



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:41 AM UTC

PDB ID : 6WP4 / pdb_00006wp4
Title : Pyruvate Kinase M2 mutant-S37E
Authors : Nandi, S.; Razzaghi, M.; Srivastava, D.; Dey, M.
Deposited on : 2020-04-26
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

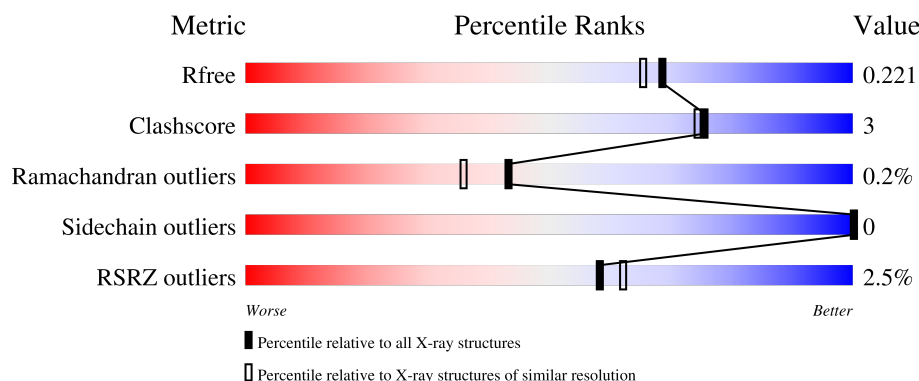
1 Overall quality at a glance

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	550	0% 85% 7% 7%
1	B	550	2% 86% 7% 7%
1	C	550	2% 85% 7% 8%
1	D	550	4% 85% 7% 8%

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
5	OXL	A	606	-	X	-	-
5	OXL	D	607	-	X	-	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 15950 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 PKM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	511	Total	C	N	O	S	0	0	0
			3849	2424	679	721	25			
1	B	511	Total	C	N	O	S	0	0	0
			3828	2420	671	712	25			
1	C	504	Total	C	N	O	S	0	0	0
			3773	2384	664	700	25			
1	D	506	Total	C	N	O	S	0	0	0
			3751	2364	656	706	25			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MET	-	initiating methionine	UNP P14618
A	-17	GLY	-	expression tag	UNP P14618
A	-16	SER	-	expression tag	UNP P14618
A	-15	SER	-	expression tag	UNP P14618
A	-14	HIS	-	expression tag	UNP P14618
A	-13	HIS	-	expression tag	UNP P14618
A	-12	HIS	-	expression tag	UNP P14618
A	-11	HIS	-	expression tag	UNP P14618
A	-10	HIS	-	expression tag	UNP P14618
A	-9	HIS	-	expression tag	UNP P14618
A	-8	SER	-	expression tag	UNP P14618
A	-7	SER	-	expression tag	UNP P14618
A	-6	GLY	-	expression tag	UNP P14618
A	-5	LEU	-	expression tag	UNP P14618
A	-4	VAL	-	expression tag	UNP P14618
A	-3	PRO	-	expression tag	UNP P14618
A	-2	ARG	-	expression tag	UNP P14618
A	-1	GLY	-	expression tag	UNP P14618
A	0	SER	-	expression tag	UNP P14618
A	37	GLU	SER	engineered mutation	UNP P14618
B	-18	MET	-	initiating methionine	UNP P14618

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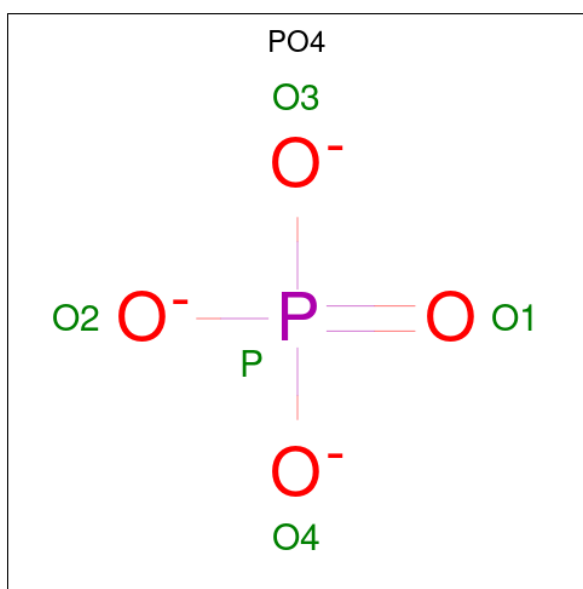
Chain	Residue	Modelled	Actual	Comment	Reference
B	-17	GLY	-	expression tag	UNP P14618
B	-16	SER	-	expression tag	UNP P14618
B	-15	SER	-	expression tag	UNP P14618
B	-14	HIS	-	expression tag	UNP P14618
B	-13	HIS	-	expression tag	UNP P14618
B	-12	HIS	-	expression tag	UNP P14618
B	-11	HIS	-	expression tag	UNP P14618
B	-10	HIS	-	expression tag	UNP P14618
B	-9	HIS	-	expression tag	UNP P14618
B	-8	SER	-	expression tag	UNP P14618
B	-7	SER	-	expression tag	UNP P14618
B	-6	GLY	-	expression tag	UNP P14618
B	-5	LEU	-	expression tag	UNP P14618
B	-4	VAL	-	expression tag	UNP P14618
B	-3	PRO	-	expression tag	UNP P14618
B	-2	ARG	-	expression tag	UNP P14618
B	-1	GLY	-	expression tag	UNP P14618
B	0	SER	-	expression tag	UNP P14618
B	37	GLU	SER	engineered mutation	UNP P14618
C	-18	MET	-	initiating methionine	UNP P14618
C	-17	GLY	-	expression tag	UNP P14618
C	-16	SER	-	expression tag	UNP P14618
C	-15	SER	-	expression tag	UNP P14618
C	-14	HIS	-	expression tag	UNP P14618
C	-13	HIS	-	expression tag	UNP P14618
C	-12	HIS	-	expression tag	UNP P14618
C	-11	HIS	-	expression tag	UNP P14618
C	-10	HIS	-	expression tag	UNP P14618
C	-9	HIS	-	expression tag	UNP P14618
C	-8	SER	-	expression tag	UNP P14618
C	-7	SER	-	expression tag	UNP P14618
C	-6	GLY	-	expression tag	UNP P14618
C	-5	LEU	-	expression tag	UNP P14618
C	-4	VAL	-	expression tag	UNP P14618
C	-3	PRO	-	expression tag	UNP P14618
C	-2	ARG	-	expression tag	UNP P14618
C	-1	GLY	-	expression tag	UNP P14618
C	0	SER	-	expression tag	UNP P14618
C	37	GLU	SER	engineered mutation	UNP P14618
D	-18	MET	-	initiating methionine	UNP P14618
D	-17	GLY	-	expression tag	UNP P14618
D	-16	SER	-	expression tag	UNP P14618

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-15	SER	-	expression tag	UNP P14618
D	-14	HIS	-	expression tag	UNP P14618
D	-13	HIS	-	expression tag	UNP P14618
D	-12	HIS	-	expression tag	UNP P14618
D	-11	HIS	-	expression tag	UNP P14618
D	-10	HIS	-	expression tag	UNP P14618
D	-9	HIS	-	expression tag	UNP P14618
D	-8	SER	-	expression tag	UNP P14618
D	-7	SER	-	expression tag	UNP P14618
D	-6	GLY	-	expression tag	UNP P14618
D	-5	LEU	-	expression tag	UNP P14618
D	-4	VAL	-	expression tag	UNP P14618
D	-3	PRO	-	expression tag	UNP P14618
D	-2	ARG	-	expression tag	UNP P14618
D	-1	GLY	-	expression tag	UNP P14618
D	0	SER	-	expression tag	UNP P14618
D	37	GLU	SER	engineered mutation	UNP P14618

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		

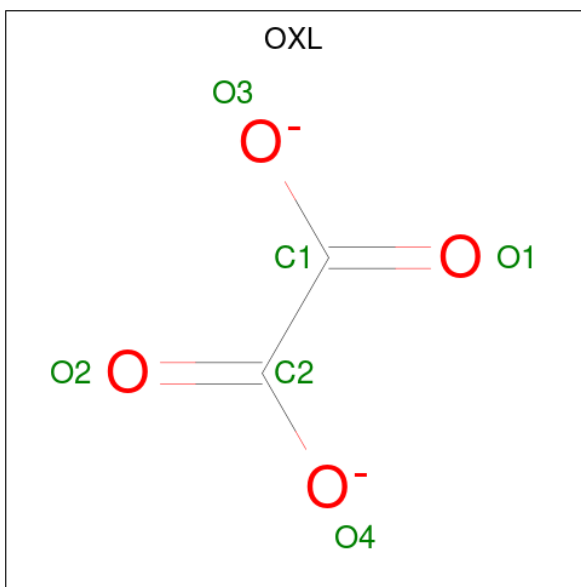
- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

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

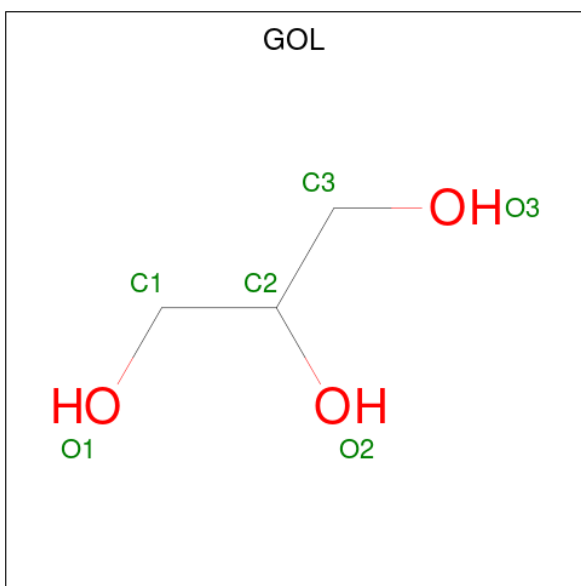
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	K	0	0
			1	1		
4	B	1	Total	K	0	0
			1	1		
4	C	1	Total	K	0	0
			1	1		
4	D	2	Total	K	0	0
			2	2		

- Molecule 5 is OXALATE ION (CCD ID: OXL) (formula: C₂O₄).



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

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

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

- Molecule 7 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	C	2	Total Cl 2 2	0	0

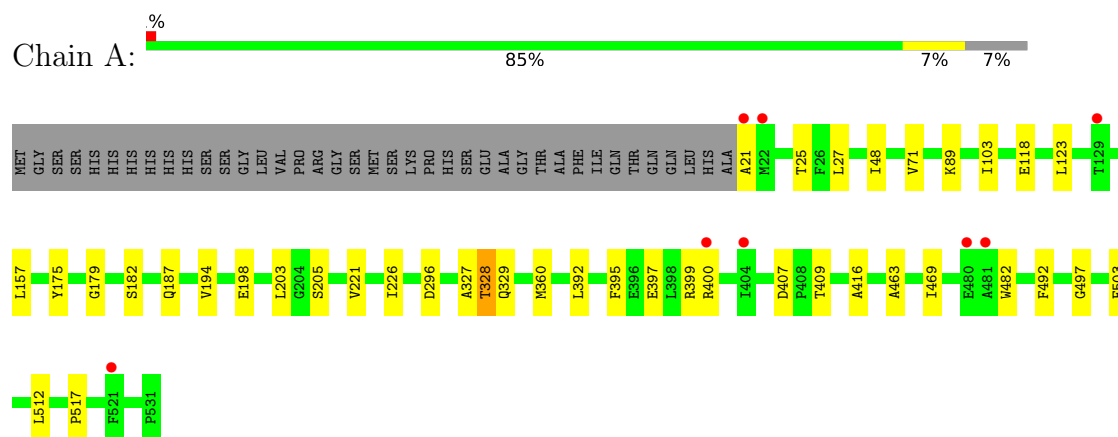
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	171	Total 171	O 171	0	0
8	B	149	Total 149	O 149	0	0
8	C	143	Total 143	O 143	0	0
8	D	100	Total 100	O 100	0	0

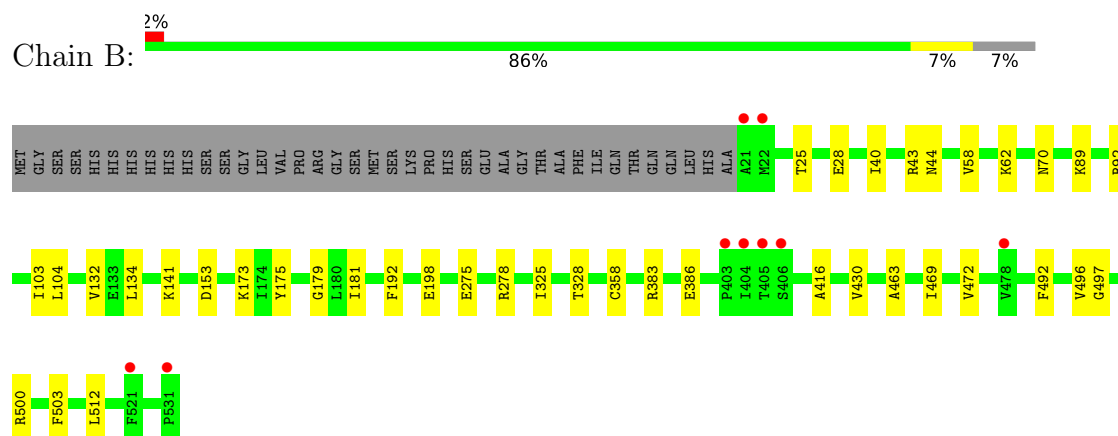
3 Residue-property plots [i](#)

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.

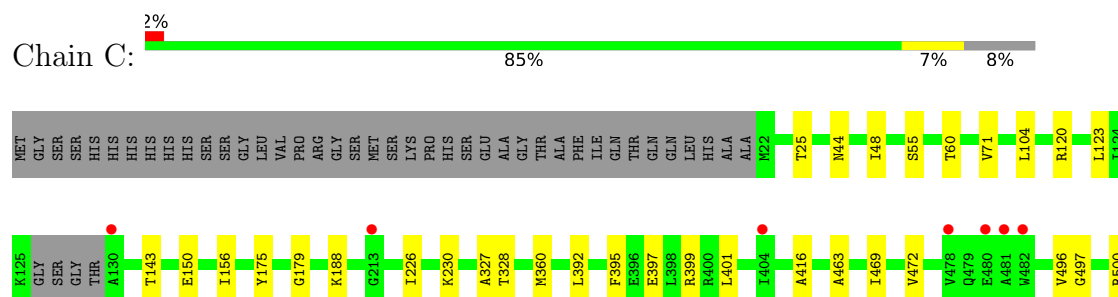
- Molecule 1: Pyruvate kinase PKM

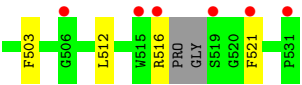


- Molecule 1: Pyruvate kinase PKM

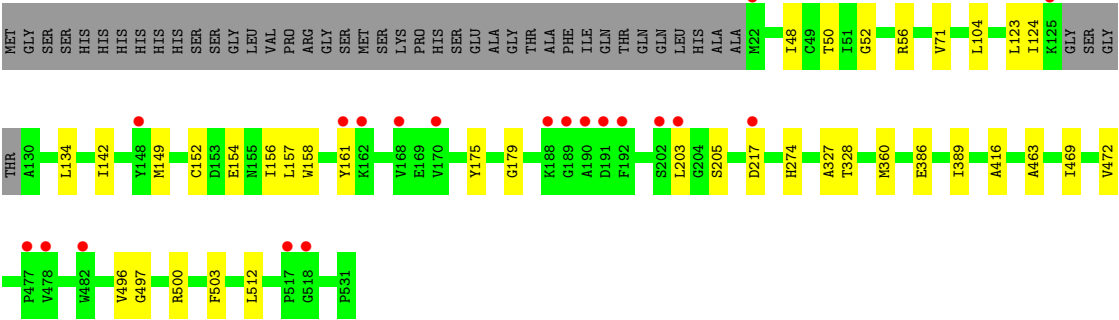
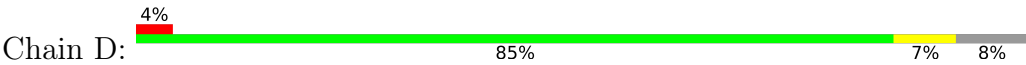


- Molecule 1: Pyruvate kinase PKM





● Molecule 1: Pyruvate kinase PKM



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	94.17Å 132.63Å 110.13Å 90.00° 112.74° 90.00°	Depositor
Resolution (Å)	55.53 – 1.90 55.53 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.9 (55.53-1.90) 97.9 (55.53-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 1.90Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.189 , 0.221 0.191 , 0.221	Depositor DCC
R_{free} test set	9566 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	24.9	Xtriage
Anisotropy	0.633	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 40.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.037 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15950	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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: MG, GOL, CL, OXL, PO4, 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	0.32	0/3909	0.49	0/5291
1	B	0.30	0/3891	0.47	0/5270
1	C	0.29	0/3831	0.47	0/5183
1	D	0.28	0/3811	0.47	0/5170
All	All	0.30	0/15442	0.47	0/20914

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	3849	0	3855	28	0
1	B	3828	0	3838	23	0
1	C	3773	0	3777	23	0
1	D	3751	0	3694	21	0
2	A	15	0	0	0	0
2	B	15	0	0	1	0
2	C	10	0	0	0	0
2	D	15	0	0	0	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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	2	0	0	0	0
5	A	6	0	0	0	0
5	D	6	0	0	0	0
6	A	36	0	48	5	0
6	B	18	0	24	3	0
6	C	30	0	40	3	0
6	D	24	0	32	2	0
7	C	2	0	0	0	0
8	A	171	0	0	3	0
8	B	149	0	0	1	0
8	C	143	0	0	0	0
8	D	100	0	0	0	0
All	All	15950	0	15308	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 3.

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:C:120:ARG:HH12	6:C:606:GOL:H2	1.34	0.92
6:B:606:GOL:O3	8:B:701:HOH:O	1.99	0.79
1:A:400:ARG:NE	8:A:701:HOH:O	2.26	0.69
6:A:609:GOL:H31	1:B:89:LYS:HA	1.75	0.68
1:C:416:ALA:HB2	1:C:512:LEU:HD21	1.83	0.61
1:A:407:ASP:OD1	1:A:409:THR:OG1	2.17	0.60
1:A:21:ALA:HB1	1:A:392:LEU:H	1.69	0.58
1:C:123:LEU:HD12	1:C:150:GLU:HG2	1.86	0.57
1:D:123:LEU:HA	1:D:205:SER:HB3	1.86	0.57
1:C:226:ILE:O	1:C:230:LYS:HG2	2.05	0.57
1:B:43:ARG:O	6:B:606:GOL:H32	2.05	0.56
1:C:175:TYR:HB3	1:C:179:GLY:HA2	1.87	0.56
1:A:27:LEU:HD23	1:C:401:LEU:HD12	1.88	0.55
1:C:104:LEU:O	1:C:500:ARG:NH1	2.39	0.55
1:B:430:VAL:HG22	1:B:512:LEU:HD12	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:LEU:HA	1:A:205:SER:HB3	1.89	0.53
1:D:175:TYR:HB3	1:D:179:GLY:HA2	1.90	0.53
1:B:134:LEU:HD12	1:B:181:ILE:HG21	1.91	0.52
1:A:400:ARG:NH1	8:A:701:HOH:O	2.43	0.52
1:C:395:PHE:CZ	1:C:399:ARG:HD2	2.44	0.52
1:A:395:PHE:CZ	1:A:399:ARG:HD3	2.46	0.51
1:D:154:GLU:CD	1:D:154:GLU:H	2.17	0.51
1:D:416:ALA:HB2	1:D:512:LEU:HD21	1.92	0.51
1:D:327:ALA:HB1	1:D:360:MET:HE2	1.93	0.51
1:C:123:LEU:HB2	1:C:150:GLU:HA	1.92	0.51
1:B:472:VAL:HG11	1:B:496:VAL:HG11	1.93	0.50
1:C:472:VAL:HG11	1:C:496:VAL:HG11	1.93	0.50
1:A:400:ARG:NH2	8:A:701:HOH:O	2.44	0.49
6:A:609:GOL:H32	1:B:92:ARG:HD2	1.93	0.49
1:D:274:HIS:HB3	6:D:611:GOL:H31	1.94	0.49
1:B:104:LEU:O	1:B:500:ARG:NH2	2.45	0.49
1:B:416:ALA:HB2	1:B:512:LEU:HD21	1.95	0.49
1:D:124:ILE:HA	1:D:152:CYS:HB2	1.94	0.48
1:A:182:SER:OG	1:A:198:GLU:HB3	2.14	0.48
1:A:416:ALA:CB	1:A:512:LEU:HD11	2.43	0.48
1:A:89:LYS:HA	6:A:608:GOL:H11	1.96	0.48
1:B:173:LYS:HE2	1:B:198:GLU:OE1	2.14	0.48
1:C:497:GLY:HA3	1:C:503:PHE:CZ	2.49	0.48
1:A:157:LEU:HD13	1:A:203:LEU:HD21	1.96	0.47
1:C:463:ALA:HB1	1:C:469:ILE:HG21	1.97	0.47
1:B:44:ASN:H	1:B:386:GLU:CD	2.23	0.47
1:C:516:ARG:HD2	1:C:521:PHE:HB3	1.97	0.46
1:D:142:ILE:HD13	1:D:157:LEU:HB3	1.97	0.46
1:A:118:GLU:HA	6:A:612:GOL:H32	1.96	0.46
1:B:40:ILE:O	1:B:383:ARG:HD2	2.16	0.46
1:C:48:ILE:HG12	1:C:71:VAL:HB	1.96	0.46
1:D:124:ILE:HD13	1:D:152:CYS:HB2	1.97	0.46
1:D:149:MET:HG3	1:D:158:TRP:CH2	2.51	0.46
1:A:327:ALA:HB1	1:A:360:MET:HE2	1.97	0.46
1:A:463:ALA:HB1	1:A:469:ILE:HG21	1.97	0.46
1:B:175:TYR:HB3	1:B:179:GLY:HA2	1.98	0.45
1:B:325:ILE:HG12	1:B:358:CYS:HB2	1.98	0.45
1:C:392:LEU:HA	6:C:609:GOL:H11	1.99	0.45
1:B:278:ARG:NH2	2:B:603:PO4:O3	2.48	0.44
1:B:275:GLU:HG3	1:B:278:ARG:NH2	2.32	0.44
1:D:156:ILE:HD12	1:D:156:ILE:HA	1.90	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:463:ALA:HB1	1:D:469:ILE:HG21	2.00	0.44
1:B:463:ALA:HB1	1:B:469:ILE:HG21	1.99	0.43
1:A:221:VAL:HG12	1:A:226:ILE:HG13	2.00	0.43
1:D:48:ILE:HG12	1:D:71:VAL:HB	2.01	0.43
1:A:296:ASP:OD1	6:A:612:GOL:H12	2.19	0.43
1:B:58:VAL:HG12	1:B:62:LYS:HE3	2.00	0.43
1:D:50:THR:HG21	6:D:609:GOL:H2	2.01	0.43
1:D:104:LEU:O	1:D:500:ARG:NH1	2.52	0.43
1:A:25:THR:HB	1:C:397:GLU:CD	2.43	0.43
1:A:103:ILE:HD13	1:A:492:PHE:CE1	2.54	0.43
1:A:497:GLY:HA3	1:A:503:PHE:CZ	2.54	0.43
1:C:44:ASN:O	6:C:605:GOL:H31	2.19	0.42
1:A:21:ALA:CB	1:A:392:LEU:H	2.31	0.42
1:D:386:GLU:HA	1:D:389:ILE:HD12	2.01	0.42
1:B:70:ASN:CG	6:B:606:GOL:H12	2.44	0.42
1:C:188:LYS:HB3	1:C:188:LYS:HE3	1.86	0.42
1:B:132:VAL:HG11	1:B:153:ASP:HA	2.02	0.42
1:B:141:LYS:HE2	1:B:192:PHE:CD2	2.55	0.42
1:D:52:GLY:O	1:D:56:ARG:HB2	2.19	0.42
1:D:472:VAL:HG11	1:D:496:VAL:HG11	2.01	0.42
1:D:497:GLY:HA3	1:D:503:PHE:CZ	2.54	0.42
1:C:143:THR:HB	1:C:156:ILE:HD11	2.01	0.42
1:D:134:LEU:HD11	1:D:203:LEU:HD22	2.01	0.42
1:B:497:GLY:HA3	1:B:503:PHE:CZ	2.55	0.42
1:D:161:TYR:CE1	1:D:217:ASP:OD2	2.73	0.42
1:A:175:TYR:HB3	1:A:179:GLY:HA2	2.02	0.42
1:A:400:ARG:NE	1:A:400:ARG:HB3	2.35	0.42
1:A:328:THR:HG22	1:A:329:GLN:HG3	2.00	0.41
1:A:48:ILE:HG12	1:A:71:VAL:HB	2.03	0.41
1:A:397:GLU:CD	1:C:25:THR:HB	2.46	0.41
1:B:103:ILE:HD13	1:B:492:PHE:CE1	2.56	0.41
1:A:482:TRP:CE3	1:A:517:PRO:HG3	2.56	0.41
1:B:25:THR:OG1	1:B:28:GLU:HG3	2.21	0.41
1:C:327:ALA:HB1	1:C:360:MET:HE2	2.03	0.41
1:A:187:GLN:HB3	1:A:194:VAL:HB	2.03	0.40
1:C:516:ARG:HH11	1:C:521:PHE:HB3	1.86	0.40
1:C:55:SER:HA	1:C:60:THR:HG21	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	509/550 (92%)	501 (98%)	7 (1%)	1 (0%)	43	36
1	B	509/550 (92%)	498 (98%)	10 (2%)	1 (0%)	43	36
1	C	498/550 (90%)	488 (98%)	9 (2%)	1 (0%)	43	36
1	D	502/550 (91%)	493 (98%)	8 (2%)	1 (0%)	43	36
All	All	2018/2200 (92%)	1980 (98%)	34 (2%)	4 (0%)	43	36

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	328	THR
1	B	328	THR
1	C	328	THR
1	D	328	THR

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	400/452 (88%)	400 (100%)	0	100	100
1	B	395/452 (87%)	395 (100%)	0	100	100
1	C	387/452 (86%)	387 (100%)	0	100	100
1	D	381/452 (84%)	381 (100%)	0	100	100
All	All	1563/1808 (86%)	1563 (100%)	0	100	100

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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	199	ASN
1	A	210	ASN
1	A	440	GLN
1	B	393	GLN
1	B	440	GLN
1	C	210	ASN
1	C	458	GLN
1	D	393	GLN
1	D	458	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 42 ligands modelled in this entry, 11 are monoatomic - leaving 31 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	$\# Z > 2$
6	GOL	D	611	-	5,5,5	1.15	0	5,5,5	0.91	0
6	GOL	D	608	-	5,5,5	1.22	1 (20%)	5,5,5	0.93	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PO4	A	601	-	4,4,4	0.96	0	6,6,6	0.90	0
2	PO4	B	602	-	4,4,4	1.03	0	6,6,6	0.83	0
2	PO4	A	603	-	4,4,4	0.88	0	6,6,6	0.51	0
6	GOL	B	606	-	5,5,5	0.84	0	5,5,5	1.58	1 (20%)
6	GOL	A	612	-	5,5,5	0.85	0	5,5,5	1.09	1 (20%)
6	GOL	C	608	-	5,5,5	1.02	0	5,5,5	1.00	0
5	OXL	A	606	3	5,5,5	1.35	0	6,6,6	2.70	2 (33%)
6	GOL	C	607	-	5,5,5	0.79	0	5,5,5	1.20	0
2	PO4	D	602	-	4,4,4	0.94	0	6,6,6	0.50	0
6	GOL	A	607	-	5,5,5	0.90	0	5,5,5	1.18	1 (20%)
6	GOL	C	605	-	5,5,5	1.62	2 (40%)	5,5,5	1.30	0
6	GOL	C	609	-	5,5,5	1.08	0	5,5,5	1.04	0
2	PO4	C	601	-	4,4,4	0.98	0	6,6,6	0.58	0
6	GOL	A	608	-	5,5,5	0.84	0	5,5,5	1.15	1 (20%)
2	PO4	B	601	-	4,4,4	0.84	0	6,6,6	0.51	0
2	PO4	C	602	-	4,4,4	0.96	0	6,6,6	0.62	0
6	GOL	A	609	-	5,5,5	0.84	0	5,5,5	0.88	0
6	GOL	B	608	-	5,5,5	0.80	0	5,5,5	1.25	1 (20%)
6	GOL	D	609	-	5,5,5	0.91	0	5,5,5	1.03	0
2	PO4	D	601	-	4,4,4	1.15	0	6,6,6	0.57	0
6	GOL	B	607	-	5,5,5	0.68	0	5,5,5	1.19	0
2	PO4	D	603	-	4,4,4	0.93	0	6,6,6	0.47	0
2	PO4	A	602	-	4,4,4	0.87	0	6,6,6	0.67	0
5	OXL	D	607	3	5,5,5	1.54	0	6,6,6	1.97	2 (33%)
6	GOL	A	611	-	5,5,5	0.49	0	5,5,5	1.13	0
2	PO4	B	603	-	4,4,4	0.85	0	6,6,6	0.64	0
6	GOL	C	606	-	5,5,5	0.87	0	5,5,5	1.01	0
6	GOL	D	610	-	5,5,5	0.85	0	5,5,5	1.14	0
6	GOL	A	610	-	5,5,5	0.65	0	5,5,5	1.54	1 (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
6	GOL	D	611	-	-	2/4/4/4	-
6	GOL	D	608	-	-	2/4/4/4	-
6	GOL	B	606	-	-	2/4/4/4	-
6	GOL	A	612	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	GOL	C	608	-	-	2/4/4/4	-
5	OXL	A	606	3	-	4/4/4/4	-
6	GOL	C	607	-	-	2/4/4/4	-
6	GOL	A	607	-	-	2/4/4/4	-
6	GOL	C	605	-	-	2/4/4/4	-
6	GOL	C	609	-	-	1/4/4/4	-
6	GOL	A	608	-	-	2/4/4/4	-
6	GOL	A	609	-	-	4/4/4/4	-
6	GOL	B	608	-	-	4/4/4/4	-
6	GOL	D	609	-	-	4/4/4/4	-
6	GOL	B	607	-	-	0/4/4/4	-
5	OXL	D	607	3	-	4/4/4/4	-
6	GOL	A	611	-	-	0/4/4/4	-
6	GOL	C	606	-	-	0/4/4/4	-
6	GOL	D	610	-	-	2/4/4/4	-
6	GOL	A	610	-	-	0/4/4/4	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	605	GOL	C3-C2	2.53	1.61	1.51
6	C	605	GOL	O2-C2	-2.17	1.37	1.43
6	D	608	GOL	O2-C2	-2.06	1.37	1.43

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	606	OXL	O3-C1-C2	4.45	121.45	112.83
5	A	606	OXL	O4-C2-C1	4.16	120.91	112.83
5	D	607	OXL	O3-C1-C2	3.31	119.25	112.83
6	B	606	GOL	C3-C2-C1	-3.20	100.04	111.80
5	D	607	OXL	O4-C2-C1	3.08	118.80	112.83
6	A	610	GOL	C3-C2-C1	-2.90	101.17	111.80
6	A	612	GOL	C3-C2-C1	-2.10	104.09	111.80
6	A	607	GOL	C3-C2-C1	-2.08	104.17	111.80
6	A	608	GOL	C3-C2-C1	-2.07	104.20	111.80
6	B	608	GOL	C3-C2-C1	-2.06	104.25	111.80

There are no chirality outliers.

All (43) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	609	GOL	C1-C2-C3-O3
6	A	612	GOL	O1-C1-C2-C3
6	A	612	GOL	C1-C2-C3-O3
6	B	608	GOL	C1-C2-C3-O3
6	B	608	GOL	O2-C2-C3-O3
6	D	609	GOL	O1-C1-C2-C3
6	D	609	GOL	C1-C2-C3-O3
6	D	610	GOL	C1-C2-C3-O3
6	A	609	GOL	O2-C2-C3-O3
6	B	606	GOL	O1-C1-C2-O2
6	A	607	GOL	O1-C1-C2-C3
6	A	608	GOL	O1-C1-C2-C3
6	A	609	GOL	O1-C1-C2-C3
6	B	606	GOL	O1-C1-C2-C3
6	B	608	GOL	O1-C1-C2-C3
6	C	605	GOL	O1-C1-C2-C3
6	C	608	GOL	O1-C1-C2-C3
6	D	608	GOL	C1-C2-C3-O3
6	D	611	GOL	C1-C2-C3-O3
6	A	607	GOL	O1-C1-C2-O2
6	B	608	GOL	O1-C1-C2-O2
6	C	605	GOL	O1-C1-C2-O2
6	D	608	GOL	O2-C2-C3-O3
6	D	609	GOL	O2-C2-C3-O3
6	D	611	GOL	O2-C2-C3-O3
6	A	612	GOL	O1-C1-C2-O2
6	D	610	GOL	O2-C2-C3-O3
6	A	608	GOL	O1-C1-C2-O2
6	A	612	GOL	O2-C2-C3-O3
6	C	607	GOL	O1-C1-C2-O2
6	C	608	GOL	O1-C1-C2-O2
6	C	609	GOL	O2-C2-C3-O3
6	D	609	GOL	O1-C1-C2-O2
5	D	607	OXL	O3-C1-C2-O4
6	C	607	GOL	O1-C1-C2-C3
5	D	607	OXL	O1-C1-C2-O2
5	A	606	OXL	O3-C1-C2-O4
5	A	606	OXL	O1-C1-C2-O2
6	A	609	GOL	O1-C1-C2-O2
5	D	607	OXL	O1-C1-C2-O4
5	D	607	OXL	O3-C1-C2-O2
5	A	606	OXL	O1-C1-C2-O4

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Mol	Chain	Res	Type	Atoms
5	A	606	OXL	O3-C1-C2-O2

There are no ring outliers.

10 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	611	GOL	1	0
6	B	606	GOL	3	0
6	A	612	GOL	2	0
6	C	605	GOL	1	0
6	C	609	GOL	1	0
6	A	608	GOL	1	0
6	A	609	GOL	2	0
6	D	609	GOL	1	0
2	B	603	PO4	1	0
6	C	606	GOL	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	511/550 (92%)	0.00	8 (1%) 70 74	16, 25, 40, 52	0
1	B	511/550 (92%)	0.09	9 (1%) 67 71	17, 26, 42, 56	0
1	C	504/550 (91%)	0.13	13 (2%) 57 61	16, 28, 43, 55	0
1	D	506/550 (92%)	0.29	20 (3%) 42 45	18, 30, 48, 63	0
All	All	2032/2200 (92%)	0.13	50 (2%) 58 62	16, 27, 44, 63	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	515	TRP	5.0
1	D	192	PHE	4.6
1	C	516	ARG	4.2
1	A	21	ALA	3.9
1	C	521	PHE	3.8
1	A	481	ALA	3.8
1	D	190	ALA	3.8
1	D	148	TYR	3.6
1	D	189	GLY	3.5
1	B	406	SER	3.5
1	B	531	PRO	3.4
1	D	478	VAL	3.4
1	C	482	TRP	3.4
1	D	482	TRP	3.3
1	C	506	GLY	3.1
1	C	478	VAL	3.1
1	B	21	ALA	3.1
1	D	191	ASP	3.0
1	B	404	ILE	2.9
1	D	217	ASP	2.9
1	D	517	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	162	LYS	2.7
1	C	481	ALA	2.6
1	D	168	VAL	2.6
1	B	403	PRO	2.5
1	C	213	GLY	2.5
1	B	22	MET	2.5
1	D	203	LEU	2.4
1	A	480	GLU	2.4
1	A	129	THR	2.4
1	C	130	ALA	2.4
1	D	170	VAL	2.4
1	B	405	THR	2.4
1	C	519	SER	2.3
1	C	404	ILE	2.3
1	C	480	GLU	2.3
1	B	478	VAL	2.3
1	D	125	LYS	2.3
1	A	404	ILE	2.2
1	A	521	PHE	2.2
1	D	202	SER	2.2
1	D	188	LYS	2.2
1	B	521	PHE	2.1
1	A	22	MET	2.1
1	D	161	TYR	2.1
1	D	518	GLY	2.1
1	C	531	PRO	2.1
1	D	477	PRO	2.1
1	D	22	MET	2.1
1	A	400	ARG	2.1

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(Å ²)	Q<0.9
6	GOL	D	611	6/6	0.72	0.16	41,47,50,53	0
6	GOL	B	606	6/6	0.81	0.16	24,28,32,38	0
6	GOL	C	608	6/6	0.84	0.14	32,37,40,40	0
2	PO4	B	603	5/5	0.84	0.12	46,47,53,54	0
6	GOL	D	608	6/6	0.85	0.18	30,32,36,36	0
4	K	D	606	1/1	0.85	0.13	73,73,73,73	0
6	GOL	C	606	6/6	0.86	0.13	32,39,41,45	0
6	GOL	A	609	6/6	0.87	0.11	33,35,38,48	0
6	GOL	D	609	6/6	0.87	0.13	36,39,44,45	0
6	GOL	A	608	6/6	0.87	0.12	31,37,42,49	0
2	PO4	D	603	5/5	0.88	0.08	45,49,58,60	0
6	GOL	A	610	6/6	0.88	0.13	24,33,36,36	0
6	GOL	A	611	6/6	0.89	0.14	27,33,35,36	0
6	GOL	A	612	6/6	0.89	0.15	24,28,34,40	0
7	CL	C	610	1/1	0.89	0.18	65,65,65,65	0
6	GOL	A	607	6/6	0.90	0.10	32,35,38,48	0
6	GOL	C	605	6/6	0.90	0.13	14,29,31,41	0
2	PO4	A	603	5/5	0.91	0.19	39,41,46,51	0
6	GOL	C	609	6/6	0.92	0.12	32,36,39,43	0
6	GOL	B	608	6/6	0.92	0.10	34,39,44,46	0
3	MG	C	603	1/1	0.92	0.14	35,35,35,35	0
2	PO4	D	602	5/5	0.92	0.08	49,50,55,58	0
2	PO4	C	602	5/5	0.92	0.08	38,45,48,65	0
6	GOL	D	610	6/6	0.93	0.09	31,34,42,43	0
6	GOL	B	607	6/6	0.93	0.09	28,35,36,38	0
5	OXL	A	606	6/6	0.93	0.08	25,26,31,32	0
7	CL	C	611	1/1	0.93	0.12	54,54,54,54	0
3	MG	B	604	1/1	0.94	0.04	27,27,27,27	0
6	GOL	C	607	6/6	0.94	0.09	30,37,38,42	0
5	OXL	D	607	6/6	0.95	0.08	25,26,33,36	0
2	PO4	A	602	5/5	0.96	0.13	29,31,36,36	0
2	PO4	B	601	5/5	0.97	0.09	28,28,35,37	0
2	PO4	A	601	5/5	0.97	0.07	26,28,29,30	0
3	MG	D	604	1/1	0.98	0.04	31,31,31,31	0
2	PO4	B	602	5/5	0.98	0.05	25,27,29,29	0
2	PO4	D	601	5/5	0.98	0.06	26,28,33,34	0
2	PO4	C	601	5/5	0.98	0.05	27,29,31,33	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	K	C	604	1/1	0.99	0.03	28,28,28,28	0
4	K	D	605	1/1	0.99	0.06	28,28,28,28	0
3	MG	A	604	1/1	0.99	0.02	24,24,24,24	0
4	K	A	605	1/1	0.99	0.03	21,21,21,21	0
4	K	B	605	1/1	0.99	0.04	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.