



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 05:23 AM UTC

PDB ID : 8IAT / pdb_00008iat
Title : Crystal structure of Streptococcus pneumoniae pyruvate kinase in complex with oxalate
Authors : Nakashima, R.; Taguchi, A.
Deposited on : 2023-02-09
Resolution : 1.80 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

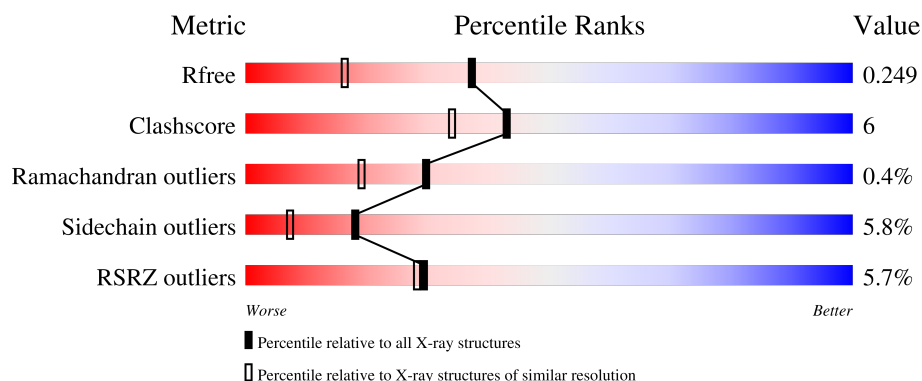
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	521	2% 85% 10% • •
1	B	521	4% 79% 15% • •
1	C	521	11% 75% 20% • •
1	D	521	6% 81% 12% • •

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	OXL	A	601	-	X	-	-
2	OXL	A	602	-	X	-	-
2	OXL	B	601	-	X	-	-
2	OXL	B	602	-	X	-	-
2	OXL	C	602	-	X	-	-
2	OXL	D	601	-	X	-	-
2	OXL	D	602	-	X	-	-
7	GOL	A	607	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 16483 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 Pyruvate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	501	Total	C	N	O	S	0	1	0
			3851	2408	665	761	17			
1	B	501	Total	C	N	O	S	0	1	0
			3851	2408	665	761	17			
1	C	501	Total	C	N	O	S	0	0	0
			3839	2399	664	759	17			
1	D	500	Total	C	N	O	S	0	0	0
			3828	2393	660	758	17			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP Q8DQ84
A	-18	GLY	-	expression tag	UNP Q8DQ84
A	-17	SER	-	expression tag	UNP Q8DQ84
A	-16	SER	-	expression tag	UNP Q8DQ84
A	-15	HIS	-	expression tag	UNP Q8DQ84
A	-14	HIS	-	expression tag	UNP Q8DQ84
A	-13	HIS	-	expression tag	UNP Q8DQ84
A	-12	HIS	-	expression tag	UNP Q8DQ84
A	-11	HIS	-	expression tag	UNP Q8DQ84
A	-10	HIS	-	expression tag	UNP Q8DQ84
A	-9	SER	-	expression tag	UNP Q8DQ84
A	-8	SER	-	expression tag	UNP Q8DQ84
A	-7	GLY	-	expression tag	UNP Q8DQ84
A	-6	LEU	-	expression tag	UNP Q8DQ84
A	-5	VAL	-	expression tag	UNP Q8DQ84
A	-4	PRO	-	expression tag	UNP Q8DQ84
A	-3	ARG	-	expression tag	UNP Q8DQ84
A	-2	GLY	-	expression tag	UNP Q8DQ84
A	-1	SER	-	expression tag	UNP Q8DQ84
A	0	HIS	-	expression tag	UNP Q8DQ84
B	-19	MET	-	initiating methionine	UNP Q8DQ84

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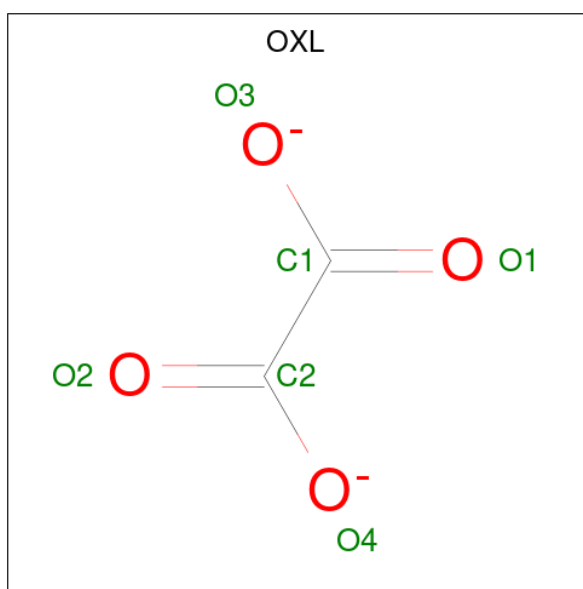
Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	expression tag	UNP Q8DQ84
B	-17	SER	-	expression tag	UNP Q8DQ84
B	-16	SER	-	expression tag	UNP Q8DQ84
B	-15	HIS	-	expression tag	UNP Q8DQ84
B	-14	HIS	-	expression tag	UNP Q8DQ84
B	-13	HIS	-	expression tag	UNP Q8DQ84
B	-12	HIS	-	expression tag	UNP Q8DQ84
B	-11	HIS	-	expression tag	UNP Q8DQ84
B	-10	HIS	-	expression tag	UNP Q8DQ84
B	-9	SER	-	expression tag	UNP Q8DQ84
B	-8	SER	-	expression tag	UNP Q8DQ84
B	-7	GLY	-	expression tag	UNP Q8DQ84
B	-6	LEU	-	expression tag	UNP Q8DQ84
B	-5	VAL	-	expression tag	UNP Q8DQ84
B	-4	PRO	-	expression tag	UNP Q8DQ84
B	-3	ARG	-	expression tag	UNP Q8DQ84
B	-2	GLY	-	expression tag	UNP Q8DQ84
B	-1	SER	-	expression tag	UNP Q8DQ84
B	0	HIS	-	expression tag	UNP Q8DQ84
C	-19	MET	-	initiating methionine	UNP Q8DQ84
C	-18	GLY	-	expression tag	UNP Q8DQ84
C	-17	SER	-	expression tag	UNP Q8DQ84
C	-16	SER	-	expression tag	UNP Q8DQ84
C	-15	HIS	-	expression tag	UNP Q8DQ84
C	-14	HIS	-	expression tag	UNP Q8DQ84
C	-13	HIS	-	expression tag	UNP Q8DQ84
C	-12	HIS	-	expression tag	UNP Q8DQ84
C	-11	HIS	-	expression tag	UNP Q8DQ84
C	-10	HIS	-	expression tag	UNP Q8DQ84
C	-9	SER	-	expression tag	UNP Q8DQ84
C	-8	SER	-	expression tag	UNP Q8DQ84
C	-7	GLY	-	expression tag	UNP Q8DQ84
C	-6	LEU	-	expression tag	UNP Q8DQ84
C	-5	VAL	-	expression tag	UNP Q8DQ84
C	-4	PRO	-	expression tag	UNP Q8DQ84
C	-3	ARG	-	expression tag	UNP Q8DQ84
C	-2	GLY	-	expression tag	UNP Q8DQ84
C	-1	SER	-	expression tag	UNP Q8DQ84
C	0	HIS	-	expression tag	UNP Q8DQ84
D	-19	MET	-	initiating methionine	UNP Q8DQ84
D	-18	GLY	-	expression tag	UNP Q8DQ84
D	-17	SER	-	expression tag	UNP Q8DQ84

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	expression tag	UNP Q8DQ84
D	-15	HIS	-	expression tag	UNP Q8DQ84
D	-14	HIS	-	expression tag	UNP Q8DQ84
D	-13	HIS	-	expression tag	UNP Q8DQ84
D	-12	HIS	-	expression tag	UNP Q8DQ84
D	-11	HIS	-	expression tag	UNP Q8DQ84
D	-10	HIS	-	expression tag	UNP Q8DQ84
D	-9	SER	-	expression tag	UNP Q8DQ84
D	-8	SER	-	expression tag	UNP Q8DQ84
D	-7	GLY	-	expression tag	UNP Q8DQ84
D	-6	LEU	-	expression tag	UNP Q8DQ84
D	-5	VAL	-	expression tag	UNP Q8DQ84
D	-4	PRO	-	expression tag	UNP Q8DQ84
D	-3	ARG	-	expression tag	UNP Q8DQ84
D	-2	GLY	-	expression tag	UNP Q8DQ84
D	-1	SER	-	expression tag	UNP Q8DQ84
D	0	HIS	-	expression tag	UNP Q8DQ84

- Molecule 2 is OXALATE ION (CCD ID: OXL) (formula: C_2O_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	2	4		
2	A	1	Total	C	O	0	0
			6	2	4		
2	B	1	Total	C	O	0	0
			6	2	4		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total C O 6 2 4	0	0
2	C	1	Total C O 6 2 4	0	0
2	C	1	Total C O 6 2 4	0	0
2	D	1	Total C O 6 2 4	0	0
2	D	1	Total C O 6 2 4	0	0

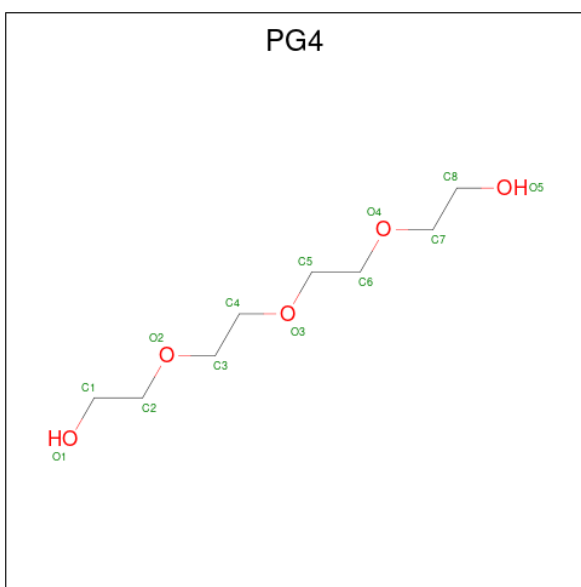
- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 4 is POTASSIUM ION (CCD ID: K) (formula: K).

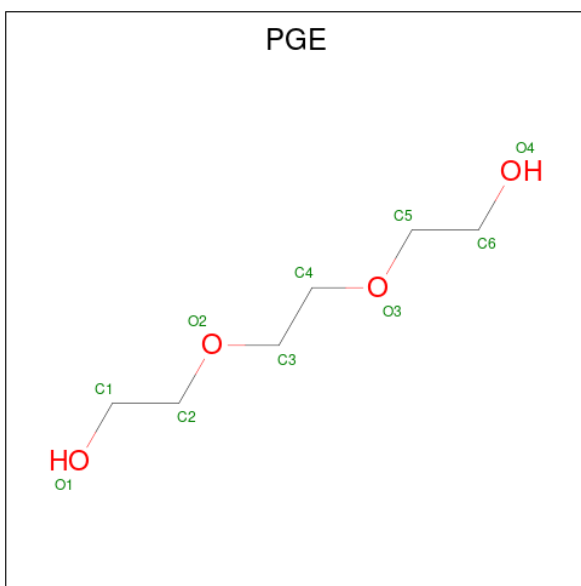
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total K 1 1	0	0
4	B	1	Total K 1 1	0	0
4	C	1	Total K 1 1	0	0
4	D	1	Total K 1 1	0	0

- Molecule 5 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



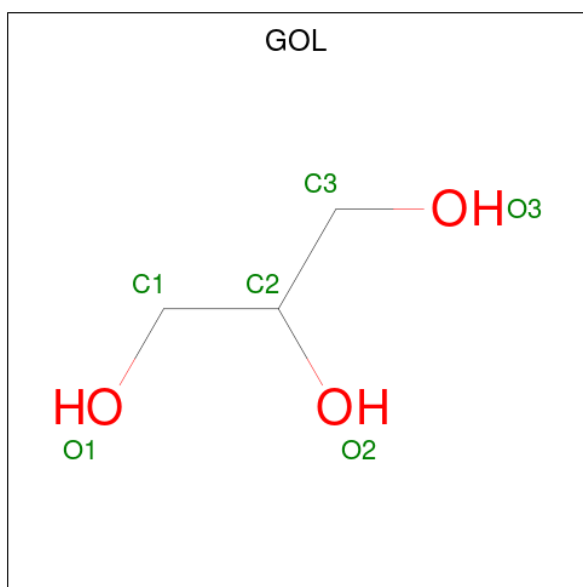
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			13	8	5		

- Molecule 6 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			10	6	4		
6	C	1	Total	C	O	0	0
			10	6	4		

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

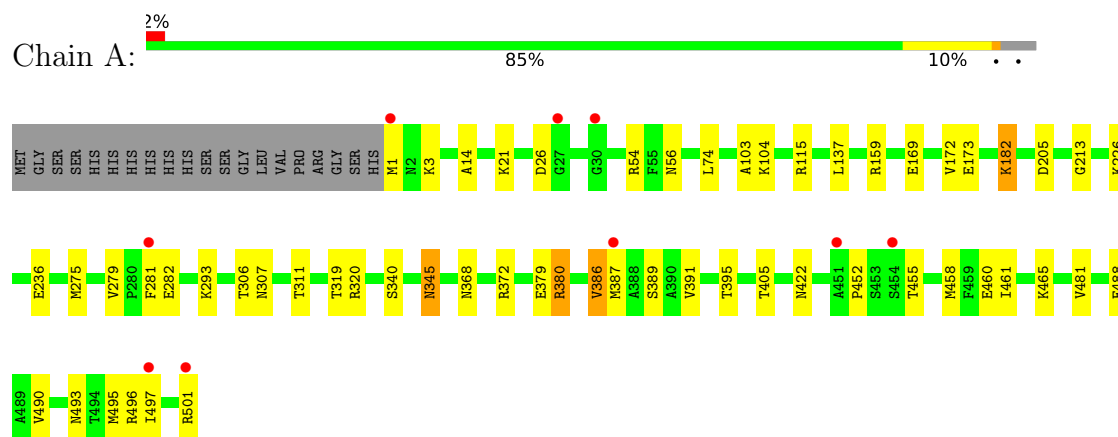
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	325	Total	O	0	0
			325	325		
8	B	322	Total	O	0	0
			322	322		
8	C	195	Total	O	0	0
			195	195		
8	D	177	Total	O	0	0
			177	177		

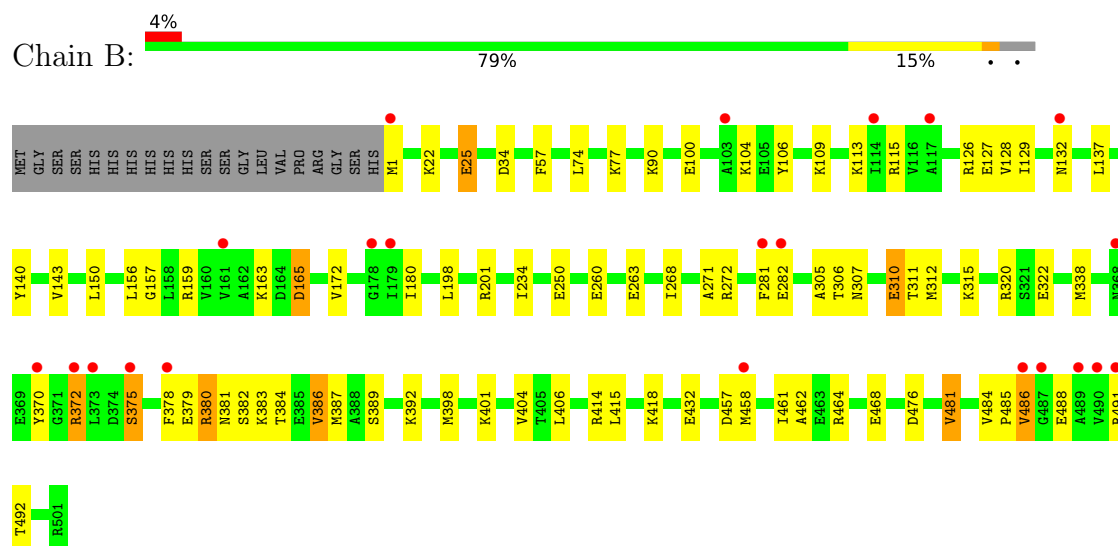
3 Residue-property plots [i](#)

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.

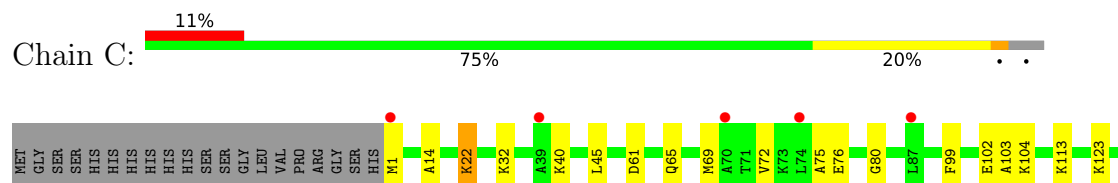
• Molecule 1: Pyruvate kinase

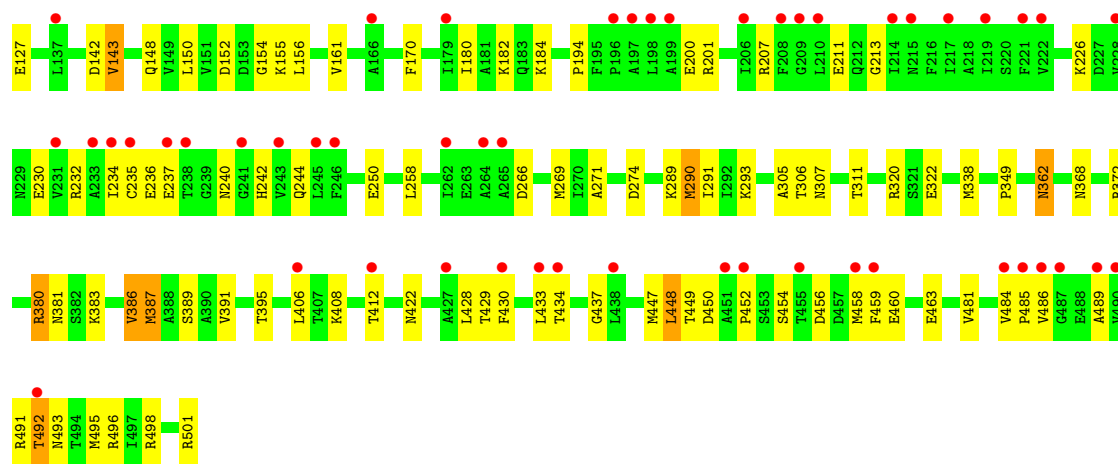


• Molecule 1: Pyruvate kinase

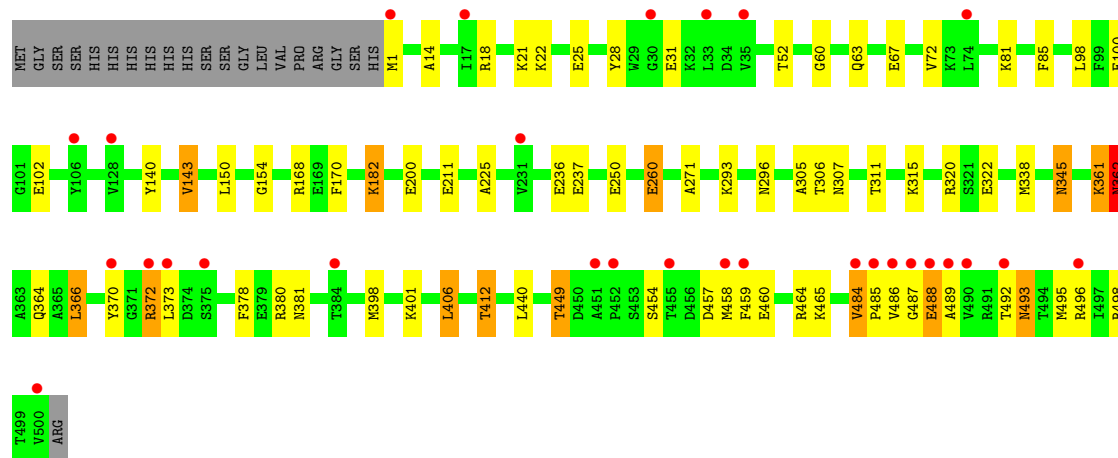
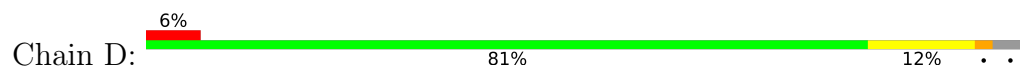


• Molecule 1: Pyruvate kinase





• Molecule 1: Pyruvate kinase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	212.05Å 118.35Å 99.82Å 90.00° 108.93° 90.00°	Depositor
Resolution (Å)	48.30 – 1.80 48.30 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.30-1.80) 99.9 (48.30-1.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, R_{free}	0.199 , 0.246 0.205 , 0.249	Depositor DCC
R_{free} test set	10764 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	36.2	Xtriage
Anisotropy	0.356	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	16483	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.00% 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: GOL, PGE, MG, PG4, OXL, K

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.07	0/3897	1.34	3/5254 (0.1%)
1	B	1.09	1/3897 (0.0%)	1.33	2/5254 (0.0%)
1	C	1.10	0/3884	1.43	8/5238 (0.2%)
1	D	1.05	0/3873	1.38	3/5224 (0.1%)
All	All	1.08	1/15551 (0.0%)	1.37	16/20970 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	268	ILE	C-O	5.51	1.29	1.24

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	362	ASN	CA-CB-CG	-7.05	105.55	112.60
1	C	61	ASP	CA-CB-CG	5.97	118.58	112.60
1	C	434	THR	CA-C-N	5.64	127.84	120.28
1	C	434	THR	C-N-CA	5.64	127.84	120.28
1	A	103	ALA	N-CA-C	-5.54	102.82	110.35
1	D	412	THR	CA-CB-OG1	-5.47	101.40	109.60
1	D	362	ASN	CA-CB-CG	-5.40	107.20	112.60
1	B	414	ARG	CG-CD-NE	-5.32	100.30	112.00
1	A	205	ASP	CA-CB-CG	5.20	117.80	112.60
1	B	34	ASP	CA-CB-CG	5.19	117.79	112.60
1	A	379	GLU	CB-CA-C	5.19	118.30	109.53
1	D	60	GLY	CA-C-O	-5.16	118.74	122.45
1	C	489	ALA	CA-C-N	5.06	127.13	120.60
1	C	489	ALA	C-N-CA	5.06	127.13	120.60
1	C	289	LYS	CA-C-N	5.04	126.99	120.44
1	C	289	LYS	C-N-CA	5.04	126.99	120.44

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	3851	0	3895	41	0
1	B	3851	0	3895	49	0
1	C	3839	0	3887	71	0
1	D	3828	0	3874	50	0
2	A	12	0	0	0	0
2	B	12	0	0	1	0
2	C	12	0	0	2	0
2	D	12	0	0	1	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	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	13	0	18	0	0
6	A	10	0	14	1	0
6	C	10	0	14	1	0
7	A	6	0	8	6	0
8	A	325	0	0	3	0
8	B	322	0	0	3	0
8	C	195	0	0	8	0
8	D	177	0	0	3	0
All	All	16483	0	15605	192	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 6.

All (192) 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:296:ASN:HB2	8:D:840:HOH:O	1.47	1.14
1:A:54:ARG:HH12	7:A:607:GOL:H11	1.27	0.95
1:C:237:GLU:HG3	8:C:707:HOH:O	1.77	0.82
1:A:54:ARG:HH12	7:A:607:GOL:C1	1.92	0.82
1:A:320:ARG:HD2	1:D:307:ASN:HD22	1.46	0.78
1:C:236:GLU:HB2	8:C:707:HOH:O	1.86	0.75
1:B:380:ARG:NH2	1:B:389:SER:OG	2.20	0.75
1:C:22:LYS:HD2	1:C:22:LYS:N	2.03	0.73
1:A:282:GLU:HG3	1:D:362:ASN:ND2	2.04	0.73
1:D:458:MET:HE1	1:D:459:PHE:CE1	2.23	0.73
1:A:282:GLU:HG3	1:D:362:ASN:HD22	1.52	0.72
1:A:54:ARG:NH1	7:A:607:GOL:H11	2.04	0.71
1:C:148:GLN:HE22	6:C:605:PGE:H3	1.58	0.69
1:C:232:ARG:HD3	8:C:822:HOH:O	1.91	0.69
1:C:142:ASP:OD2	1:C:194:PRO:CD	2.42	0.67
1:D:459:PHE:CD2	1:D:496:ARG:NH1	2.63	0.67
1:D:484:VAL:HA	1:D:485:PRO:C	2.18	0.67
1:D:143:VAL:HG22	1:D:170:PHE:CE1	2.30	0.66
1:B:281[B]:PHE:CD1	1:B:282:GLU:N	2.64	0.66
1:C:492:THR:HG21	1:D:498:ARG:NH2	2.10	0.66
1:C:387:MET:SD	1:C:493:ASN:OD1	2.55	0.65
1:A:307:ASN:HD22	1:D:320:ARG:HD2	1.61	0.64
1:C:492:THR:HG21	1:D:498:ARG:HH21	1.63	0.64
1:B:77:LYS:NZ	8:B:701:HOH:O	2.32	0.63
1:C:498:ARG:NH1	1:C:501:ARG:HG3	2.15	0.61
1:A:455:THR:O	1:A:458:MET:HB3	2.00	0.61
1:D:493:ASN:C	1:D:493:ASN:HD22	2.09	0.60
1:A:391:VAL:O	1:A:395:THR:HG23	2.01	0.60
1:B:320:ARG:HD2	1:C:307:ASN:HD22	1.66	0.60
1:A:340:SER:HB3	7:A:607:GOL:H12	1.83	0.60
1:C:258:LEU:CD2	1:C:290:MET:HE1	2.32	0.60
1:D:21:LYS:HD2	1:D:31:GLU:HG3	1.83	0.59
1:D:406:LEU:HD13	1:D:458:MET:HG2	1.85	0.59
1:C:491:ARG:C	1:C:492:THR:HG23	2.28	0.58
1:B:115:ARG:HB2	1:B:128:VAL:HG22	1.85	0.58
1:C:142:ASP:OD2	1:C:194:PRO:HD2	2.03	0.58
1:B:281[B]:PHE:CE1	1:B:282:GLU:HB3	2.38	0.58
1:C:412:THR:HG23	2:C:602:OXL:O2	2.03	0.57
1:A:387:MET:HE2	1:A:493:ASN:OD1	2.05	0.57
1:A:372:ARG:HH21	1:D:372:ARG:NH1	2.03	0.56
1:B:106:TYR:CG	1:B:126:ARG:HG3	2.41	0.56
1:C:463:GLU:OE2	1:C:498:ARG:NE	2.32	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:458:MET:HB2	1:D:486:VAL:HG21	1.88	0.56
1:C:290:MET:HE2	1:C:290:MET:C	2.31	0.56
1:B:406:LEU:HD13	1:B:458:MET:HB3	1.87	0.56
1:C:240:ASN:C	1:C:242:HIS:HD2	2.13	0.55
1:A:14:ALA:HB2	1:A:345:ASN:HA	1.88	0.55
1:B:281[B]:PHE:CG	1:B:282:GLU:N	2.74	0.55
1:B:379:GLU:CD	1:B:381:ASN:HD21	2.15	0.55
1:B:384:THR:HG22	1:B:415:LEU:HD12	1.88	0.55
1:D:492:THR:HG21	8:D:743:HOH:O	2.07	0.55
1:B:198:LEU:HD21	1:B:234:ILE:HD11	1.88	0.54
1:B:464:ARG:O	1:B:468:GLU:HG3	2.07	0.54
1:A:380:ARG:HB3	1:A:386:VAL:HG12	1.90	0.54
1:D:63:GLN:NE2	1:D:67:GLU:OE1	2.41	0.54
1:C:460:GLU:HG3	8:C:701:HOH:O	2.07	0.54
1:A:490:VAL:HG23	8:A:873:HOH:O	2.07	0.54
1:C:429:THR:O	1:C:448:LEU:HD12	2.07	0.54
1:B:401:LYS:HE2	1:B:476:ASP:OD2	2.08	0.53
1:C:383:LYS:HA	1:C:386:VAL:CG1	2.38	0.53
1:C:452:PRO:HG3	1:C:486:VAL:HG21	1.90	0.53
1:C:305:ALA:HB1	1:C:338:MET:HE2	1.90	0.53
1:D:361:LYS:O	1:D:364:GLN:HG2	2.08	0.53
1:A:387:MET:CE	1:A:493:ASN:OD1	2.56	0.53
1:D:449:THR:OG1	1:D:465:LYS:HE3	2.08	0.53
1:C:458:MET:HE1	1:C:459:PHE:CZ	2.44	0.53
1:B:305:ALA:HB1	1:B:338:MET:HE2	1.90	0.53
1:D:412:THR:HG21	1:D:492:THR:HG21	1.91	0.52
1:C:65:GLN:O	1:C:69:MET:HG3	2.09	0.52
1:A:21:LYS:HE3	1:A:26:ASP:O	2.09	0.52
1:B:22:LYS:O	1:B:25:GLU:HB2	2.09	0.52
1:D:311:THR:HG23	1:D:322:GLU:OE1	2.09	0.52
1:B:404:VAL:HG11	1:B:462:ALA:HB1	1.92	0.52
1:B:150:LEU:HD23	1:B:157:GLY:HA2	1.92	0.51
1:D:366:LEU:HD23	1:D:370:TYR:CD2	2.45	0.51
1:B:156:LEU:HD23	1:B:157:GLY:N	2.26	0.51
1:C:240:ASN:HA	1:C:242:HIS:CD2	2.45	0.51
1:C:406:LEU:HD22	1:C:458:MET:HG2	1.91	0.51
1:A:380:ARG:NH2	1:A:389:SER:OG	2.45	0.50
1:B:156:LEU:HD23	1:B:156:LEU:C	2.35	0.50
1:C:387:MET:CE	1:D:495:MET:HE2	2.40	0.50
1:A:452:PRO:HB3	1:A:461:ILE:HD12	1.92	0.50
1:B:481:VAL:HG13	1:B:492:THR:HG21	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:405:THR:HA	1:A:481:VAL:O	2.12	0.50
1:C:456:ASP:O	8:C:701:HOH:O	2.20	0.50
1:D:225:ALA:HB2	1:D:260:GLU:HG3	1.93	0.50
1:A:391:VAL:HG21	1:A:481:VAL:HG21	1.94	0.49
1:A:452:PRO:HB3	1:A:461:ILE:CD1	2.42	0.49
1:B:307:ASN:HD22	1:C:320:ARG:HD2	1.77	0.49
1:D:22:LYS:O	1:D:28:TYR:CD2	2.65	0.49
1:D:449:THR:OG1	1:D:465:LYS:CE	2.60	0.49
1:C:240:ASN:O	1:C:242:HIS:CD2	2.65	0.49
1:D:14:ALA:HB2	1:D:345:ASN:HA	1.94	0.49
1:B:281[B]:PHE:CD1	1:B:281[B]:PHE:C	2.91	0.49
1:C:458:MET:HE1	1:C:459:PHE:CE1	2.48	0.49
1:C:22:LYS:N	1:C:22:LYS:CD	2.75	0.49
1:C:386:VAL:HG11	1:D:398:MET:SD	2.53	0.49
1:C:311:THR:HG23	1:C:322:GLU:OE1	2.13	0.48
1:A:104:LYS:HD2	1:A:182:LYS:HE3	1.94	0.48
1:A:387:MET:SD	1:A:495:MET:HB2	2.53	0.48
1:B:372:ARG:HH12	1:C:372:ARG:HG2	1.77	0.48
1:C:391:VAL:HG21	1:C:481:VAL:HG21	1.96	0.48
1:B:250:GLU:HG2	1:B:271:ALA:HB3	1.94	0.48
1:B:370:TYR:O	1:B:372:ARG:NH1	2.45	0.48
1:C:496:ARG:HA	1:D:493:ASN:O	2.14	0.48
1:C:213:GLY:HA2	8:C:796:HOH:O	2.13	0.47
1:B:375:SER:OG	1:B:392:LYS:NZ	2.46	0.47
1:B:271:ALA:HB1	2:B:601:OXL:C1	2.44	0.47
1:C:242:HIS:CE1	1:C:433:LEU:HD22	2.49	0.47
1:D:492:THR:CG2	8:D:743:HOH:O	2.62	0.47
1:C:240:ASN:C	1:C:242:HIS:CD2	2.92	0.47
1:C:207:ARG:O	1:C:211:GLU:HG3	2.15	0.47
1:C:230:GLU:O	1:C:234:ILE:HD13	2.14	0.47
1:C:428:LEU:HD23	1:C:447:MET:HB3	1.97	0.47
1:D:305:ALA:HB1	1:D:338:MET:HE2	1.95	0.47
1:B:457:ASP:O	1:B:461:ILE:HG13	2.15	0.46
1:A:372:ARG:HG2	1:D:372:ARG:HH12	1.79	0.46
1:D:296:ASN:HD21	1:D:378:PHE:HZ	1.62	0.46
1:B:159:ARG:O	1:B:172:VAL:HA	2.16	0.46
1:B:312:MET:HA	1:B:315:LYS:O	2.16	0.46
1:B:129:ILE:HD12	1:B:180:ILE:HG21	1.98	0.46
1:C:244:GLN:NE2	1:C:437:GLY:O	2.48	0.46
1:C:152:ASP:O	1:C:155:LYS:HG2	2.15	0.46
1:A:56:ASN:HD22	7:A:607:GOL:H31	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:76:GLU:O	1:C:80:GLY:N	2.46	0.45
1:D:406:LEU:HD13	1:D:458:MET:CG	2.46	0.45
1:A:213:GLY:C	6:A:606:PGE:H3	2.41	0.45
1:B:140:TYR:OH	1:B:163:LYS:HB3	2.16	0.45
1:B:165:ASP:OD2	1:B:165:ASP:N	2.49	0.45
1:B:90:LYS:HD3	1:B:201:ARG:HD2	1.98	0.45
1:C:391:VAL:O	1:C:395:THR:HG23	2.16	0.45
1:C:150:LEU:HB3	1:C:154:GLY:HA2	1.99	0.45
1:B:310:GLU:HG3	8:B:752:HOH:O	2.17	0.44
1:D:140:TYR:CE2	1:D:168:ARG:HA	2.52	0.44
1:D:1:MET:HE2	1:D:1:MET:HA	1.99	0.44
1:D:487:GLY:C	1:D:489:ALA:H	2.25	0.44
1:D:250:GLU:HB3	1:D:271:ALA:HB3	1.99	0.44
1:C:368:ASN:HD21	1:C:422:ASN:ND2	2.15	0.44
1:B:382:SER:O	1:B:386:VAL:HG13	2.17	0.44
1:B:57:PHE:O	1:B:201:ARG:NH2	2.50	0.44
1:A:386:VAL:HG11	1:B:398:MET:HG3	1.99	0.43
1:B:90:LYS:HD3	1:B:201:ARG:CD	2.48	0.43
1:D:98:LEU:HD23	1:D:182:LYS:HG3	2.00	0.43
1:A:3:LYS:NZ	8:A:714:HOH:O	2.49	0.43
1:A:461:ILE:O	1:A:465:LYS:HG2	2.18	0.43
1:A:159:ARG:O	1:A:172:VAL:HA	2.19	0.43
1:A:173:GLU:HG2	8:A:876:HOH:O	2.18	0.43
1:A:496:ARG:NH2	1:B:484:VAL:HG11	2.33	0.43
1:C:184:LYS:NZ	8:C:716:HOH:O	2.52	0.43
1:A:497:ILE:HG13	1:B:387:MET:HE1	2.00	0.43
1:B:310:GLU:CG	8:B:752:HOH:O	2.67	0.43
1:B:485:PRO:HB2	1:B:486:VAL:HG22	1.99	0.43
1:D:18:ARG:HB3	1:D:31:GLU:HB2	2.01	0.43
1:C:430:PHE:HZ	1:C:486:VAL:HG11	1.83	0.43
1:D:22:LYS:O	1:D:28:TYR:HD2	1.99	0.43
1:B:311:THR:HG23	1:B:322:GLU:OE1	2.19	0.43
1:C:143:VAL:HG13	1:C:170:PHE:CZ	2.54	0.42
1:D:143:VAL:CG2	1:D:170:PHE:CE1	3.01	0.42
1:B:106:TYR:CD1	1:B:126:ARG:HG3	2.54	0.42
1:C:156:LEU:HD23	1:C:180:ILE:CG1	2.49	0.42
1:A:275:MET:O	1:A:279:VAL:HG22	2.20	0.42
1:C:142:ASP:OD2	1:C:194:PRO:HD3	2.17	0.42
1:C:484:VAL:HA	1:C:485:PRO:HA	1.94	0.42
1:C:99:PHE:HB2	1:C:103:ALA:O	2.18	0.42
1:C:380:ARG:NH2	1:C:389:SER:OG	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:387:MET:HE2	1:C:495:MET:HB2	2.00	0.42
1:C:430:PHE:C	1:C:448:LEU:HD11	2.45	0.42
1:D:52:THR:HB	1:D:440:LEU:HD21	2.01	0.42
1:D:72:VAL:HG12	1:D:85:PHE:CZ	2.55	0.42
1:C:250:GLU:HB3	1:C:274:ASP:HB2	2.02	0.41
1:A:368:ASN:OD1	1:A:422:ASN:ND2	2.53	0.41
1:C:14:ALA:HB1	1:C:349:PRO:HG3	2.02	0.41
1:C:72:VAL:O	1:C:75:ALA:HB3	2.20	0.41
1:C:271:ALA:HB1	2:C:601:OXL:C2	2.50	0.41
1:C:290:MET:HE2	1:C:291:ILE:N	2.35	0.41
1:D:150:LEU:HB3	1:D:154:GLY:HA2	2.01	0.41
1:D:271:ALA:HB1	2:D:601:OXL:C1	2.50	0.41
1:B:384:THR:CG2	1:B:415:LEU:HD12	2.51	0.41
1:C:387:MET:HE1	1:D:495:MET:HE2	2.01	0.41
1:C:236:GLU:CB	8:C:707:HOH:O	2.58	0.41
1:A:56:ASN:HD22	7:A:607:GOL:C3	2.33	0.41
1:A:104:LYS:HD2	1:A:182:LYS:CE	2.50	0.41
1:A:115:ARG:HD3	1:A:169:GLU:OE2	2.20	0.41
1:C:269:MET:HE2	1:C:338:MET:SD	2.61	0.41
1:D:454:SER:H	1:D:457:ASP:HB3	1.86	0.41
1:C:244:GLN:HA	1:C:266:ASP:OD2	2.21	0.41
1:B:378:PHE:CE2	1:B:418:LYS:HG2	2.55	0.41
1:C:235:CYS:O	1:C:240:ASN:C	2.64	0.40
1:B:415:LEU:HD23	1:B:415:LEU:HA	1.89	0.40
1:C:258:LEU:HD23	1:C:290:MET:HE1	2.03	0.40
1:A:282:GLU:CG	1:D:362:ASN:ND2	2.81	0.40
1:A:311:THR:HG21	1:A:319:THR:CG2	2.52	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	500/521 (96%)	491 (98%)	8 (2%)	1 (0%)	43	31
1	B	500/521 (96%)	480 (96%)	18 (4%)	2 (0%)	30	19
1	C	499/521 (96%)	481 (96%)	15 (3%)	3 (1%)	21	11
1	D	498/521 (96%)	480 (96%)	16 (3%)	2 (0%)	30	19
All	All	1997/2084 (96%)	1932 (97%)	57 (3%)	8 (0%)	30	19

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	488	GLU
1	C	306	THR
1	D	306	THR
1	A	306	THR
1	B	306	THR
1	C	492	THR
1	D	488	GLU
1	C	381	ASN

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	412/428 (96%)	397 (96%)	15 (4%)	31	18
1	B	412/428 (96%)	387 (94%)	25 (6%)	17	6
1	C	411/428 (96%)	384 (93%)	27 (7%)	15	5
1	D	410/428 (96%)	381 (93%)	29 (7%)	13	4
All	All	1645/1712 (96%)	1549 (94%)	96 (6%)	18	7

All (96) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	74	LEU
1	A	137	LEU

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Mol	Chain	Res	Type
1	A	182	LYS
1	A	226	LYS
1	A	236	GLU
1	A	281[A]	PHE
1	A	281[B]	PHE
1	A	293	LYS
1	A	345	ASN
1	A	380	ARG
1	A	386	VAL
1	A	460	GLU
1	A	488	GLU
1	A	501	ARG
1	B	1	MET
1	B	25	GLU
1	B	74	LEU
1	B	100	GLU
1	B	104	LYS
1	B	109	LYS
1	B	113	LYS
1	B	127	GLU
1	B	132	ASN
1	B	137	LEU
1	B	143	VAL
1	B	165	ASP
1	B	260	GLU
1	B	263	GLU
1	B	272	ARG
1	B	310	GLU
1	B	372	ARG
1	B	375	SER
1	B	380	ARG
1	B	383	LYS
1	B	386	VAL
1	B	432	GLU
1	B	481	VAL
1	B	486	VAL
1	B	491	ARG
1	C	1	MET
1	C	22	LYS
1	C	32	LYS
1	C	40	LYS
1	C	45	LEU

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Mol	Chain	Res	Type
1	C	102	GLU
1	C	104	LYS
1	C	113	LYS
1	C	123	LYS
1	C	127	GLU
1	C	143	VAL
1	C	161	VAL
1	C	182	LYS
1	C	200	GLU
1	C	201	ARG
1	C	226	LYS
1	C	290	MET
1	C	293	LYS
1	C	362	ASN
1	C	380	ARG
1	C	386	VAL
1	C	387	MET
1	C	408	LYS
1	C	448	LEU
1	C	449	THR
1	C	450	ASP
1	C	454	SER
1	D	25	GLU
1	D	81	LYS
1	D	100	GLU
1	D	102	GLU
1	D	143	VAL
1	D	182	LYS
1	D	200	GLU
1	D	211	GLU
1	D	236	GLU
1	D	237	GLU
1	D	260	GLU
1	D	293	LYS
1	D	315	LYS
1	D	345	ASN
1	D	361	LYS
1	D	362	ASN
1	D	366	LEU
1	D	372	ARG
1	D	373	LEU
1	D	380	ARG

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Mol	Chain	Res	Type
1	D	381	ASN
1	D	401	LYS
1	D	406	LEU
1	D	449	THR
1	D	460	GLU
1	D	464	ARG
1	D	484	VAL
1	D	488	GLU
1	D	493	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	HIS
1	A	242	HIS
1	A	253	GLN
1	A	307	ASN
1	A	345	ASN
1	A	381	ASN
1	A	422	ASN
1	B	242	HIS
1	B	296	ASN
1	B	307	ASN
1	B	362	ASN
1	B	368	ASN
1	B	381	ASN
1	C	148	GLN
1	C	212	GLN
1	C	242	HIS
1	C	307	ASN
1	C	368	ASN
1	D	183	GLN
1	D	296	ASN
1	D	307	ASN
1	D	345	ASN
1	D	362	ASN
1	D	493	ASN

5.3.3 RNA ⓘ

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 20 ligands modelled in this entry, 8 are monoatomic - leaving 12 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	OXL	A	602	-	5,5,5	1.57	2 (40%)	6,6,6	1.94	4 (66%)
6	PGE	A	606	-	9,9,9	0.20	0	8,8,8	0.26	0
2	OXL	A	601	3	5,5,5	2.04	2 (40%)	6,6,6	1.38	1 (16%)
2	OXL	D	601	3	5,5,5	1.33	1 (20%)	6,6,6	2.57	4 (66%)
2	OXL	C	601	3	5,5,5	1.20	0	6,6,6	1.64	1 (16%)
7	GOL	A	607	-	5,5,5	0.22	0	5,5,5	0.52	0
2	OXL	C	602	-	5,5,5	1.41	1 (20%)	6,6,6	1.90	2 (33%)
2	OXL	D	602	-	5,5,5	1.53	2 (40%)	6,6,6	1.73	2 (33%)
2	OXL	B	601	3	5,5,5	1.84	2 (40%)	6,6,6	1.87	4 (66%)
2	OXL	B	602	-	5,5,5	1.67	2 (40%)	6,6,6	1.62	2 (33%)
6	PGE	C	605	-	9,9,9	0.29	0	8,8,8	0.19	0
5	PG4	A	605	-	12,12,12	0.32	0	11,11,11	0.15	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
2	OXL	A	602	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	PGE	A	606	-	-	6/7/7/7	-
2	OXL	A	601	3	-	4/4/4/4	-
2	OXL	D	601	3	-	4/4/4/4	-
2	OXL	C	601	3	-	4/4/4/4	-
7	GOL	A	607	-	-	4/4/4/4	-
2	OXL	C	602	-	-	4/4/4/4	-
2	OXL	D	602	-	-	4/4/4/4	-
2	OXL	B	601	3	-	4/4/4/4	-
2	OXL	B	602	-	-	4/4/4/4	-
6	PGE	C	605	-	-	2/7/7/7	-
5	PG4	A	605	-	-	5/10/10/10	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	OXL	O4-C2	-3.14	1.22	1.30
2	B	602	OXL	O3-C1	-2.69	1.23	1.30
2	B	601	OXL	O2-C2	2.63	1.29	1.22
2	D	602	OXL	O4-C2	-2.56	1.23	1.30
2	B	601	OXL	O3-C1	-2.54	1.23	1.30
2	B	602	OXL	O4-C2	-2.43	1.24	1.30
2	A	601	OXL	O2-C2	2.40	1.28	1.22
2	D	601	OXL	O3-C1	-2.35	1.24	1.30
2	C	602	OXL	O3-C1	-2.33	1.24	1.30
2	A	602	OXL	O4-C2	-2.29	1.24	1.30
2	D	602	OXL	O3-C1	-2.17	1.24	1.30
2	A	602	OXL	O3-C1	-2.09	1.24	1.30

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	601	OXL	O4-C2-C1	3.54	119.69	112.83
2	D	601	OXL	O3-C1-C2	3.33	119.29	112.83
2	D	601	OXL	O2-C2-C1	-2.98	114.43	120.63
2	C	602	OXL	O2-C2-C1	-2.63	115.15	120.63
2	D	601	OXL	O1-C1-C2	-2.52	115.38	120.63
2	C	602	OXL	O4-C2-C1	2.52	117.71	112.83
2	D	602	OXL	O3-C1-C2	2.37	117.43	112.83
2	A	602	OXL	O1-C1-C2	-2.31	115.81	120.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	602	OXL	O3-C1-C2	2.29	117.27	112.83
2	D	602	OXL	O4-C2-C1	2.23	117.16	112.83
2	B	602	OXL	O3-C1-C2	2.23	117.15	112.83
2	B	601	OXL	O3-C1-C2	2.17	117.04	112.83
2	A	602	OXL	O4-C2-C1	2.17	117.03	112.83
2	A	602	OXL	O2-C2-C1	-2.16	116.13	120.63
2	B	601	OXL	O2-C2-C1	-2.16	116.14	120.63
2	C	601	OXL	O4-C2-C1	2.12	116.95	112.83
2	B	601	OXL	O1-C1-C2	-2.09	116.27	120.63
2	A	601	OXL	O3-C1-C2	2.05	116.81	112.83
2	B	602	OXL	O4-C2-C1	2.05	116.81	112.83
2	B	601	OXL	O4-C2-C1	2.03	116.77	112.83

There are no chirality outliers.

All (49) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	602	OXL	O1-C1-C2-O2
2	B	602	OXL	O1-C1-C2-O4
2	B	602	OXL	O3-C1-C2-O2
2	B	602	OXL	O3-C1-C2-O4
7	A	607	GOL	O1-C1-C2-O2
7	A	607	GOL	O1-C1-C2-C3
7	A	607	GOL	C1-C2-C3-O3
2	A	602	OXL	O1-C1-C2-O2
2	D	602	OXL	O1-C1-C2-O2
2	A	602	OXL	O3-C1-C2-O4
6	A	606	PGE	O2-C3-C4-O3
6	C	605	PGE	C3-C4-O3-C5
6	C	605	PGE	O2-C3-C4-O3
2	A	602	OXL	O1-C1-C2-O4
2	D	602	OXL	O1-C1-C2-O4
2	D	602	OXL	O3-C1-C2-O2
2	D	602	OXL	O3-C1-C2-O4
2	A	602	OXL	O3-C1-C2-O2
6	A	606	PGE	O3-C5-C6-O4
2	C	602	OXL	O3-C1-C2-O4
7	A	607	GOL	O2-C2-C3-O3
5	A	605	PG4	O1-C1-C2-O2
6	A	606	PGE	O1-C1-C2-O2
2	C	602	OXL	O1-C1-C2-O2
2	C	602	OXL	O3-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
5	A	605	PG4	O3-C5-C6-O4
2	C	602	OXL	O1-C1-C2-O4
2	C	601	OXL	O3-C1-C2-O4
2	D	601	OXL	O3-C1-C2-O4
6	A	606	PGE	C6-C5-O3-C4
5	A	605	PG4	C8-C7-O4-C6
2	B	601	OXL	O3-C1-C2-O4
2	C	601	OXL	O1-C1-C2-O2
2	D	601	OXL	O1-C1-C2-O2
2	A	601	OXL	O3-C1-C2-O4
2	A	601	OXL	O1-C1-C2-O2
2	B	601	OXL	O1-C1-C2-O2
6	A	606	PGE	C1-C2-O2-C3
6	A	606	PGE	C3-C4-O3-C5
5	A	605	PG4	O2-C3-C4-O3
2	A	601	OXL	O1-C1-C2-O4
5	A	605	PG4	C1-C2-O2-C3
2	A	601	OXL	O3-C1-C2-O2
2	B	601	OXL	O3-C1-C2-O2
2	C	601	OXL	O1-C1-C2-O4
2	D	601	OXL	O1-C1-C2-O4
2	B	601	OXL	O1-C1-C2-O4
2	D	601	OXL	O3-C1-C2-O2
2	C	601	OXL	O3-C1-C2-O2

There are no ring outliers.

7 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	606	PGE	1	0
2	D	601	OXL	1	0
2	C	601	OXL	1	0
7	A	607	GOL	6	0
2	C	602	OXL	1	0
2	B	601	OXL	1	0
6	C	605	PGE	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	501/521 (96%)	0.28	9 (1%) 67 68	19, 38, 63, 98	1 (0%)
1	B	501/521 (96%)	0.36	22 (4%) 39 38	19, 39, 70, 123	1 (0%)
1	C	501/521 (96%)	0.92	55 (10%) 10 9	29, 51, 75, 111	0
1	D	500/521 (95%)	0.69	29 (5%) 29 27	32, 52, 85, 120	0
All	All	2003/2084 (96%)	0.56	115 (5%) 29 28	19, 45, 74, 123	2 (0%)

All (115) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	281[A]	PHE	6.8
1	B	486	VAL	5.9
1	B	487	GLY	5.2
1	B	372	ARG	4.9
1	C	486	VAL	4.4
1	C	489	ALA	4.3
1	D	485	PRO	4.2
1	C	485	PRO	3.9
1	C	231	VAL	3.8
1	D	486	VAL	3.8
1	C	490	VAL	3.8
1	D	500	VAL	3.8
1	B	373	LEU	3.7
1	C	241	GLY	3.6
1	B	489	ALA	3.5
1	C	74	LEU	3.4
1	D	375	SER	3.3
1	C	235	CYS	3.3
1	C	233	ALA	3.3
1	D	489	ALA	3.3
1	C	234	ILE	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	427	ALA	3.2
1	D	487	GLY	3.2
1	B	370	TYR	3.2
1	D	372	ARG	3.1
1	B	1	MET	3.1
1	C	492	THR	3.0
1	C	455	THR	2.9
1	B	103	ALA	2.9
1	D	451	ALA	2.9
1	C	237	GLU	2.9
1	C	208	PHE	2.8
1	C	137	LEU	2.8
1	C	219	ILE	2.8
1	A	1	MET	2.8
1	D	484	VAL	2.8
1	D	490	VAL	2.8
1	C	166	ALA	2.7
1	D	373	LEU	2.7
1	D	452	PRO	2.7
1	B	378	PHE	2.6
1	D	492	THR	2.6
1	C	217	ILE	2.6
1	B	281[A]	PHE	2.6
1	A	30	GLY	2.6
1	C	487	GLY	2.6
1	B	491	ARG	2.6
1	D	17	ILE	2.6
1	C	430	PHE	2.5
1	B	282	GLU	2.5
1	C	406	LEU	2.5
1	A	451	ALA	2.5
1	C	243	VAL	2.5
1	C	458	MET	2.5
1	D	458	MET	2.5
1	C	434	THR	2.5
1	D	455	THR	2.5
1	C	1	MET	2.5
1	C	452	PRO	2.5
1	C	210	LEU	2.5
1	C	433	LEU	2.5
1	C	214	ILE	2.5
1	D	128	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	370	TYR	2.4
1	C	245	LEU	2.4
1	C	484	VAL	2.4
1	B	178	GLY	2.4
1	D	106	TYR	2.4
1	C	459	PHE	2.4
1	C	87	LEU	2.4
1	D	459	PHE	2.4
1	A	497	ILE	2.4
1	A	501	ARG	2.4
1	A	454	SER	2.4
1	D	496	ARG	2.3
1	A	387	MET	2.3
1	C	246	PHE	2.3
1	C	196	PRO	2.3
1	C	209	GLY	2.3
1	D	1	MET	2.3
1	C	199	ALA	2.3
1	D	35	VAL	2.3
1	B	132	ASN	2.3
1	C	198	LEU	2.3
1	B	161	VAL	2.2
1	C	197	ALA	2.2
1	C	215	ASN	2.2
1	B	490	VAL	2.2
1	C	39	ALA	2.2
1	B	179	ILE	2.2
1	C	412	THR	2.2
1	C	438	LEU	2.2
1	D	74	LEU	2.2
1	C	262	ILE	2.1
1	D	384	THR	2.1
1	D	30	GLY	2.1
1	B	117	ALA	2.1
1	C	451	ALA	2.1
1	C	238	THR	2.1
1	C	228	VAL	2.1
1	D	488	GLU	2.1
1	B	114	ILE	2.1
1	C	222	VAL	2.1
1	D	33	LEU	2.1
1	C	70	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	264	ALA	2.1
1	B	375	SER	2.0
1	C	179	ILE	2.0
1	C	206	ILE	2.0
1	A	27	GLY	2.0
1	B	368	ASN	2.0
1	D	231	VAL	2.0
1	C	265	ALA	2.0
1	B	458	MET	2.0
1	C	221	PHE	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	OXL	C	602	6/6	0.75	0.14	60,67,69,74	0
6	PGE	C	605	10/10	0.77	0.16	65,75,83,83	0
5	PG4	A	605	13/13	0.83	0.15	56,69,79,79	0
2	OXL	D	602	6/6	0.83	0.12	58,72,83,85	0
7	GOL	A	607	6/6	0.87	0.15	42,54,58,61	0
2	OXL	B	602	6/6	0.88	0.11	44,56,79,82	0
4	K	C	604	1/1	0.89	0.13	50,50,50,50	0
2	OXL	A	602	6/6	0.90	0.09	42,49,52,55	0
6	PGE	A	606	10/10	0.91	0.11	52,55,60,66	0
2	OXL	D	601	6/6	0.93	0.07	37,45,48,52	0
2	OXL	C	601	6/6	0.94	0.06	34,41,44,46	0
2	OXL	A	601	6/6	0.95	0.07	32,36,43,44	0
2	OXL	B	601	6/6	0.97	0.05	28,35,39,39	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	K	D	604	1/1	0.97	0.10	48,48,48,48	0
3	MG	C	603	1/1	0.97	0.08	37,37,37,37	0
3	MG	A	603	1/1	0.99	0.05	32,32,32,32	0
4	K	A	604	1/1	0.99	0.06	37,37,37,37	0
4	K	B	604	1/1	0.99	0.21	48,48,48,48	0
3	MG	B	603	1/1	1.00	0.04	30,30,30,30	0
3	MG	D	603	1/1	1.00	0.02	41,41,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.