2.3
1	D	186	ASN	2.3
1	C	330	LEU	2.3
1	D	277	LEU	2.3
1	B	94	LYS	2.3
1	D	112	ASP	2.3
1	C	22	ALA	2.3
1	C	45	GLU	2.3
1	D	116	LYS	2.3
1	C	119	LEU	2.2
1	C	82	ARG	2.2
1	D	26	ASP	2.2
1	C	74	SER	2.2
1	B	51	ALA	2.2
1	C	290	THR	2.2
1	A	108	GLU	2.1
1	C	115	SER	2.1
1	D	296	ARG	2.1
1	D	87	GLY	2.1
1	A	46	ALA	2.1
1	D	164	GLY	2.1
1	C	20	ASN	2.1
1	D	20	ASN	2.1
1	C	70	GLY	2.1
1	D	205	PHE	2.0
1	B	330	LEU	2.0
1	D	99	ASP	2.0
1	D	297	ASP	2.0
1	D	21	ILE	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 [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NA	B	343	1/1	0.91	0.11	42,42,42,42	0
2	NA	A	343	1/1	0.93	0.06	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.