



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:23 PM UTC

PDB ID : 4EAC / pdb_00004eac
Title : Crystal structure of mannonate dehydratase from Escherichia coli strain K12
Authors : Qiu, X.; Zhu, Y.; Yuan, Y.; Zhang, Y.; Liu, H.; Gao, Y.; Teng, M.; Niu, L.
Deposited on : 2012-03-22
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
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

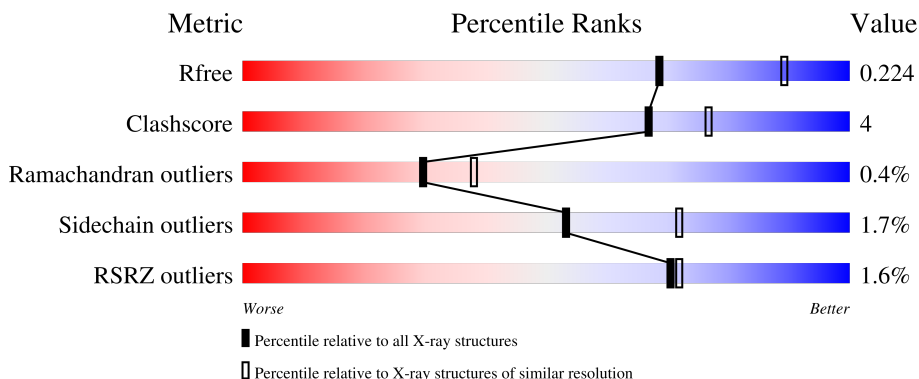
1 Overall quality at a glance

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	414	 % 85% 10% 5%
1	B	414	 86% 8% 5%
1	C	414	 2% 88% 7% 3%
1	D	414	 3% 87% 8% 5%

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 13491 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 Mannonate dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	395	3151	1991	553	590	17	0	0	0
1	B	395	3148	1989	553	589	17	0	0	0
1	C	396	3149	1989	554	589	17	0	0	0
1	D	395	3146	1987	553	589	17	0	0	0

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP P24215
A	-18	GLY	-	expression tag	UNP P24215
A	-17	SER	-	expression tag	UNP P24215
A	-16	SER	-	expression tag	UNP P24215
A	-15	HIS	-	expression tag	UNP P24215
A	-14	HIS	-	expression tag	UNP P24215
A	-13	HIS	-	expression tag	UNP P24215
A	-12	HIS	-	expression tag	UNP P24215
A	-11	HIS	-	expression tag	UNP P24215
A	-10	HIS	-	expression tag	UNP P24215
A	-9	SER	-	expression tag	UNP P24215
A	-8	SER	-	expression tag	UNP P24215
A	-7	GLY	-	expression tag	UNP P24215
A	-6	LEU	-	expression tag	UNP P24215
A	-5	VAL	-	expression tag	UNP P24215
A	-4	PRO	-	expression tag	UNP P24215
A	-3	ARG	-	expression tag	UNP P24215
A	-2	GLY	-	expression tag	UNP P24215
A	-1	SER	-	expression tag	UNP P24215
A	0	HIS	-	expression tag	UNP P24215
B	-19	MET	-	expression tag	UNP P24215

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	expression tag	UNP P24215
B	-17	SER	-	expression tag	UNP P24215
B	-16	SER	-	expression tag	UNP P24215
B	-15	HIS	-	expression tag	UNP P24215
B	-14	HIS	-	expression tag	UNP P24215
B	-13	HIS	-	expression tag	UNP P24215
B	-12	HIS	-	expression tag	UNP P24215
B	-11	HIS	-	expression tag	UNP P24215
B	-10	HIS	-	expression tag	UNP P24215
B	-9	SER	-	expression tag	UNP P24215
B	-8	SER	-	expression tag	UNP P24215
B	-7	GLY	-	expression tag	UNP P24215
B	-6	LEU	-	expression tag	UNP P24215
B	-5	VAL	-	expression tag	UNP P24215
B	-4	PRO	-	expression tag	UNP P24215
B	-3	ARG	-	expression tag	UNP P24215
B	-2	GLY	-	expression tag	UNP P24215
B	-1	SER	-	expression tag	UNP P24215
B	0	HIS	-	expression tag	UNP P24215
C	-19	MET	-	expression tag	UNP P24215
C	-18	GLY	-	expression tag	UNP P24215
C	-17	SER	-	expression tag	UNP P24215
C	-16	SER	-	expression tag	UNP P24215
C	-15	HIS	-	expression tag	UNP P24215
C	-14	HIS	-	expression tag	UNP P24215
C	-13	HIS	-	expression tag	UNP P24215
C	-12	HIS	-	expression tag	UNP P24215
C	-11	HIS	-	expression tag	UNP P24215
C	-10	HIS	-	expression tag	UNP P24215
C	-9	SER	-	expression tag	UNP P24215
C	-8	SER	-	expression tag	UNP P24215
C	-7	GLY	-	expression tag	UNP P24215
C	-6	LEU	-	expression tag	UNP P24215
C	-5	VAL	-	expression tag	UNP P24215
C	-4	PRO	-	expression tag	UNP P24215
C	-3	ARG	-	expression tag	UNP P24215
C	-2	GLY	-	expression tag	UNP P24215
C	-1	SER	-	expression tag	UNP P24215
C	0	HIS	-	expression tag	UNP P24215
D	-19	MET	-	expression tag	UNP P24215
D	-18	GLY	-	expression tag	UNP P24215
D	-17	SER	-	expression tag	UNP P24215

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	expression tag	UNP P24215
D	-15	HIS	-	expression tag	UNP P24215
D	-14	HIS	-	expression tag	UNP P24215
D	-13	HIS	-	expression tag	UNP P24215
D	-12	HIS	-	expression tag	UNP P24215
D	-11	HIS	-	expression tag	UNP P24215
D	-10	HIS	-	expression tag	UNP P24215
D	-9	SER	-	expression tag	UNP P24215
D	-8	SER	-	expression tag	UNP P24215
D	-7	GLY	-	expression tag	UNP P24215
D	-6	LEU	-	expression tag	UNP P24215
D	-5	VAL	-	expression tag	UNP P24215
D	-4	PRO	-	expression tag	UNP P24215
D	-3	ARG	-	expression tag	UNP P24215
D	-2	GLY	-	expression tag	UNP P24215
D	-1	SER	-	expression tag	UNP P24215
D	0	HIS	-	expression tag	UNP P24215

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Cl 1 1	0	0
2	B	1	Total Cl 1 1	0	0
2	C	1	Total Cl 1 1	0	0
2	D	1	Total Cl 1 1	0	0

- Molecule 3 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mn 1 1	0	0
3	B	1	Total Mn 1 1	0	0
3	C	1	Total Mn 1 1	0	0
3	D	1	Total Mn 1 1	0	0

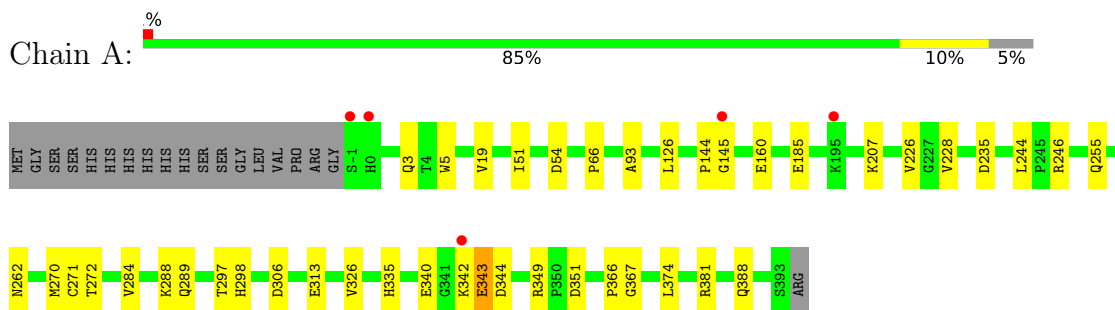
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	273	Total 273	O 273	0	0
4	B	256	Total 256	O 256	0	0
4	C	202	Total 202	O 202	0	0
4	D	158	Total 158	O 158	0	0

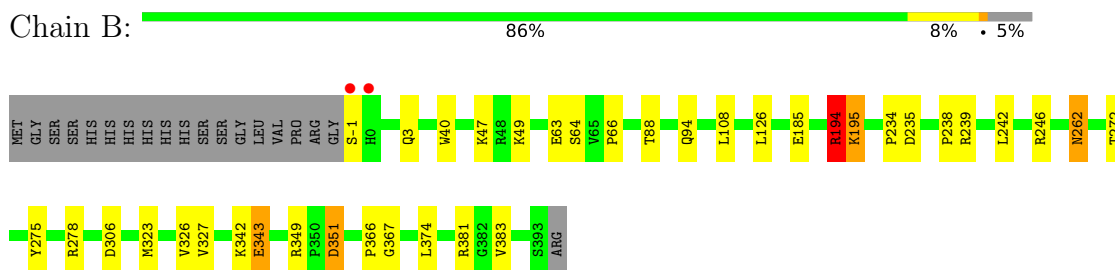
3 Residue-property plots [i](#)

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.

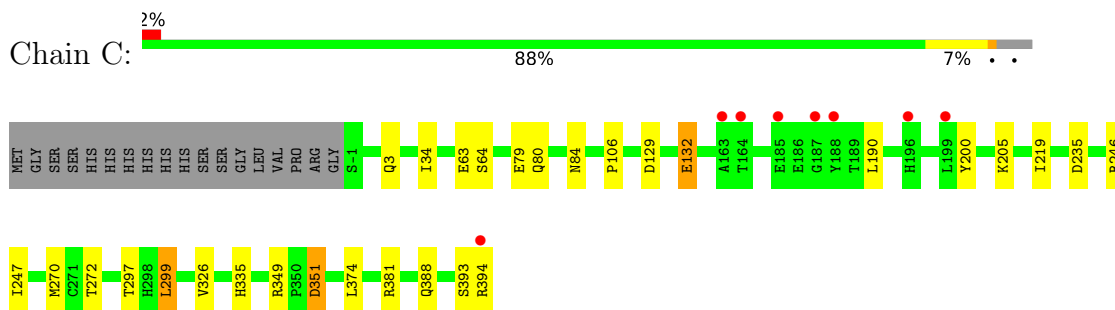
- Molecule 1: Mannonate dehydratase



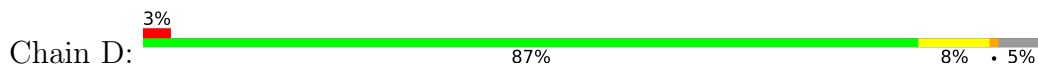
- Molecule 1: Mannonate dehydratase

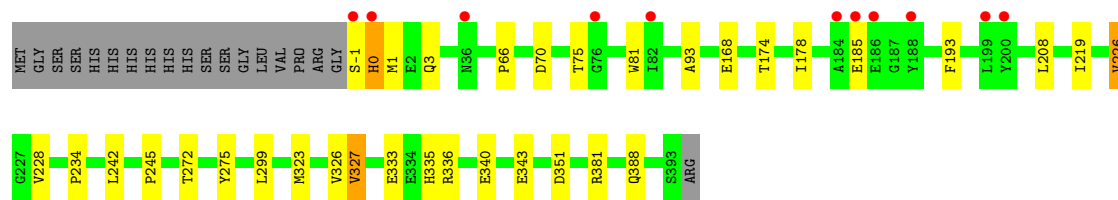


- Molecule 1: Mannonate dehydratase



- Molecule 1: Mannonate dehydratase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	158.96Å 238.46Å 54.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.42 – 2.30 48.42 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.42-2.30) 99.8 (48.42-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.18 (at 2.29Å)	Xtrriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.183 , 0.219 0.188 , 0.224	Depositor DCC
R_{free} test set	4630 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	26.2	Xtrriage
Anisotropy	0.505	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 32.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13491	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.87% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.57	0/3223	0.82	0/4369
1	B	0.54	0/3220	0.87	4/4366 (0.1%)
1	C	0.54	0/3221	0.82	2/4368 (0.0%)
1	D	0.50	0/3218	0.78	0/4364
All	All	0.54	0/12882	0.82	6/17467 (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	B	0	2
All	All	0	3

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	194	ARG	N-CA-C	14.07	126.34	111.14
1	B	194	ARG	CA-C-N	7.16	134.59	121.70
1	B	194	ARG	C-N-CA	7.16	134.59	121.70
1	C	393	SER	N-CA-C	-6.23	105.70	113.55
1	B	262	ASN	N-CA-C	5.49	118.19	111.82
1	C	394	ARG	N-CA-C	-5.33	96.06	111.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	366	PRO	Peptide
1	B	342	LYS	Peptide
1	B	366	PRO	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	3151	0	3078	33	0
1	B	3148	0	3071	27	0
1	C	3149	0	3070	15	0
1	D	3146	0	3067	23	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	273	0	0	4	0
4	B	256	0	0	3	0
4	C	202	0	0	0	0
4	D	158	0	0	0	0
All	All	13491	0	12286	93	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (93) 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:D:323:MET:O	1:D:327:VAL:HG12	1.65	0.97
1:B:323:MET:O	1:B:327:VAL:HG12	1.80	0.80
1:A:93:ALA:CB	1:A:226:VAL:HG12	2.14	0.78
1:A:3:GLN:HE22	1:A:381:ARG:HE	1.30	0.77
1:B:327:VAL:HG11	1:B:383:VAL:CG1	2.17	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:327:VAL:HG11	1:B:383:VAL:HG12	1.71	0.72
1:A:289:GLN:HE22	1:B:94:GLN:HE22	1.38	0.72
1:A:255:GLN:HE21	1:B:94:GLN:HE21	1.37	0.71
1:B:3:GLN:HE22	1:B:381:ARG:HE	1.40	0.69
1:A:342:LYS:O	1:A:343:GLU:HB2	1.92	0.66
1:A:289:GLN:NE2	1:B:94:GLN:HE22	1.93	0.66
1:D:93:ALA:CB	1:D:226:VAL:HG13	2.26	0.65
1:A:367:GLY:HA2	4:A:504:HOH:O	1.96	0.64
1:D:3:GLN:HE22	1:D:381:ARG:HE	1.45	0.64
1:C:3:GLN:HE22	1:C:381:ARG:HE	1.45	0.63
1:A:93:ALA:CB	1:A:226:VAL:CG1	2.77	0.63
1:A:93:ALA:HB1	1:A:226:VAL:HG12	1.82	0.62
1:D:93:ALA:CB	1:D:226:VAL:CG1	2.78	0.62
1:C:335:HIS:HE1	1:C:388:GLN:OE1	1.84	0.61
1:B:126:LEU:HD11	1:B:367:GLY:H	1.66	0.60
1:D:-1:SER:O	1:D:0:HIS:HB2	2.03	0.59
1:A:93:ALA:HB3	1:A:226:VAL:CG1	2.34	0.58
1:A:126:LEU:HD11	1:A:367:GLY:H	1.68	0.58
1:A:66:PRO:HB3	1:A:185:GLU:HG3	1.84	0.57
1:D:335:HIS:HE1	1:D:388:GLN:OE1	1.89	0.56
1:B:272:THR:HG21	1:B:326:VAL:HG11	1.88	0.54
1:B:238:PRO:HG2	1:B:278:ARG:HG3	1.90	0.54
1:D:333:GLU:HA	1:D:336:ARG:HG2	1.90	0.53
1:B:327:VAL:HG11	1:B:383:VAL:HG11	1.88	0.53
1:A:235:ASP:HB3	1:A:246:ARG:HD3	1.91	0.53
1:D:234:PRO:HB3	1:D:275:TYR:CZ	2.45	0.52
4:A:534:HOH:O	1:B:49:LYS:HE3	2.10	0.52
1:A:93:ALA:HB3	1:A:226:VAL:HG11	1.92	0.52
1:A:3:GLN:NE2	1:A:381:ARG:HE	2.05	0.52
1:B:3:GLN:NE2	1:B:381:ARG:HE	2.08	0.52
1:C:3:GLN:NE2	1:C:381:ARG:HH21	2.09	0.51
1:C:349:ARG:NH1	1:C:351:ASP:HA	2.26	0.51
1:B:194:ARG:N	1:B:195:LYS:HB3	2.25	0.51
1:D:174:THR:O	1:D:178:ILE:HG12	2.11	0.50
1:B:66:PRO:HB3	1:B:185:GLU:HG3	1.94	0.49
1:A:226:VAL:HG12	1:A:226:VAL:O	2.12	0.49
1:B:367:GLY:HA2	4:B:745:HOH:O	2.12	0.48
1:B:126:LEU:HD11	1:B:367:GLY:N	2.27	0.48
1:B:235:ASP:HB3	1:B:246:ARG:HD3	1.95	0.48
1:D:272:THR:HG21	1:D:326:VAL:HG11	1.95	0.47
1:C:106:PRO:HB3	1:C:200:TYR:CE1	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:208:LEU:HD22	1:D:245:PRO:HG2	1.97	0.47
1:A:51:ILE:HG13	4:A:766:HOH:O	2.15	0.46
1:B:349:ARG:NH1	1:B:351:ASP:HA	2.31	0.46
1:C:235:ASP:HB3	1:C:246:ARG:HD3	1.97	0.46
1:C:272:THR:HG21	1:C:326:VAL:HG11	1.96	0.46
1:A:270:MET:HB3	1:A:297:THR:HG22	1.98	0.46
1:A:144:PRO:HA	1:A:145:GLY:HA2	1.78	0.46
1:C:80:GLN:HE21	1:C:84:ASN:HD21	1.63	0.46
1:B:40:TRP:HB2	1:B:88:THR:HG23	1.98	0.45
1:C:190:LEU:HD23	1:D:340:GLU:CD	2.42	0.45
1:D:93:ALA:HB3	1:D:226:VAL:HG13	1.98	0.45
1:A:289:GLN:HE22	1:B:94:GLN:NE2	2.12	0.45
1:D:226:VAL:HG12	1:D:228:VAL:HG23	1.99	0.44
1:B:63:GLU:HA	1:B:64:SER:HA	1.76	0.44
1:A:272:THR:HG21	1:A:326:VAL:HG11	1.98	0.44
1:B:306:ASP:HB2	4:B:579:HOH:O	2.16	0.44
1:C:270:MET:HB3	1:C:297:THR:HG22	1.99	0.44
1:A:349:ARG:NH1	1:A:351:ASP:HA	2.32	0.44
1:D:70:ASP:HB2	1:D:81:TRP:CD1	2.54	0.43
1:C:205:LYS:HG3	1:C:247:ILE:HG22	2.01	0.43
1:C:374:LEU:C	1:C:374:LEU:HD23	2.43	0.43
1:D:3:GLN:NE2	1:D:381:ARG:HH21	2.17	0.43
1:C:297:THR:HB	1:C:299:LEU:HD13	2.01	0.43
1:A:5:TRP:CD1	1:A:19:VAL:HG22	2.54	0.43
1:A:126:LEU:HB2	1:A:313:GLU:OE1	2.18	0.43
1:A:342:LYS:HB3	1:A:343:GLU:H	1.49	0.42
1:A:342:LYS:HD3	1:A:344:ASP:HB2	2.00	0.42
1:B:239:ARG:HD3	4:B:551:HOH:O	2.19	0.42
1:D:336:ARG:O	1:D:340:GLU:HG3	2.20	0.42
1:B:108:LEU:HD12	1:B:242:LEU:HD11	2.02	0.42
1:C:63:GLU:HA	1:C:64:SER:HA	1.81	0.42
1:A:126:LEU:HD11	1:A:367:GLY:N	2.34	0.42
1:A:271:CYS:HA	1:A:298:HIS:HB2	2.02	0.41
1:B:234:PRO:HB3	1:B:275:TYR:CZ	2.56	0.41
1:A:335:HIS:HE1	1:A:388:GLN:OE1	2.03	0.41
1:A:374:LEU:C	1:A:374:LEU:HD23	2.46	0.41
1:A:93:ALA:HB2	1:A:228:VAL:HG23	2.02	0.41
1:A:284:VAL:HG12	1:A:288:LYS:HD2	2.01	0.41
1:D:1:MET:HE3	1:D:3:GLN:HG2	2.03	0.41
1:D:66:PRO:HG3	1:D:185:GLU:HG2	2.03	0.41
1:D:193:PHE:HZ	1:D:242:LEU:HD13	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:GLU:HG2	4:A:678:HOH:O	2.20	0.40
1:D:3:GLN:NE2	1:D:381:ARG:HE	2.15	0.40
1:B:374:LEU:C	1:B:374:LEU:HD23	2.47	0.40
1:C:129:ASP:HB3	1:C:132:GLU:HB2	2.03	0.40
1:D:93:ALA:HB2	1:D:226:VAL:CG1	2.51	0.40
1:D:333:GLU:OE2	1:D:336:ARG:HD3	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	393/414 (95%)	379 (96%)	13 (3%)	1 (0%)	36	46
1	B	393/414 (95%)	378 (96%)	12 (3%)	3 (1%)	16	20
1	C	394/414 (95%)	383 (97%)	10 (2%)	1 (0%)	36	46
1	D	393/414 (95%)	378 (96%)	13 (3%)	2 (0%)	24	31
All	All	1573/1656 (95%)	1518 (96%)	48 (3%)	7 (0%)	30	38

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	343	GLU
1	B	195	LYS
1	B	343	GLU
1	D	0	HIS
1	B	351	ASP
1	D	351	ASP
1	C	351	ASP

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	332/351 (95%)	326 (98%)	6 (2%)	51 70
1	B	331/351 (94%)	326 (98%)	5 (2%)	57 75
1	C	331/351 (94%)	326 (98%)	5 (2%)	57 75
1	D	331/351 (94%)	324 (98%)	7 (2%)	47 66
All	All	1325/1404 (94%)	1302 (98%)	23 (2%)	53 72

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	ASP
1	A	207	LYS
1	A	244	LEU
1	A	262	ASN
1	A	306	ASP
1	A	340	GLU
1	B	-1	SER
1	B	47	LYS
1	B	194	ARG
1	B	262	ASN
1	B	343	GLU
1	C	34	ILE
1	C	79	GLU
1	C	132	GLU
1	C	219	ILE
1	C	299	LEU
1	D	75	THR
1	D	168	GLU
1	D	219	ILE
1	D	226	VAL
1	D	299	LEU
1	D	327	VAL
1	D	343	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	74	HIS
1	A	77	ASN
1	A	87	GLN
1	A	255	GLN
1	A	289	GLN
1	A	335	HIS
1	B	3	GLN
1	B	74	HIS
1	B	87	GLN
1	B	289	GLN
1	B	335	HIS
1	C	3	GLN
1	C	74	HIS
1	C	80	GLN
1	C	87	GLN
1	C	196	HIS
1	C	289	GLN
1	C	335	HIS
1	D	3	GLN
1	D	87	GLN
1	D	157	GLN
1	D	196	HIS
1	D	289	GLN
1	D	335	HIS
1	D	388	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 8 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	395/414 (95%)	-0.35	5 (1%) 75 76	9, 22, 42, 50	0
1	B	395/414 (95%)	-0.32	2 (0%) 87 87	9, 23, 45, 56	0
1	C	396/414 (95%)	-0.10	8 (2%) 65 66	9, 27, 56, 71	0
1	D	395/414 (95%)	0.20	11 (2%) 55 57	14, 37, 65, 82	0
All	All	1581/1656 (95%)	-0.14	26 (1%) 70 72	9, 27, 57, 82	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	394	ARG	5.0
1	D	0	HIS	3.7
1	C	185	GLU	3.6
1	C	187	GLY	3.5
1	D	185	GLU	3.5
1	D	-1	SER	3.3
1	A	342	LYS	3.3
1	B	-1	SER	3.1
1	A	-1	SER	2.9
1	C	188	TYR	2.5
1	A	0	HIS	2.5
1	C	163	ALA	2.5
1	D	76	GLY	2.5
1	D	200	TYR	2.4
1	C	164	THR	2.3
1	D	186	GLU	2.3
1	D	199	LEU	2.3
1	A	145	GLY	2.3
1	C	199	LEU	2.2
1	A	195	LYS	2.1
1	D	36	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	0	HIS	2.1
1	C	196	HIS	2.1
1	D	188	TYR	2.1
1	D	82	ILE	2.0
1	D	184	ALA	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 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(Å ²)	Q<0.9
2	CL	A	401	1/1	0.99	0.02	10,10,10,10	0
3	MN	C	402	1/1	0.99	0.02	16,16,16,16	0
2	CL	C	401	1/1	1.00	0.04	14,14,14,14	0
2	CL	D	401	1/1	1.00	0.07	21,21,21,21	0
3	MN	A	402	1/1	1.00	0.01	11,11,11,11	0
3	MN	B	402	1/1	1.00	0.01	12,12,12,12	0
2	CL	B	401	1/1	1.00	0.04	10,10,10,10	0
3	MN	D	402	1/1	1.00	0.02	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.