



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:00 PM UTC

PDB ID : 3ES7 / pdb_00003es7
Title : Crystal structure of divergent enolase from *Oceanobacillus Iheyensis* complexed with Mg and L-malate.
Authors : Fedorov, A.A.; Fedorov, E.V.; Sauder, J.M.; Burley, S.K.; Gerlt, J.A.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2008-10-04
Resolution : 1.90 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

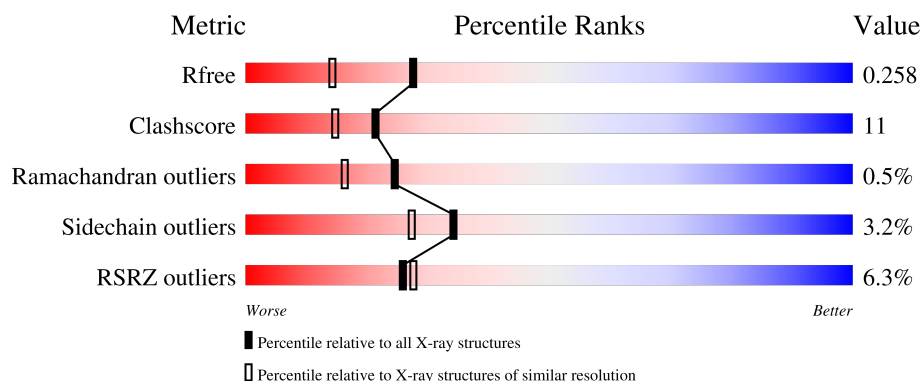
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	391	5% 76% 20% ..
1	B	391	8% 75% 21% ..

2 Entry composition [i](#)

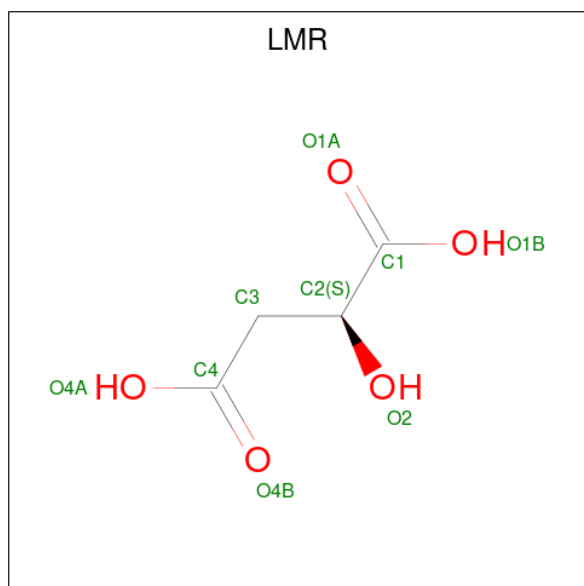
There are 4 unique types of molecules in this entry. The entry contains 6460 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Muconate cycloisomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	387	Total	C	N	O	S	0	0	0
			3097	1976	525	586	10			
1	B	387	Total	C	N	O	S	0	0	0
			3097	1976	525	586	10			

- Molecule 2 is (2S)-2-hydroxybutanedioic acid (CCD ID: LMR) (formula: $C_4H_6O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			9	4	5		
2	B	1	Total	C	O	0	0
			9	4	5		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Mg 1	0	0
3	B	1	Total 1	Mg 1	0	0

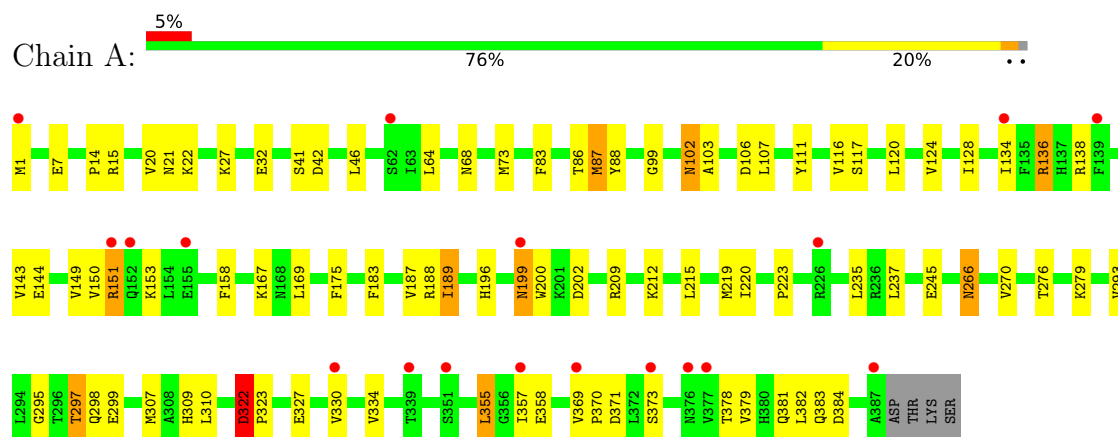
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	131	Total 131	O 131	0	0
4	B	115	Total 115	O 115	0	0

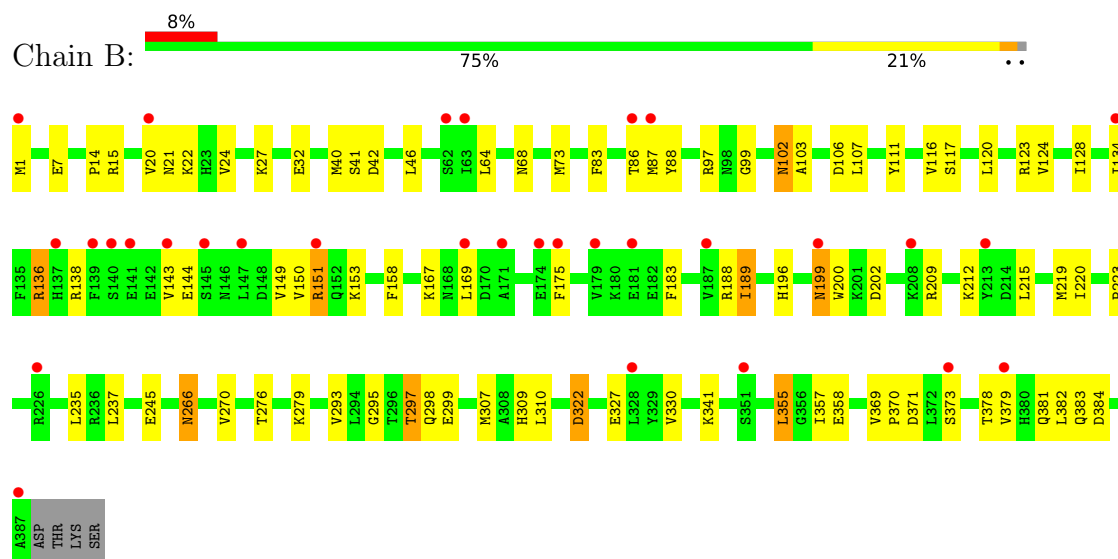
3 Residue-property plots

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: Muconate cycloisomerase



• Molecule 1: Muconate cycloisomerase



4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	121.16Å 121.16Å 122.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.86 – 1.90 24.86 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (24.86-1.90) 99.9 (24.86-1.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.12 (at 1.89Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.227 , 0.259 0.226 , 0.258	Depositor DCC
R_{free} test set	3503 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	31.2	Xtriage
Anisotropy	0.310	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 28.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for l,-k,h 0.008 for -l,-k,-h 0.007 for -h,-l,-k 0.000 for -h,l,k 0.023 for -h,k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6460	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 3.76% 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: LMR, MG

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	0.39	1/3163 (0.0%)	0.93	15/4276 (0.4%)
1	B	0.39	1/3163 (0.0%)	0.93	15/4276 (0.4%)
All	All	0.39	2/6326 (0.0%)	0.93	30/8552 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	307	MET	SD-CE	7.45	1.98	1.79
1	B	307	MET	SD-CE	6.71	1.96	1.79

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	124	VAL	N-CA-C	-7.99	104.97	112.96
1	A	212	LYS	N-CA-C	-7.98	102.81	112.54
1	B	212	LYS	N-CA-C	-7.92	102.87	112.54
1	A	297	THR	N-CA-C	-7.56	102.27	113.61
1	B	297	THR	N-CA-C	-7.46	102.42	113.61
1	A	64	LEU	N-CA-C	7.25	120.61	111.69
1	B	41	SER	N-CA-C	7.18	120.91	110.28
1	A	41	SER	N-CA-C	7.16	120.88	110.28
1	B	124	VAL	N-CA-C	-7.05	105.17	113.42
1	A	138	ARG	N-CA-C	-6.76	105.19	113.50
1	B	138	ARG	N-CA-C	-6.70	105.26	113.50
1	A	200	TRP	N-CA-C	6.64	118.52	111.28
1	A	322	ASP	N-CA-C	6.58	124.34	109.81
1	B	200	TRP	N-CA-C	6.57	118.44	111.28
1	B	199	ASN	N-CA-C	-6.57	101.86	110.53
1	B	322	ASP	N-CA-C	6.55	124.30	109.81
1	B	64	LEU	N-CA-C	6.51	120.44	112.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	199	ASN	N-CA-C	-6.42	102.06	110.53
1	B	136	ARG	N-CA-C	6.28	119.33	110.23
1	A	136	ARG	N-CA-C	6.16	119.16	110.23
1	B	270	VAL	N-CA-C	6.16	116.90	110.62
1	B	143	VAL	N-CA-C	5.85	115.97	110.30
1	A	270	VAL	N-CA-C	5.54	116.94	110.62
1	A	143	VAL	N-CA-C	5.49	115.69	110.42
1	B	83	PHE	CA-C-N	5.49	125.49	119.89
1	B	83	PHE	C-N-CA	5.49	125.49	119.89
1	A	220	ILE	N-CA-C	-5.47	100.45	108.11
1	A	83	PHE	CA-C-N	5.38	125.37	119.89
1	A	83	PHE	C-N-CA	5.38	125.37	119.89
1	B	220	ILE	N-CA-C	-5.10	100.30	107.75

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	3097	0	3076	69	0
1	B	3097	0	3076	69	0
2	A	9	0	4	0	0
2	B	9	0	3	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	131	0	0	1	0
4	B	115	0	0	4	0
All	All	6460	0	6159	135	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 11.

All (135) 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:266:ASN:HD21	1:B:295:GLY:HA3	1.20	1.05
1:A:266:ASN:HD21	1:A:295:GLY:HA3	1.22	1.04
1:B:151:ARG:HB2	1:B:151:ARG:NH1	1.86	0.90
1:A:151:ARG:NH1	1:A:151:ARG:HB2	1.86	0.90
1:A:102:ASN:HD22	1:A:103:ALA:N	1.81	0.78
1:B:102:ASN:HD22	1:B:103:ALA:N	1.81	0.78
1:B:266:ASN:ND2	1:B:295:GLY:HA3	2.01	0.74
1:A:117:SER:OG	1:A:309:HIS:HD2	1.71	0.74
1:B:136:ARG:HD2	1:B:384:ASP:OD1	1.86	0.74
1:B:117:SER:OG	1:B:309:HIS:HD2	1.71	0.73
1:B:151:ARG:HB2	1:B:151:ARG:HH11	1.51	0.73
1:A:151:ARG:HB2	1:A:151:ARG:HH11	1.51	0.72
1:A:136:ARG:HD2	1:A:384:ASP:OD1	1.88	0.72
1:B:134:ILE:HD13	1:B:175:PHE:CZ	2.26	0.70
1:B:134:ILE:HD13	1:B:175:PHE:HZ	1.56	0.69
1:A:266:ASN:HD22	1:A:266:ASN:C	2.01	0.69
1:A:134:ILE:HD13	1:A:175:PHE:CZ	2.28	0.69
1:A:134:ILE:HD13	1:A:175:PHE:HZ	1.57	0.69
1:B:266:ASN:C	1:B:266:ASN:HD22	2.01	0.69
1:A:266:ASN:ND2	1:A:295:GLY:HA3	2.04	0.69
1:A:169:LEU:HD13	1:A:209:ARG:HD3	1.75	0.67
1:B:169:LEU:HD13	1:B:209:ARG:HD3	1.75	0.67
1:B:189:ILE:HD12	1:B:189:ILE:N	2.12	0.64
1:A:189:ILE:N	1:A:189:ILE:HD12	2.12	0.64
1:A:86:THR:O	1:B:86:THR:O	2.17	0.61
1:B:189:ILE:HD13	1:B:215:LEU:HD12	1.81	0.61
1:B:309:HIS:HE1	1:B:355:LEU:O	1.83	0.61
1:A:189:ILE:HD13	1:A:215:LEU:HD12	1.82	0.60
1:A:309:HIS:HE1	1:A:355:LEU:O	1.83	0.60
1:B:22:LYS:HE2	1:B:369:VAL:HG11	1.84	0.60
1:A:22:LYS:HE2	1:A:369:VAL:HG11	1.84	0.59
1:A:187:VAL:HG22	4:A:440:HOH:O	2.05	0.55
1:A:219:MET:HE1	1:A:293:VAL:HG11	1.87	0.55
1:A:73:MET:HE1	1:A:276:THR:HG23	1.88	0.55
1:B:219:MET:HE1	1:B:293:VAL:HG11	1.87	0.55
1:A:102:ASN:HD22	1:A:103:ALA:H	1.52	0.53
1:A:188:ARG:C	1:A:189:ILE:HD12	2.34	0.53
1:B:15:ARG:NE	1:B:21:ASN:HD21	2.06	0.52
1:A:369:VAL:HG13	1:A:370:PRO:HD2	1.91	0.52
1:B:73:MET:HE1	1:B:276:THR:HG23	1.90	0.52
1:B:151:ARG:HB2	1:B:151:ARG:CZ	2.39	0.52
1:A:15:ARG:NE	1:A:21:ASN:HD21	2.07	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:ARG:HB2	1:A:151:ARG:CZ	2.39	0.51
1:B:14:PRO:HB3	1:B:20:VAL:HG22	1.92	0.51
1:B:102:ASN:HD22	1:B:103:ALA:H	1.54	0.51
1:B:369:VAL:HG13	1:B:370:PRO:HD2	1.93	0.51
1:B:188:ARG:C	1:B:189:ILE:HD12	2.36	0.51
1:A:7:GLU:HB2	1:A:27:LYS:HB2	1.91	0.50
1:A:14:PRO:HB3	1:A:20:VAL:HG22	1.93	0.50
1:A:88:TYR:HB2	1:B:86:THR:HB	1.94	0.50
1:A:378:THR:OG1	1:A:381:GLN:HG3	2.11	0.50
1:A:68:ASN:HD22	1:A:68:ASN:C	2.18	0.50
1:B:46:LEU:HD23	4:B:456:HOH:O	2.11	0.50
1:B:378:THR:OG1	1:B:381:GLN:HG3	2.12	0.49
1:B:102:ASN:HD22	1:B:102:ASN:C	2.20	0.49
1:A:102:ASN:HD22	1:A:102:ASN:C	2.20	0.49
1:B:7:GLU:HB2	1:B:27:LYS:HB2	1.92	0.49
1:B:120:LEU:HD22	1:B:279:LYS:HD3	1.94	0.49
1:B:68:ASN:C	1:B:68:ASN:HD22	2.19	0.48
1:B:116:VAL:HG21	1:B:355:LEU:HD13	1.95	0.48
1:A:116:VAL:HG21	1:A:355:LEU:HD13	1.95	0.48
1:A:369:VAL:CG1	1:A:370:PRO:HD2	2.44	0.47
1:B:73:MET:HA	1:B:73:MET:HE2	1.95	0.47
1:A:73:MET:HA	1:A:73:MET:HE2	1.95	0.47
1:A:120:LEU:HD22	1:A:279:LYS:HD3	1.96	0.47
1:A:266:ASN:HD21	1:A:295:GLY:CA	2.11	0.47
1:A:196:HIS:HE1	1:A:245:GLU:OE1	1.99	0.46
1:B:14:PRO:HG2	1:B:330:VAL:HG21	1.98	0.46
1:B:369:VAL:CG1	1:B:370:PRO:HD2	2.46	0.46
1:A:14:PRO:HG2	1:A:330:VAL:HG21	1.98	0.46
1:A:149:VAL:O	1:A:153:LYS:HG2	2.15	0.46
1:B:149:VAL:O	1:B:153:LYS:HG2	2.16	0.46
1:A:297:THR:O	1:A:298:GLN:HB2	2.15	0.46
1:A:134:ILE:HD11	1:A:150:VAL:HG22	1.98	0.46
1:A:199:ASN:ND2	1:A:202:ASP:OD2	2.49	0.46
1:A:86:THR:HB	1:B:88:TYR:HB2	1.99	0.45
1:B:196:HIS:HE1	1:B:245:GLU:OE1	2.00	0.45
1:B:199:ASN:ND2	1:B:202:ASP:OD2	2.49	0.45
1:B:266:ASN:HD21	1:B:295:GLY:CA	2.08	0.45
1:A:266:ASN:HA	1:A:293:VAL:HG23	1.98	0.45
1:B:15:ARG:HE	1:B:21:ASN:ND2	2.14	0.45
1:B:136:ARG:CD	1:B:167:LYS:HD3	2.47	0.45
1:B:151:ARG:HH11	1:B:183:PHE:HZ	1.65	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:ILE:HD11	1:B:150:VAL:HG22	1.98	0.44
1:B:14:PRO:HG2	1:B:330:VAL:CG2	2.48	0.44
1:A:151:ARG:HH11	1:A:183:PHE:HZ	1.66	0.44
1:B:223:PRO:HB2	1:B:235:LEU:CD2	2.47	0.44
1:A:14:PRO:HG2	1:A:330:VAL:CG2	2.48	0.44
1:A:15:ARG:HE	1:A:21:ASN:ND2	2.15	0.44
1:A:266:ASN:C	1:A:266:ASN:ND2	2.72	0.44
1:A:136:ARG:CD	1:A:167:LYS:HD3	2.48	0.43
1:B:297:THR:O	1:B:298:GLN:HB2	2.17	0.43
1:A:189:ILE:N	1:A:189:ILE:CD1	2.80	0.43
1:B:123:ARG:HG2	4:B:467:HOH:O	2.17	0.43
1:A:99:GLY:O	1:A:102:ASN:ND2	2.52	0.43
1:B:189:ILE:N	1:B:189:ILE:CD1	2.80	0.43
1:A:1:MET:HE3	1:A:111:TYR:CD1	2.54	0.43
1:B:99:GLY:O	1:B:102:ASN:ND2	2.52	0.43
1:B:1:MET:HE3	1:B:111:TYR:CD1	2.54	0.43
1:B:371:ASP:OD1	1:B:373:SER:HB2	2.19	0.43
1:A:223:PRO:HB2	1:A:235:LEU:CD2	2.49	0.43
1:B:341:LYS:HD2	4:B:532:HOH:O	2.18	0.43
1:A:334:VAL:HG11	1:A:357:ILE:CG2	2.49	0.43
1:A:371:ASP:OD1	1:A:373:SER:HB2	2.19	0.43
1:B:223:PRO:HB2	1:B:235:LEU:HD21	2.01	0.43
1:B:266:ASN:HA	1:B:293:VAL:HG23	1.99	0.42
1:A:128:ILE:O	1:A:128:ILE:HG13	2.19	0.42
1:B:379:VAL:O	1:B:383:GLN:HG3	2.19	0.42
1:B:106:ASP:HB2	1:B:355:LEU:HD22	2.00	0.42
1:B:24:VAL:HB	1:B:40:MET:HB2	2.01	0.42
1:B:189:ILE:HD13	1:B:215:LEU:CD1	2.49	0.42
1:A:189:ILE:HD13	1:A:215:LEU:CD1	2.49	0.41
1:A:357:ILE:HG22	1:A:358:GLU:N	2.34	0.41
1:A:223:PRO:HB2	1:A:235:LEU:HD21	2.02	0.41
1:B:128:ILE:O	1:B:128:ILE:HG13	2.19	0.41
1:A:334:VAL:HG11	1:A:357:ILE:HG21	2.03	0.41
1:B:357:ILE:HG22	1:B:358:GLU:N	2.34	0.41
1:A:158:PHE:HZ	1:A:327:GLU:HG3	1.86	0.41
1:A:196:HIS:HE1	1:A:245:GLU:CD	2.28	0.41
1:B:46:LEU:CD2	1:B:382:LEU:HB3	2.51	0.41
1:B:266:ASN:ND2	1:B:266:ASN:C	2.73	0.41
1:A:73:MET:HE1	1:A:276:THR:CG2	2.51	0.41
1:A:106:ASP:HB2	1:A:355:LEU:HD22	2.01	0.41
1:A:322:ASP:N	1:A:323:PRO:HD2	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:ARG:HD3	1:B:167:LYS:HD3	2.02	0.41
1:A:136:ARG:HD3	1:A:167:LYS:HD3	2.02	0.41
1:A:1:MET:HE1	1:A:32:GLU:HB2	2.02	0.41
1:B:97:ARG:HD2	4:B:476:HOH:O	2.20	0.40
1:B:158:PHE:HZ	1:B:327:GLU:HG3	1.86	0.40
1:B:297:THR:HA	2:B:392:LMR:O2	2.21	0.40
1:A:379:VAL:O	1:A:383:GLN:HG3	2.21	0.40
1:A:46:LEU:CD2	1:A:382:LEU:HB3	2.52	0.40
1:B:1:MET:HE1	1:B:32:GLU:HB2	2.03	0.40
1:B:196:HIS:HE1	1:B:245:GLU:CD	2.29	0.40
1:A:86:THR:O	1:A:87:MET:CB	2.69	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	385/391 (98%)	374 (97%)	9 (2%)	2 (0%)	24	16
1	B	385/391 (98%)	374 (97%)	9 (2%)	2 (0%)	24	16
All	All	770/782 (98%)	748 (97%)	18 (2%)	4 (0%)	24	16

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	42	ASP
1	B	42	ASP
1	A	322	ASP
1	B	322	ASP

5.3.2 Protein sidechains ⓘ

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	341/345 (99%)	330 (97%)	11 (3%)	34	27
1	B	341/345 (99%)	330 (97%)	11 (3%)	34	27
All	All	682/690 (99%)	660 (97%)	22 (3%)	34	27

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	87	MET
1	A	102	ASN
1	A	107	LEU
1	A	144	GLU
1	A	151	ARG
1	A	189	ILE
1	A	237	LEU
1	A	266	ASN
1	A	299	GLU
1	A	310	LEU
1	A	355	LEU
1	B	87	MET
1	B	102	ASN
1	B	107	LEU
1	B	144	GLU
1	B	151	ARG
1	B	189	ILE
1	B	237	LEU
1	B	266	ASN
1	B	299	GLU
1	B	310	LEU
1	B	355	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	68	ASN
1	A	82	ASN
1	A	102	ASN
1	A	146	ASN
1	A	196	HIS
1	A	254	GLN
1	A	266	ASN
1	A	309	HIS
1	A	368	GLN
1	A	376	ASN
1	A	380	HIS
1	B	21	ASN
1	B	68	ASN
1	B	102	ASN
1	B	146	ASN
1	B	196	HIS
1	B	254	GLN
1	B	266	ASN
1	B	309	HIS
1	B	368	GLN
1	B	376	ASN
1	B	380	HIS

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 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	LMR	A	392	3	8,8,8	1.18	0	10,10,10	2.14	2 (20%)
2	LMR	B	392	3	8,8,8	1.29	0	10,10,10	2.16	2 (20%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	LMR	A	392	3	-	2/8/8/8	-
2	LMR	B	392	3	-	2/8/8/8	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	392	LMR	O1B-C1-C2	4.81	122.91	112.74
2	A	392	LMR	O1B-C1-C2	4.75	122.79	112.74
2	B	392	LMR	O1B-C1-O1A	-2.70	117.94	124.08
2	A	392	LMR	O1B-C1-O1A	-2.70	117.96	124.08

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	392	LMR	O1A-C1-C2-O2
2	A	392	LMR	O1B-C1-C2-O2
2	B	392	LMR	O1A-C1-C2-O2
2	B	392	LMR	O1B-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	392	LMR	1	0

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	387/391 (98%)	0.47	18 (4%) 36 39	22, 32, 44, 54	0
1	B	387/391 (98%)	0.63	31 (8%) 18 19	22, 33, 49, 61	0
All	All	774/782 (98%)	0.55	49 (6%) 26 27	22, 33, 47, 61	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	MET	4.1
1	B	387	ALA	3.9
1	A	387	ALA	3.6
1	B	379	VAL	3.1
1	B	147	LEU	3.0
1	A	1	MET	3.0
1	A	351	SER	3.0
1	B	139	PHE	2.9
1	B	351	SER	2.8
1	B	86	THR	2.7
1	A	357	ILE	2.7
1	A	151	ARG	2.5
1	B	134	ILE	2.5
1	B	171	ALA	2.5
1	B	199	ASN	2.5
1	A	134	ILE	2.5
1	B	140	SER	2.5
1	A	330	VAL	2.4
1	A	339	THR	2.4
1	B	328	LEU	2.3
1	B	63	ILE	2.3
1	A	199	ASN	2.3
1	B	20	VAL	2.3
1	A	155	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	169	LEU	2.3
1	B	87	MET	2.3
1	B	137	HIS	2.3
1	B	175	PHE	2.3
1	B	226	ARG	2.3
1	B	143	VAL	2.3
1	A	62	SER	2.2
1	B	208	LYS	2.2
1	A	373	SER	2.2
1	B	213	TYR	2.2
1	B	179	VAL	2.2
1	B	174	GLU	2.2
1	B	373	SER	2.2
1	A	376	ASN	2.2
1	B	181	GLU	2.2
1	A	369	VAL	2.2
1	A	377	VAL	2.2
1	A	139	PHE	2.1
1	B	145	SER	2.1
1	A	226	ARG	2.1
1	B	62	SER	2.1
1	B	151	ARG	2.1
1	B	141	GLU	2.0
1	A	152	GLN	2.0
1	B	187	VAL	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 oligosaccharides 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, 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	LMR	B	392	9/9	0.80	0.13	34,38,41,42	0
2	LMR	A	392	9/9	0.86	0.11	30,32,38,40	0
3	MG	A	393	1/1	0.99	0.08	19,19,19,19	0
3	MG	B	393	1/1	0.99	0.07	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.