



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:44 PM UTC

PDB ID : 3BLF / pdb_00003blf
Title : Crystal structure of the catalytic domain of LiCMS in complexed with pyruvate
Authors : Zhang, P.; Ma, J.
Deposited on : 2007-12-11
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 : 4-5-2 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

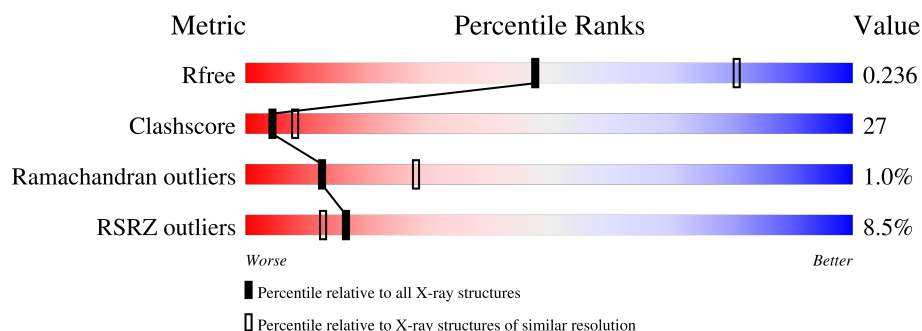
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (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 $\leq 5\%$. 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	337	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2481 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 Citramalate synthase from *Leptospira interrogans*.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	306	Total	C	N	O	S	0	0	0
			2415	1532	418	464	1			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	GLY	-	expression tag	UNP Q8F3Q1
A	-10	SER	-	expression tag	UNP Q8F3Q1
A	-9	HIS	-	expression tag	UNP Q8F3Q1
A	-8	MET	-	expression tag	UNP Q8F3Q1
A	-7	GLY	-	expression tag	UNP Q8F3Q1
A	-6	ARG	-	expression tag	UNP Q8F3Q1
A	-5	SER	-	expression tag	UNP Q8F3Q1
A	-4	GLN	-	expression tag	UNP Q8F3Q1
A	-3	LYS	-	expression tag	UNP Q8F3Q1
A	-2	VAL	-	expression tag	UNP Q8F3Q1
A	-1	SER	-	expression tag	UNP Q8F3Q1
A	0	GLN	-	expression tag	UNP Q8F3Q1

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is PYRUVIC ACID (CCD ID: PYR) (formula: C₃H₄O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		

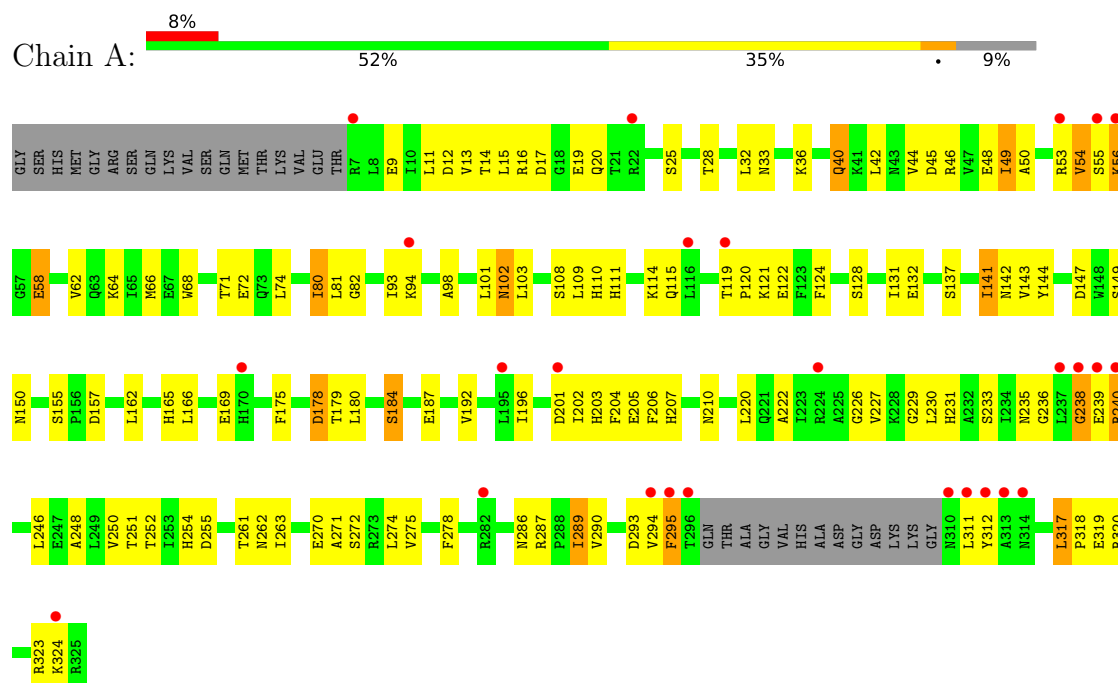
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	59	Total	O	0	0
			59	59		

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: Citramalate synthase from *Leptospira interrogans*



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	85.14Å 85.14Å 116.06Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.97 – 2.60 39.97 – 2.60	Depositor EDS
% Data completeness (in resolution range)	96.9 (39.97-2.60) 99.7 (39.97-2.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.223 , 0.259 0.228 , 0.236	Depositor DCC
R_{free} test set	777 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	(Not available)	Xtriage
Anisotropy	(Not available)	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 54.5	EDS
L-test for twinning ¹	$\langle L \rangle =$ (Not available), $\langle L^2 \rangle =$ (Not available)	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2481	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *(Not available)*

¹Theoretical 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: PYR, ZN

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	0.58	3/2457 (0.1%)	1.07	21/3318 (0.6%)

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

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	238	GLY	C-N	10.95	1.49	1.33
1	A	295	PHE	C-N	-8.73	1.21	1.33
1	A	233	SER	C-N	5.68	1.40	1.33

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	54	VAL	CA-C-O	-10.21	111.37	121.58
1	A	210	ASN	N-CA-C	7.84	121.28	112.97
1	A	53	ARG	CA-C-N	-7.79	109.52	120.35
1	A	53	ARG	C-N-CA	-7.79	109.52	120.35
1	A	293	ASP	CA-C-N	7.49	130.72	120.46
1	A	293	ASP	C-N-CA	7.49	130.72	120.46
1	A	184	SER	N-CA-C	-7.28	100.87	110.40
1	A	178	ASP	N-CA-C	-7.25	101.51	110.65
1	A	240	ARG	O-C-N	7.13	132.08	122.59
1	A	141	ILE	N-CA-C	6.44	117.39	108.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	58	GLU	N-CA-C	-6.15	104.63	111.71
1	A	49	ILE	N-CA-C	6.00	120.85	112.50
1	A	50	ALA	N-CA-C	5.94	116.26	107.88
1	A	143	VAL	N-CA-C	5.84	116.28	108.11
1	A	236	GLY	CA-C-N	-5.81	112.28	120.82
1	A	236	GLY	C-N-CA	-5.81	112.28	120.82
1	A	226	GLY	N-CA-C	5.79	121.36	114.48
1	A	40	GLN	N-CA-C	5.34	117.91	111.40
1	A	317	LEU	CA-C-N	5.30	125.37	119.32
1	A	317	LEU	C-N-CA	5.30	125.37	119.32
1	A	80	ILE	N-CA-C	5.20	115.34	107.75

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	238	GLY	Mainchain

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	2415	0	2425	129	0
2	A	1	0	0	0	0
3	A	6	0	0	0	0
4	A	59	0	0	4	0
All	All	2481	0	2425	129	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 27.

All (129) 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:A:102:ASN:HD21	1:A:142:ASN:HD22	1.18	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:261:THR:HG22	1:A:263:ILE:H	1.38	0.89
1:A:207:HIS:HD1	1:A:231:HIS:HD2	1.21	0.88
1:A:101:LEU:O	1:A:102:ASN:HB2	1.76	0.84
1:A:248:ALA:HA	1:A:289:ILE:HD12	1.61	0.83
1:A:207:HIS:HD1	1:A:231:HIS:CD2	1.98	0.80
1:A:42:LEU:HD22	1:A:263:ILE:HG23	1.64	0.79
1:A:102:ASN:ND2	1:A:142:ASN:HD22	1.81	0.79
1:A:124:PHE:HB3	1:A:165:HIS:CD2	2.18	0.79
1:A:102:ASN:ND2	1:A:142:ASN:HB2	2.01	0.76
1:A:261:THR:CG2	1:A:263:ILE:H	2.01	0.73
1:A:295:PHE:CZ	1:A:318:PRO:HA	2.24	0.73
1:A:80:ILE:O	1:A:101:LEU:O	2.12	0.68
1:A:295:PHE:CD2	1:A:318:PRO:HD3	2.27	0.68
1:A:16:ARG:HH21	1:A:20:GLN:HE22	1.41	0.67
1:A:16:ARG:HD3	1:A:48:GLU:OE1	1.95	0.67
1:A:13:VAL:HG12	1:A:17:ASP:HB2	1.77	0.65
1:A:49:ILE:HD11	1:A:62:VAL:HG13	1.79	0.64
1:A:110:HIS:O	1:A:114:LYS:HG2	1.97	0.63
1:A:147:ASP:OD2	1:A:150:ASN:HB2	1.98	0.63
1:A:287:ARG:O	1:A:294:VAL:HG21	1.99	0.62
1:A:28:THR:HG21	1:A:64:LYS:HD2	1.81	0.62
1:A:102:ASN:HD22	1:A:142:ASN:HB2	1.64	0.62
1:A:20:GLN:HB3	1:A:239:GLU:HG3	1.80	0.62
1:A:295:PHE:CE2	1:A:318:PRO:N	2.68	0.62
1:A:13:VAL:CG1	1:A:17:ASP:HB2	2.30	0.61
1:A:141:ILE:HD12	1:A:141:ILE:N	2.15	0.61
1:A:119:THR:HB	1:A:122:GLU:HB2	1.83	0.61
1:A:16:ARG:HD2	1:A:81:LEU:HB2	1.85	0.59
1:A:131:ILE:CD1	1:A:169:GLU:HG2	2.33	0.59
1:A:19:GLU:OE2	1:A:25:SER:HA	2.03	0.59
1:A:179:THR:O	1:A:240:ARG:NH2	2.33	0.59
1:A:207:HIS:ND1	1:A:231:HIS:HD2	1.97	0.58
1:A:36:LYS:HE2	1:A:68:TRP:CD1	2.38	0.58
1:A:58:GLU:O	1:A:62:VAL:HG23	2.04	0.57
1:A:40:GLN:NE2	1:A:72:GLU:HG2	2.19	0.57
1:A:192:VAL:HG11	1:A:227:VAL:HG23	1.87	0.57
1:A:240:ARG:NH2	4:A:1039:HOH:O	2.38	0.56
1:A:192:VAL:O	1:A:196:ILE:HD13	2.05	0.56
1:A:206:PHE:HB3	1:A:230:LEU:HD23	1.88	0.56
1:A:227:VAL:HG12	1:A:229:GLY:H	1.71	0.55
1:A:131:ILE:HD12	1:A:169:GLU:HG2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:235:ASN:ND2	1:A:275:VAL:HG22	2.22	0.54
1:A:235:ASN:HD22	1:A:275:VAL:HG22	1.73	0.54
1:A:32:LEU:HD22	1:A:36:LYS:HE3	1.91	0.53
1:A:32:LEU:CD2	1:A:36:LYS:HE3	2.39	0.53
1:A:157:ASP:HB2	4:A:1033:HOH:O	2.08	0.53
1:A:131:ILE:HG21	1:A:169:GLU:CG	2.38	0.52
1:A:205:GLU:HA	1:A:227:VAL:HG13	1.92	0.52
1:A:312:TYR:N	1:A:312:TYR:CD1	2.77	0.52
1:A:270:GLU:O	1:A:274:LEU:HB2	2.10	0.52
1:A:131:ILE:HG21	1:A:169:GLU:HG3	1.91	0.51
1:A:287:ARG:O	1:A:294:VAL:CG2	2.58	0.51
1:A:42:LEU:HD22	1:A:263:ILE:CG2	2.38	0.51
1:A:62:VAL:O	1:A:66:MET:HG2	2.11	0.51
1:A:204:PHE:O	1:A:227:VAL:HG13	2.10	0.51
1:A:254:HIS:NE2	1:A:261:THR:HB	2.27	0.50
1:A:124:PHE:HB3	1:A:165:HIS:HD2	1.75	0.50
1:A:33:ASN:ND2	1:A:274:LEU:HD21	2.27	0.50
1:A:72:GLU:HB3	1:A:74:LEU:HD22	1.92	0.50
1:A:286:ASN:ND2	4:A:1055:HOH:O	2.41	0.50
1:A:205:GLU:HA	1:A:227:VAL:CG1	2.42	0.50
1:A:323:ARG:HG2	1:A:324:LYS:H	1.77	0.49
1:A:144:TYR:HA	1:A:175:PHE:HB2	1.94	0.49
1:A:207:HIS:HB2	1:A:231:HIS:CD2	2.46	0.49
1:A:94:LYS:HE2	1:A:137:SER:HB3	1.95	0.49
1:A:317:LEU:O	1:A:320:ARG:HB3	2.13	0.49
1:A:101:LEU:O	1:A:102:ASN:CB	2.51	0.48
1:A:149:SER:HB3	1:A:178:ASP:HA	1.95	0.48
1:A:261:THR:HG22	1:A:263:ILE:N	2.17	0.48
1:A:11:LEU:HD13	1:A:13:VAL:HG22	1.95	0.48
1:A:11:LEU:CD1	1:A:13:VAL:HG22	2.44	0.48
1:A:272:SER:OG	1:A:290:VAL:HG11	2.14	0.47
1:A:207:HIS:O	1:A:207:HIS:CD2	2.67	0.47
1:A:261:THR:CG2	1:A:263:ILE:HG12	2.44	0.47
1:A:55:SER:O	1:A:56:LYS:C	2.58	0.47
1:A:54:VAL:HB	1:A:58:GLU:HG2	1.97	0.47
1:A:323:ARG:HG2	1:A:324:LYS:N	2.30	0.47
1:A:295:PHE:HD1	1:A:323:ARG:NH1	2.13	0.47
1:A:16:ARG:HG2	4:A:1021:HOH:O	2.15	0.46
1:A:207:HIS:HA	1:A:231:HIS:HB2	1.97	0.46
1:A:184:SER:OG	1:A:187:GLU:HG3	2.15	0.46
1:A:202:ILE:HB	1:A:204:PHE:CE1	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:255:ASP:CG	1:A:320:ARG:HH22	2.23	0.46
1:A:94:LYS:HE2	1:A:137:SER:CB	2.45	0.46
1:A:94:LYS:HE2	1:A:137:SER:OG	2.15	0.46
1:A:128:SER:O	1:A:132:GLU:HB2	2.14	0.46
1:A:261:THR:CG2	1:A:262:ASN:N	2.79	0.46
1:A:205:GLU:OE1	1:A:231:HIS:HE1	1.99	0.46
1:A:11:LEU:HD23	1:A:46:ARG:HB2	1.98	0.46
1:A:150:ASN:ND2	1:A:180:LEU:HD12	2.30	0.46
1:A:311:LEU:HB2	1:A:312:TYR:CE1	2.50	0.46
1:A:108:SER:OG	1:A:111:HIS:HB2	2.16	0.46
1:A:222:ALA:HB1	1:A:227:VAL:HG21	1.98	0.46
1:A:36:LYS:HE2	1:A:68:TRP:NE1	2.32	0.45
1:A:202:ILE:HG22	1:A:203:HIS:N	2.30	0.45
1:A:220:LEU:HD11	1:A:252:THR:CG2	2.47	0.44
1:A:40:GLN:HE22	1:A:72:GLU:HG2	1.81	0.44
1:A:201:ASP:OD2	1:A:201:ASP:N	2.46	0.44
1:A:206:PHE:HB3	1:A:230:LEU:CD2	2.47	0.44
1:A:119:THR:O	1:A:120:PRO:C	2.61	0.44
1:A:109:LEU:HB2	1:A:155:SER:CB	2.48	0.44
1:A:251:THR:OG1	1:A:289:ILE:HD13	2.17	0.43
1:A:33:ASN:HD21	1:A:274:LEU:HD21	1.82	0.43
1:A:9:GLU:HA	1:A:45:ASP:OD2	2.17	0.43
1:A:36:LYS:HG2	1:A:68:TRP:CZ2	2.53	0.43
1:A:162:LEU:O	1:A:162:LEU:HD23	2.19	0.43
1:A:246:LEU:HD13	1:A:250:VAL:HG23	2.01	0.42
1:A:295:PHE:HD1	1:A:323:ARG:HH11	1.67	0.42
1:A:82:GLY:HA3	1:A:101:LEU:HD21	2.01	0.42
1:A:15:LEU:O	1:A:19:GLU:HB2	2.19	0.42
1:A:124:PHE:CG	1:A:165:HIS:HD2	2.38	0.42
1:A:271:ALA:O	1:A:275:VAL:HG13	2.20	0.42
1:A:132:GLU:HA	1:A:132:GLU:OE1	2.20	0.42
1:A:68:TRP:O	1:A:71:THR:HB	2.20	0.42
1:A:33:ASN:ND2	1:A:278:PHE:HE1	2.17	0.41
1:A:93:ILE:HG23	1:A:98:ALA:HB3	2.01	0.41
1:A:111:HIS:CD2	1:A:115:GLN:HE22	2.37	0.41
1:A:44:VAL:HA	1:A:263:ILE:HD11	2.02	0.41
1:A:121:LYS:HB3	1:A:121:LYS:HE2	1.86	0.41
1:A:103:LEU:N	1:A:142:ASN:O	2.51	0.41
1:A:124:PHE:CB	1:A:165:HIS:CD2	3.00	0.41
1:A:207:HIS:ND1	1:A:231:HIS:CD2	2.78	0.41
1:A:12:ASP:OD2	1:A:14:THR:OG1	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:162:LEU:HD23	1:A:166:LEU:HG	2.03	0.40
1:A:202:ILE:CG2	1:A:203:HIS:N	2.84	0.40
1:A:227:VAL:CG1	1:A:229:GLY:O	2.69	0.40
1:A:295:PHE:CE2	1:A:317:LEU:C	3.00	0.40
1:A:317:LEU:HB3	1:A:319:GLU:OE2	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	302/337 (90%)	286 (95%)	13 (4%)	3 (1%)	12 28

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	56	LYS
1	A	289	ILE
1	A	102	ASN

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\# Z > 2$
3	PYR	A	1004	2	5,5,5	2.66	1 (20%)	3,6,6	0.82	0

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
3	PYR	A	1004	2	-	0/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1004	PYR	CB-CA	-5.52	1.39	1.50

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	306/337 (90%)	0.58	26 (8%) 16 13	24, 49, 100, 164	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	296	THR	5.5
1	A	295	PHE	4.8
1	A	311	LEU	4.6
1	A	310	ASN	4.3
1	A	312	TYR	4.3
1	A	239	GLU	4.1
1	A	313	ALA	3.8
1	A	53	ARG	3.8
1	A	7	ARG	3.7
1	A	56	LYS	3.5
1	A	294	VAL	3.0
1	A	116	LEU	2.6
1	A	282	ARG	2.5
1	A	201	ASP	2.4
1	A	170	HIS	2.4
1	A	195	LEU	2.4
1	A	119	THR	2.3
1	A	314	ASN	2.3
1	A	22	ARG	2.3
1	A	238	GLY	2.3
1	A	324	LYS	2.3
1	A	55	SER	2.3
1	A	224	ARG	2.3
1	A	94	LYS	2.2
1	A	237	LEU	2.0
1	A	240	ARG	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($\text{\AA}^2$)	Q<0.9
3	PYR	A	1004	6/6	0.93	0.15	54,56,58,59	0
2	ZN	A	1003	1/1	0.98	0.10	37,37,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.