



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 30, 2026 – 11:50 AM UTC

PDB ID : 5WS8 / pdb_00005ws8
Title : Pyruvate kinase (PYK) from Mycobacterium tuberculosis in complex with Oxalate
Authors : Zhong, W.; Cai, Q.; El Sahili, A.; Lescar, J.; Dedon, P.C.
Deposited on : 2016-12-06
Resolution : 2.62 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

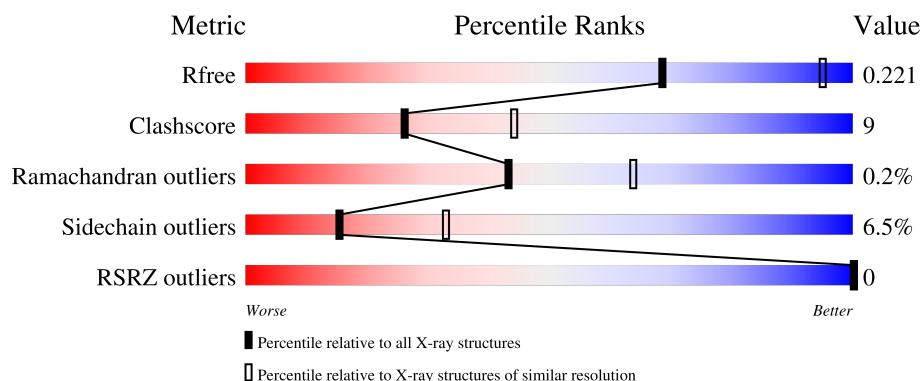
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.62 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	4951 (2.64-2.60)
Clashscore	190562	5303 (2.64-2.60)
Ramachandran outliers	187476	5217 (2.64-2.60)
Sidechain outliers	187428	5217 (2.64-2.60)
RSRZ outliers	180081	4950 (2.64-2.60)

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	475	
1	B	475	
1	C	475	
1	D	475	

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
3	OXL	A	502	-	X	-	-
3	OXL	B	502	-	X	-	-
3	OXL	C	502	-	X	-	-
3	OXL	D	502	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14566 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	464	Total	C	N	O	S	0	1	0
			3511	2185	635	673	18			
1	B	464	Total	C	N	O	S	0	1	0
			3511	2185	635	673	18			
1	C	464	Total	C	N	O	S	0	1	0
			3511	2185	635	673	18			
1	D	464	Total	C	N	O	S	0	1	0
			3511	2185	635	673	18			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P9WKE5
A	-1	GLY	-	expression tag	UNP P9WKE5
A	0	HIS	-	expression tag	UNP P9WKE5
B	-2	GLY	-	expression tag	UNP P9WKE5
B	-1	GLY	-	expression tag	UNP P9WKE5
B	0	HIS	-	expression tag	UNP P9WKE5
C	-2	GLY	-	expression tag	UNP P9WKE5
C	-1	GLY	-	expression tag	UNP P9WKE5
C	0	HIS	-	expression tag	UNP P9WKE5
D	-2	GLY	-	expression tag	UNP P9WKE5
D	-1	GLY	-	expression tag	UNP P9WKE5
D	0	HIS	-	expression tag	UNP P9WKE5

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

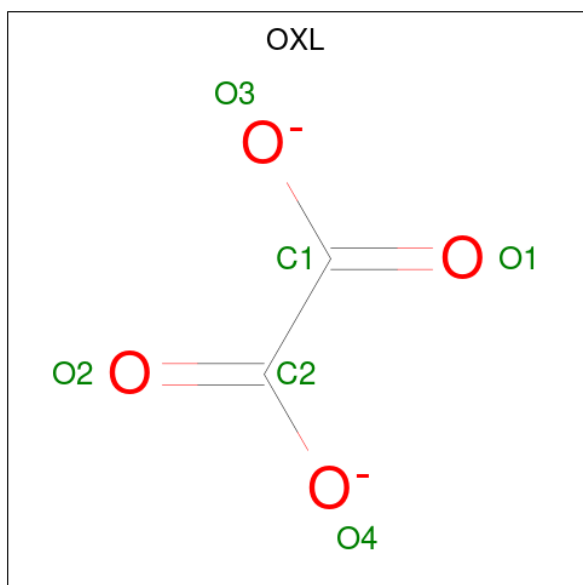
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	1	Total	Mg	0	0
			1	1		
2	D	1	Total	Mg	0	0
			1	1		

- Molecule 3 is OXALATE ION (CCD ID: OXL) (formula: C_2O_4).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	130	Total	O	0	0
			130	130		
4	B	110	Total	O	0	0
			110	110		
4	C	123	Total	O	0	0
			123	123		

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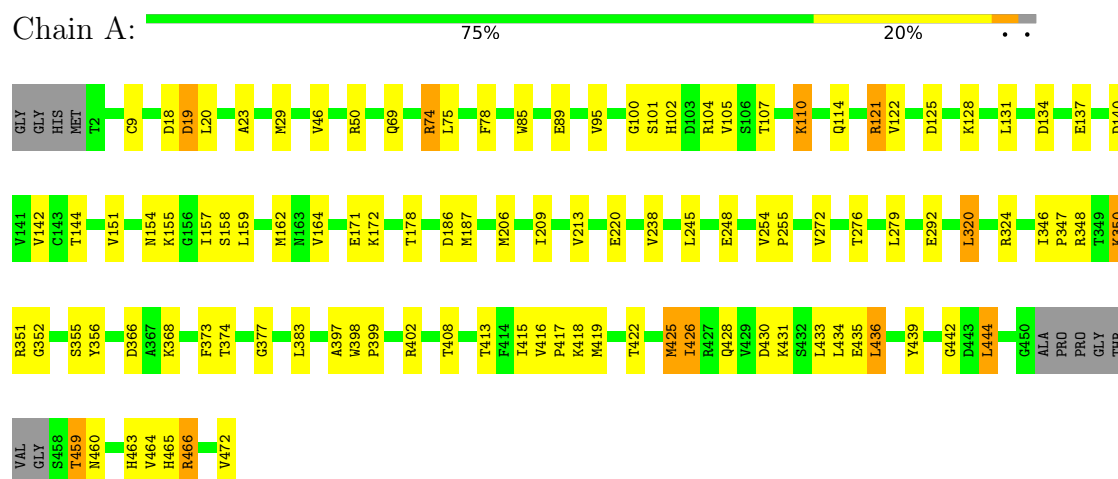
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	131	Total 131	O 131	0	0

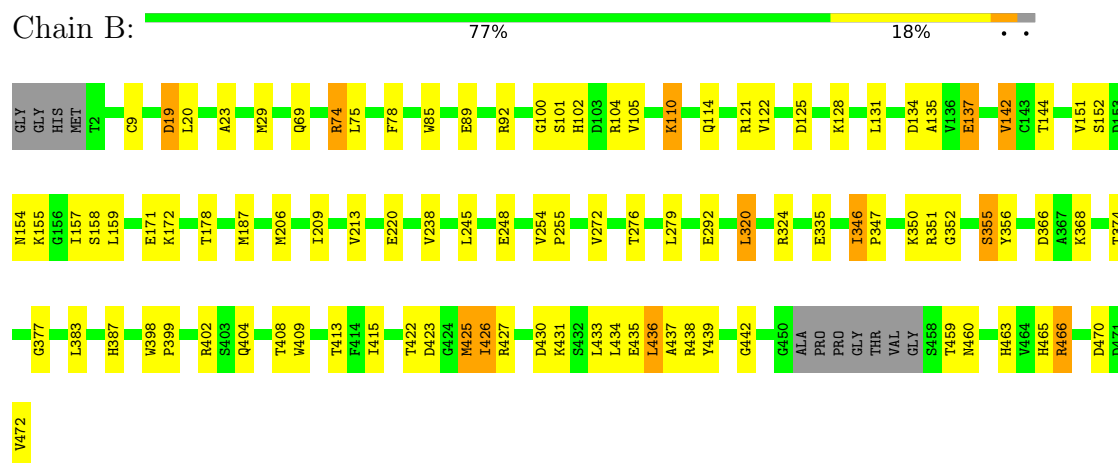
3 Residue-property plots

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.

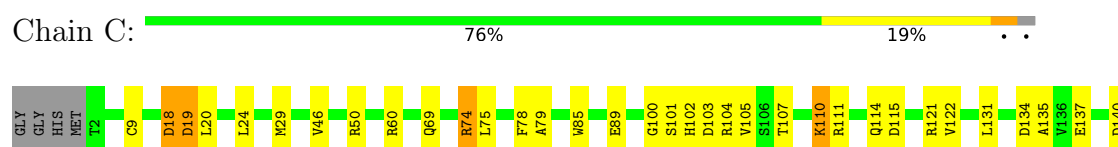
• Molecule 1: Pyruvate kinase

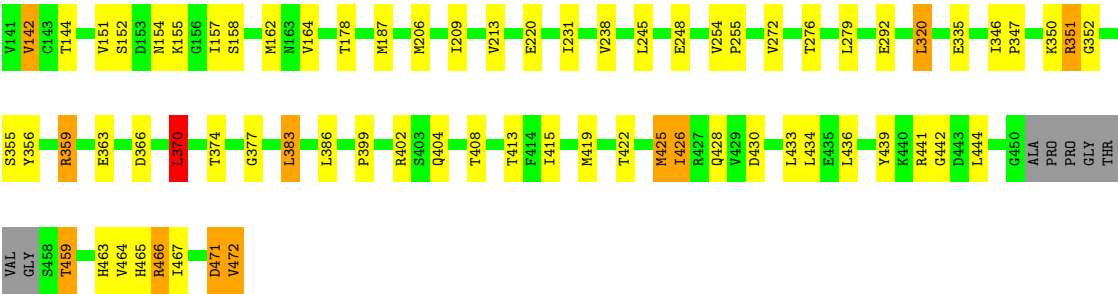


• Molecule 1: Pyruvate kinase

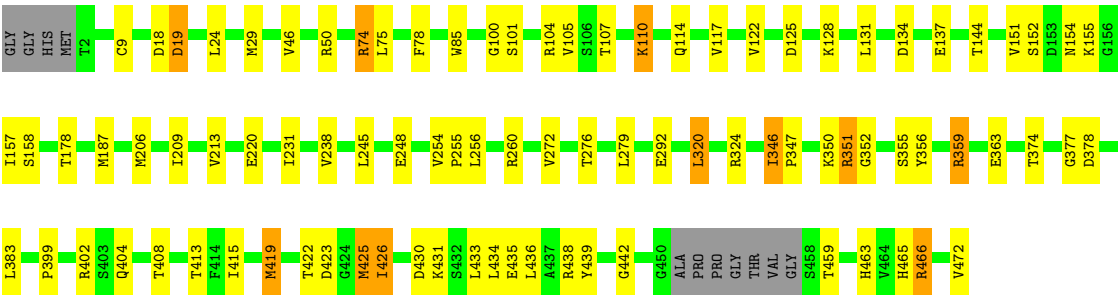
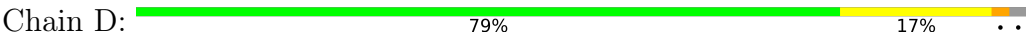


• Molecule 1: Pyruvate kinase





● Molecule 1: Pyruvate kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	127.01Å 127.01Å 144.65Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	63.51 – 2.62 63.51 – 2.62	Depositor EDS
% Data completeness (in resolution range)	100.0 (63.51-2.62) 100.0 (63.51-2.62)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.173 , 0.218 0.181 , 0.221	Depositor DCC
R_{free} test set	4012 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	49.6	Xtriage
Anisotropy	0.703	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 36.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.477 for -h,-k,l 0.477 for h,-h-k,-l 0.480 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14566	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% 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: OXL, 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	0.87	0/3558	0.82	0/4829
1	B	0.87	1/3558 (0.0%)	0.82	0/4829
1	C	0.89	1/3558 (0.0%)	0.83	0/4829
1	D	0.88	0/3558	0.82	0/4829
All	All	0.88	2/14232 (0.0%)	0.83	0/19316

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	370	LEU	N-CA	-5.29	1.39	1.45
1	B	346	ILE	N-CA	-5.03	1.41	1.46

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	3511	0	3573	76	0
1	B	3511	0	3573	58	0
1	C	3511	0	3573	70	0
1	D	3511	0	3573	64	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	6	0	0	0	0
3	B	6	0	0	0	0
3	C	6	0	0	0	0
3	D	6	0	0	0	0
4	A	130	0	0	5	0
4	B	110	0	0	3	0
4	C	123	0	0	8	0
4	D	131	0	0	7	0
All	All	14566	0	14292	267	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 9.

All (267) 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:351:ARG:NH2	1:D:378:ASP:OD1	1.82	1.12
1:C:430:ASP:O	1:C:434:LEU:HD12	1.58	1.03
1:B:320:LEU:HD23	1:B:324:ARG:NH1	1.82	0.94
1:D:347:PRO:HG3	1:D:356:TYR:CE2	2.04	0.91
1:D:125:ASP:OD2	1:D:128:LYS:HD2	1.70	0.91
1:A:122:VAL:HG22	1:A:159:LEU:HD11	1.52	0.89
1:B:423:ASP:O	1:B:427:ARG:HG3	1.76	0.85
1:B:92:ARG:NH1	1:B:137:GLU:OE2	2.10	0.84
1:A:95:VAL:HG13	1:A:110:LYS:HA	1.58	0.83
1:A:320:LEU:HD23	1:A:324:ARG:NH1	1.94	0.81
1:C:135:ALA:O	1:C:142:VAL:HG23	1.82	0.79
1:A:69:GLN:OE1	1:A:74:ARG:NH1	2.17	0.77
1:A:418:LYS:H	1:A:418:LYS:HD2	1.50	0.77
1:C:85:TRP:CE3	1:C:151:VAL:HG11	2.21	0.76
1:C:347:PRO:HG3	1:C:356:TYR:CE2	2.21	0.75
1:B:85:TRP:CE3	1:B:151:VAL:HG11	2.22	0.75
1:A:347:PRO:HG3	1:A:356:TYR:CE2	2.21	0.75
1:D:85:TRP:CE3	1:D:151:VAL:HG11	2.22	0.75
1:A:85:TRP:CE3	1:A:151:VAL:HG11	2.22	0.75
1:B:347:PRO:HG3	1:B:356:TYR:CE2	2.22	0.74
1:B:436:LEU:HD12	1:B:436:LEU:N	2.03	0.73
1:A:122:VAL:HG22	1:A:159:LEU:CD1	2.17	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:18:ASP:HA	1:C:50:ARG:HH12	1.54	0.73
1:A:436:LEU:N	1:A:436:LEU:HD12	2.05	0.71
1:A:402:ARG:HG3	1:A:413:THR:HB	1.72	0.71
1:C:402:ARG:HG3	1:C:413:THR:HB	1.73	0.70
1:D:402:ARG:HG3	1:D:413:THR:HB	1.73	0.70
1:B:402:ARG:HG3	1:B:413:THR:HB	1.75	0.69
1:A:419:MET:SD	1:A:428:GLN:HG2	2.33	0.68
1:D:74:ARG:NH1	1:D:154:ASN:O	2.26	0.68
1:B:426:ILE:O	1:B:430:ASP:OD2	2.13	0.66
1:D:426:ILE:O	1:D:430:ASP:OD2	2.13	0.66
1:A:158:SER:OG	1:A:248:GLU:OE1	2.14	0.66
1:D:158:SER:OG	1:D:248:GLU:OE1	2.14	0.66
1:B:158:SER:OG	1:B:248:GLU:OE1	2.14	0.66
1:C:426:ILE:O	1:C:430:ASP:OD2	2.13	0.66
1:C:436:LEU:N	1:C:436:LEU:HD12	2.10	0.66
1:C:158:SER:OG	1:C:248:GLU:OE1	2.14	0.66
1:A:426:ILE:O	1:A:430:ASP:OD2	2.14	0.66
1:B:134:ASP:OD2	1:B:144:THR:OG1	2.15	0.65
1:D:125:ASP:OD2	1:D:128:LYS:CD	2.44	0.65
1:A:178:THR:OG1	1:A:209:ILE:HD11	1.97	0.65
1:C:430:ASP:C	1:C:434:LEU:HD12	2.22	0.65
1:A:134:ASP:OD2	1:A:144:THR:OG1	2.16	0.64
1:C:430:ASP:O	1:C:434:LEU:CD1	2.40	0.64
1:C:134:ASP:OD2	1:C:144:THR:OG1	2.15	0.64
1:D:134:ASP:OD2	1:D:144:THR:OG1	2.15	0.64
1:D:436:LEU:N	1:D:436:LEU:HD12	2.12	0.64
1:C:140:ASP:HB2	4:C:621:HOH:O	1.97	0.63
1:A:238:VAL:HG23	1:A:272:VAL:HG23	1.81	0.62
1:C:178:THR:OG1	1:C:209:ILE:HD11	1.99	0.62
1:C:104:ARG:NH2	4:C:601:HOH:O	2.32	0.62
1:B:238:VAL:HG23	1:B:272:VAL:HG23	1.82	0.61
1:C:238:VAL:HG23	1:C:272:VAL:HG23	1.81	0.61
1:D:260:ARG:HD3	4:D:675:HOH:O	2.01	0.61
1:D:178:THR:OG1	1:D:209:ILE:HD11	2.01	0.61
1:D:238:VAL:HG23	1:D:272:VAL:HG23	1.81	0.61
1:B:69:GLN:OE1	1:B:74:ARG:NH1	2.34	0.61
1:B:178:THR:OG1	1:B:209:ILE:HD11	2.01	0.61
1:C:441:ARG:HA	1:C:467:ILE:HG22	1.81	0.61
1:D:346:ILE:H	1:D:346:ILE:HD12	1.66	0.60
1:A:140:ASP:HB2	4:A:628:HOH:O	2.03	0.59
1:C:78:PHE:CZ	1:C:105:VAL:HG21	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:78:PHE:CZ	1:B:105:VAL:HG21	2.38	0.58
1:A:416:VAL:HG21	1:A:419:MET:HE1	1.84	0.58
1:C:69:GLN:OE1	1:C:74:ARG:NH2	2.36	0.58
1:A:78:PHE:CZ	1:A:105:VAL:HG21	2.39	0.58
1:D:347:PRO:CG	1:D:356:TYR:CE2	2.83	0.57
1:C:29:MET:O	1:C:60:ARG:NH2	2.37	0.57
1:B:125:ASP:OD2	1:B:128:LYS:HG3	2.04	0.57
1:D:78:PHE:CZ	1:D:105:VAL:HG21	2.39	0.57
1:A:419:MET:CE	1:A:428:GLN:HG2	2.36	0.56
1:D:78:PHE:HA	1:D:100:GLY:O	2.07	0.55
1:D:104:ARG:NH2	4:D:602:HOH:O	2.39	0.55
1:A:422:THR:HA	1:A:425:MET:HG3	1.88	0.55
1:A:95:VAL:HG13	1:A:110:LYS:HD2	1.89	0.55
1:A:399:PRO:HA	1:A:415:ILE:HD12	1.87	0.55
1:D:347:PRO:CG	1:D:356:TYR:CD2	2.90	0.55
1:C:399:PRO:HA	1:C:415:ILE:HD12	1.88	0.55
1:D:355:SER:HB3	1:D:383:LEU:HD13	1.87	0.55
1:A:151:VAL:HG23	1:A:151:VAL:O	2.07	0.54
1:C:370:LEU:HD11	1:C:383:LEU:HD22	1.89	0.54
1:A:122:VAL:HB	1:A:131:LEU:HB2	1.89	0.54
1:A:460:ASN:ND2	1:C:471:ASP:HB2	2.23	0.54
1:B:347:PRO:CG	1:B:356:TYR:CE2	2.89	0.54
1:C:121:ARG:HD2	4:D:661:HOH:O	2.08	0.54
1:A:347:PRO:CG	1:A:356:TYR:CE2	2.90	0.54
1:B:347:PRO:HG2	1:B:356:TYR:CD2	2.42	0.54
1:C:347:PRO:HG2	1:C:356:TYR:CD2	2.43	0.54
1:B:351:ARG:O	1:B:355:SER:OG	2.25	0.54
1:B:399:PRO:HA	1:B:415:ILE:HD12	1.89	0.53
1:C:422:THR:HA	1:C:425:MET:HG3	1.89	0.53
1:C:19:ASP:HA	4:C:622:HOH:O	2.07	0.53
1:D:347:PRO:HG3	1:D:356:TYR:CD2	2.43	0.53
1:D:399:PRO:HA	1:D:415:ILE:HD12	1.90	0.53
1:A:78:PHE:HA	1:A:100:GLY:O	2.09	0.53
1:A:19:ASP:HB2	4:A:610:HOH:O	2.09	0.53
1:C:347:PRO:CG	1:C:356:TYR:CE2	2.90	0.52
1:A:347:PRO:HG2	1:A:356:TYR:CD2	2.44	0.52
1:D:75:LEU:O	1:D:154:ASN:HA	2.10	0.52
1:B:436:LEU:N	1:B:436:LEU:CD1	2.72	0.52
1:A:347:PRO:CG	1:A:356:TYR:CD2	2.93	0.52
1:B:78:PHE:HA	1:B:100:GLY:O	2.10	0.52
1:C:75:LEU:O	1:C:154:ASN:HA	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:422:THR:HA	1:D:425:MET:HG3	1.91	0.52
1:B:347:PRO:CG	1:B:356:TYR:CD2	2.94	0.51
1:A:75:LEU:O	1:A:154:ASN:HA	2.10	0.51
1:A:436:LEU:N	1:A:436:LEU:CD1	2.73	0.51
1:B:422:THR:HA	1:B:425:MET:HG3	1.92	0.51
1:D:46:VAL:O	1:D:50:ARG:HG3	2.11	0.51
1:A:95:VAL:HG11	1:A:110:LYS:O	2.10	0.51
1:A:125:ASP:OD2	1:A:128:LYS:HG3	2.10	0.51
1:B:75:LEU:O	1:B:154:ASN:HA	2.10	0.51
1:D:18:ASP:HA	1:D:50:ARG:HH22	1.74	0.51
1:D:122:VAL:HB	1:D:131:LEU:HB2	1.91	0.51
1:C:347:PRO:CG	1:C:356:TYR:CD2	2.93	0.51
1:C:79:ALA:HB2	4:C:604:HOH:O	2.10	0.51
1:B:151:VAL:HG23	1:B:151:VAL:O	2.10	0.51
1:C:151:VAL:HG23	1:C:151:VAL:O	2.10	0.50
1:C:436:LEU:N	1:C:436:LEU:CD1	2.75	0.50
1:D:151:VAL:HG23	1:D:151:VAL:O	2.11	0.50
1:C:89:GLU:OE2	1:C:102:HIS:CE1	2.64	0.50
1:A:23:ALA:HB1	1:A:320:LEU:HD11	1.94	0.50
1:C:78:PHE:HA	1:C:100:GLY:O	2.12	0.49
1:D:78:PHE:HB3	4:D:687:HOH:O	2.11	0.49
1:A:373:PHE:HE2	1:A:419:MET:CE	2.26	0.49
1:D:419:MET:HB3	1:D:425:MET:HE2	1.94	0.49
1:B:352:GLY:HA3	4:B:644:HOH:O	2.13	0.49
1:C:352:GLY:HA3	4:C:666:HOH:O	2.13	0.49
1:D:436:LEU:N	1:D:436:LEU:CD1	2.76	0.49
1:D:430:ASP:O	1:D:434:LEU:HD12	2.14	0.48
1:A:426:ILE:HD13	1:A:472:VAL:HG22	1.95	0.48
1:B:135:ALA:O	1:B:142:VAL:HG23	2.14	0.48
1:C:374:THR:HG21	1:C:377:GLY:HA2	1.95	0.48
1:C:101:SER:O	1:C:104:ARG:O	2.32	0.48
1:B:101:SER:O	1:B:104:ARG:O	2.32	0.48
1:D:374:THR:HG21	1:D:377:GLY:HA2	1.96	0.48
1:D:101:SER:O	1:D:104:ARG:O	2.31	0.48
1:B:19:ASP:HA	4:B:619:HOH:O	2.14	0.48
1:D:433:LEU:O	1:D:439:TYR:HB2	2.14	0.48
1:D:347:PRO:HG2	1:D:356:TYR:CD2	2.48	0.47
1:A:397:ALA:HB1	1:A:418:LYS:HE3	1.95	0.47
1:B:426:ILE:HD13	1:B:472:VAL:HG22	1.96	0.47
1:C:350:LYS:NZ	1:C:459:THR:O	2.47	0.47
1:D:426:ILE:HD13	1:D:472:VAL:HG22	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:VAL:HB	1:B:131:LEU:HB2	1.96	0.47
1:C:89:GLU:OE2	1:C:102:HIS:NE2	2.47	0.47
1:B:430:ASP:O	1:B:434:LEU:HD12	2.14	0.47
1:C:419:MET:CE	1:C:428:GLN:HG2	2.44	0.47
1:A:186:ASP:HB2	1:A:187:MET:HE3	1.97	0.47
1:A:374:THR:HG21	1:A:377:GLY:HA2	1.95	0.46
1:A:110:LYS:HA	1:A:110:LYS:HD2	1.66	0.46
1:A:121:ARG:HG2	4:B:662:HOH:O	2.14	0.46
1:D:320:LEU:HD23	1:D:324:ARG:HH12	1.80	0.46
1:D:350:LYS:NZ	4:D:605:HOH:O	2.42	0.46
1:A:430:ASP:O	1:A:434:LEU:HD12	2.14	0.46
1:B:374:THR:HG21	1:B:377:GLY:HA2	1.96	0.46
1:A:19:ASP:HA	4:A:616:HOH:O	2.15	0.46
1:D:9:CYS:HB2	1:D:29:MET:SD	2.56	0.46
1:B:89:GLU:OE1	1:B:102:HIS:NE2	2.49	0.46
1:C:351:ARG:O	1:C:355:SER:OG	2.33	0.46
1:D:351:ARG:CZ	1:D:378:ASP:OD1	2.59	0.46
1:B:171:GLU:OE2	1:B:171:GLU:HA	2.16	0.46
1:C:9:CYS:HB2	1:C:29:MET:SD	2.56	0.46
1:C:370:LEU:HD11	1:C:383:LEU:CD2	2.45	0.46
1:A:463:HIS:CD2	1:A:465:HIS:HB2	2.50	0.46
1:B:426:ILE:HD13	1:B:472:VAL:HG13	1.98	0.46
1:A:442:GLY:O	1:A:466:ARG:HG3	2.16	0.45
1:C:433:LEU:O	1:C:439:TYR:HB2	2.16	0.45
1:D:110:LYS:HD2	1:D:110:LYS:HA	1.70	0.45
1:B:431:LYS:O	1:B:435:GLU:HG3	2.16	0.45
4:C:708:HOH:O	1:D:256:LEU:HD12	2.17	0.45
1:D:351:ARG:O	1:D:355:SER:OG	2.34	0.45
1:A:433:LEU:O	1:A:439:TYR:HB2	2.17	0.45
1:C:279:LEU:HD23	1:C:292:GLU:HB3	1.99	0.45
1:D:426:ILE:HD13	1:D:472:VAL:HG13	1.99	0.45
1:B:463:HIS:CD2	1:B:465:HIS:HB2	2.51	0.45
1:D:431:LYS:O	1:D:435:GLU:HG3	2.16	0.45
1:C:463:HIS:CD2	1:C:465:HIS:HB2	2.52	0.44
1:D:279:LEU:HD23	1:D:292:GLU:HB3	1.98	0.44
1:A:89:GLU:OE1	1:A:102:HIS:CE1	2.70	0.44
1:A:347:PRO:HG3	1:A:356:TYR:CD2	2.52	0.44
1:C:110:LYS:HD2	1:C:110:LYS:HA	1.68	0.44
1:C:122:VAL:HB	1:C:131:LEU:HB2	1.98	0.44
1:C:351:ARG:HD3	4:C:689:HOH:O	2.17	0.44
1:A:431:LYS:O	1:A:435:GLU:HG3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:101:SER:N	1:B:104:ARG:O	2.51	0.44
1:C:442:GLY:O	1:C:466:ARG:HG3	2.18	0.44
1:A:9:CYS:HB2	1:A:29:MET:SD	2.57	0.44
1:B:9:CYS:HB2	1:B:29:MET:SD	2.58	0.44
1:B:433:LEU:O	1:B:439:TYR:HB2	2.18	0.44
1:D:435:GLU:C	1:D:436:LEU:HD12	2.42	0.44
1:B:254:VAL:N	1:B:255:PRO:CD	2.81	0.44
1:B:279:LEU:HD23	1:B:292:GLU:HB3	1.99	0.44
1:C:101:SER:N	1:C:104:ARG:O	2.51	0.44
1:C:24:LEU:HD23	1:C:320:LEU:HD12	2.00	0.43
1:B:442:GLY:O	1:B:466:ARG:HG3	2.18	0.43
1:A:426:ILE:HD13	1:A:472:VAL:HG13	2.00	0.43
1:C:162:MET:HE3	1:C:164:VAL:HG22	2.01	0.43
1:D:320:LEU:HA	1:D:320:LEU:HD12	1.63	0.43
1:A:254:VAL:N	1:A:255:PRO:CD	2.82	0.43
1:A:350:LYS:HE2	1:A:459:THR:O	2.18	0.43
1:B:206:MET:HE2	1:B:213:VAL:HG22	2.01	0.43
1:C:346:ILE:HD12	1:C:347:PRO:HD3	2.01	0.43
1:A:46:VAL:O	1:A:50:ARG:HG3	2.18	0.43
1:A:78:PHE:HB3	4:A:687:HOH:O	2.19	0.43
1:B:423:ASP:OD1	1:B:472:VAL:HG11	2.18	0.43
1:B:110:LYS:HD2	1:B:110:LYS:HA	1.69	0.43
1:B:346:ILE:HD12	1:B:346:ILE:HA	1.91	0.43
1:B:346:ILE:HA	1:B:347:PRO:HD3	1.87	0.42
1:C:206:MET:HE2	1:C:213:VAL:HG22	2.00	0.42
1:D:125:ASP:O	1:D:128:LYS:HG3	2.19	0.42
1:A:95:VAL:CG1	1:A:110:LYS:HA	2.40	0.42
1:A:206:MET:HE2	1:A:213:VAL:HG22	2.01	0.42
1:A:279:LEU:HD23	1:A:292:GLU:HB3	2.01	0.42
1:C:347:PRO:HG3	1:C:356:TYR:CD2	2.53	0.42
1:C:419:MET:HE1	1:C:428:GLN:HG2	2.01	0.42
1:D:463:HIS:CD2	1:D:465:HIS:HB2	2.54	0.42
1:A:171:GLU:HA	1:A:171:GLU:OE2	2.17	0.42
1:A:346:ILE:HA	1:A:347:PRO:HD3	1.87	0.42
1:B:437:ALA:O	1:B:438:ARG:HB2	2.19	0.42
1:C:444:LEU:HD11	1:C:464:VAL:CG1	2.49	0.42
1:D:423:ASP:OD1	1:D:472:VAL:HG11	2.19	0.42
1:A:220:GLU:HA	1:A:245:LEU:HD13	2.00	0.42
1:A:348:ARG:CZ	1:A:348:ARG:HB2	2.49	0.42
1:C:187:MET:HE2	1:C:404:GLN:HA	2.00	0.42
1:B:220:GLU:HA	1:B:245:LEU:HD13	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:350:LYS:HE3	1:B:460:ASN:HB3	2.02	0.42
1:B:398:TRP:HA	1:B:399:PRO:HD3	1.85	0.42
1:C:335:GLU:HB2	4:C:633:HOH:O	2.19	0.42
1:D:254:VAL:N	1:D:255:PRO:CD	2.82	0.42
1:A:351:ARG:O	1:A:355:SER:OG	2.34	0.42
1:D:442:GLY:O	1:D:466:ARG:HG3	2.20	0.42
1:A:134:ASP:HB2	1:A:142:VAL:HG13	2.02	0.42
1:C:346:ILE:HA	1:C:347:PRO:HD3	1.86	0.42
1:D:107:THR:O	1:D:110:LYS:NZ	2.47	0.42
1:D:352:GLY:HA3	4:D:673:HOH:O	2.20	0.42
1:A:352:GLY:HA3	4:A:690:HOH:O	2.20	0.41
1:C:254:VAL:N	1:C:255:PRO:CD	2.82	0.41
1:A:463:HIS:HD2	1:A:465:HIS:HB2	1.85	0.41
1:A:23:ALA:CB	1:A:320:LEU:HD11	2.50	0.41
1:A:398:TRP:HA	1:A:399:PRO:HD3	1.84	0.41
1:C:107:THR:O	1:C:110:LYS:NZ	2.48	0.41
1:D:101:SER:N	1:D:104:ARG:O	2.53	0.41
1:D:220:GLU:HA	1:D:245:LEU:HD13	2.00	0.41
1:A:444:LEU:HD21	1:A:464:VAL:CG1	2.51	0.41
1:D:19:ASP:HA	4:D:613:HOH:O	2.19	0.41
1:D:187:MET:HE2	1:D:404:GLN:HA	2.01	0.41
1:B:23:ALA:CB	1:B:320:LEU:HD11	2.51	0.41
1:B:187:MET:HE2	1:B:404:GLN:HA	2.01	0.41
1:C:220:GLU:HA	1:C:245:LEU:HD13	2.01	0.41
1:D:206:MET:HE2	1:D:213:VAL:HG22	2.02	0.41
1:A:417:PRO:O	1:A:419:MET:SD	2.78	0.41
1:C:359:ARG:HD3	1:C:363:GLU:CD	2.46	0.41
1:C:426:ILE:HD13	1:C:472:VAL:HB	2.03	0.41
1:D:359:ARG:HD3	1:D:363:GLU:CD	2.46	0.41
1:A:162:MET:HE3	1:A:164:VAL:HG22	2.02	0.40
1:C:19:ASP:H	1:C:50:ARG:NH1	2.19	0.40
1:C:111:ARG:O	1:C:115:ASP:HB2	2.21	0.40
1:A:107:THR:O	1:A:110:LYS:NZ	2.46	0.40
1:A:121:ARG:O	1:A:159:LEU:HD13	2.21	0.40
1:B:320:LEU:HD23	1:B:324:ARG:CZ	2.48	0.40
1:C:46:VAL:O	1:C:50:ARG:HG3	2.21	0.40
1:D:24:LEU:HD23	1:D:320:LEU:HD12	2.02	0.40
1:A:101:SER:N	1:A:104:ARG:O	2.55	0.40
1:B:346:ILE:HD12	1:B:347:PRO:HD3	2.04	0.40
1:B:387:HIS:HA	1:B:409:TRP:CZ3	2.57	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	461/475 (97%)	445 (96%)	15 (3%)	1 (0%)	43	64
1	B	461/475 (97%)	446 (97%)	14 (3%)	1 (0%)	43	64
1	C	461/475 (97%)	444 (96%)	16 (4%)	1 (0%)	43	64
1	D	461/475 (97%)	447 (97%)	13 (3%)	1 (0%)	43	64
All	All	1844/1900 (97%)	1782 (97%)	58 (3%)	4 (0%)	43	64

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	276	THR
1	A	276	THR
1	B	276	THR
1	D	276	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	376/381 (99%)	353 (94%)	23 (6%)	17	35
1	B	376/381 (99%)	350 (93%)	26 (7%)	14	30
1	C	376/381 (99%)	349 (93%)	27 (7%)	13	28
1	D	376/381 (99%)	355 (94%)	21 (6%)	19	39
All	All	1504/1524 (99%)	1407 (94%)	97 (6%)	15	33

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ASP
1	A	19	ASP
1	A	20	LEU
1	A	74	ARG
1	A	110	LYS
1	A	114	GLN
1	A	121	ARG
1	A	137	GLU
1	A	155	LYS
1	A	157	ILE
1	A	172	LYS
1	A	320	LEU
1	A	350	LYS
1	A	366	ASP
1	A	368	LYS
1	A	383	LEU
1	A	408	THR
1	A	425	MET
1	A	426	ILE
1	A	436	LEU
1	A	444	LEU
1	A	459	THR
1	A	466	ARG
1	B	19	ASP
1	B	20	LEU
1	B	74	ARG
1	B	110	LYS
1	B	114	GLN
1	B	121	ARG
1	B	137	GLU
1	B	142	VAL
1	B	152	SER
1	B	155	LYS
1	B	157	ILE
1	B	159	LEU
1	B	172	LYS
1	B	320	LEU
1	B	335	GLU
1	B	355	SER
1	B	366	ASP
1	B	368	LYS
1	B	383	LEU

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Mol	Chain	Res	Type
1	B	408	THR
1	B	425	MET
1	B	426	ILE
1	B	436	LEU
1	B	459	THR
1	B	466	ARG
1	B	470	ASP
1	C	18	ASP
1	C	19	ASP
1	C	20	LEU
1	C	74	ARG
1	C	103	ASP
1	C	110	LYS
1	C	114	GLN
1	C	137	GLU
1	C	142	VAL
1	C	152	SER
1	C	155	LYS
1	C	157	ILE
1	C	231	ILE
1	C	320	LEU
1	C	351	ARG
1	C	359	ARG
1	C	366	ASP
1	C	370	LEU
1	C	383	LEU
1	C	386	LEU
1	C	408	THR
1	C	425	MET
1	C	426	ILE
1	C	459	THR
1	C	466	ARG
1	C	471	ASP
1	C	472	VAL
1	D	19	ASP
1	D	74	ARG
1	D	110	LYS
1	D	114	GLN
1	D	117	VAL
1	D	137	GLU
1	D	152	SER
1	D	155	LYS

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Mol	Chain	Res	Type
1	D	157	ILE
1	D	231	ILE
1	D	320	LEU
1	D	346	ILE
1	D	351	ARG
1	D	359	ARG
1	D	408	THR
1	D	419	MET
1	D	425	MET
1	D	426	ILE
1	D	438	ARG
1	D	459	THR
1	D	466	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	336	ASN
1	A	387	HIS
1	A	404	GLN
1	A	428	GLN
1	A	463	HIS
1	B	336	ASN
1	B	387	HIS
1	B	404	GLN
1	B	428	GLN
1	B	463	HIS
1	C	336	ASN
1	C	387	HIS
1	C	404	GLN
1	C	428	GLN
1	C	463	HIS
1	D	336	ASN
1	D	387	HIS
1	D	404	GLN
1	D	428	GLN

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 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	OXL	A	502	2	5,5,5	2.13	1 (20%)	6,6,6	2.66	2 (33%)
3	OXL	C	502	2	5,5,5	2.33	2 (40%)	6,6,6	2.65	2 (33%)
3	OXL	D	502	2	5,5,5	2.32	1 (20%)	6,6,6	1.71	2 (33%)
3	OXL	B	502	2	5,5,5	2.34	1 (20%)	6,6,6	1.66	2 (33%)

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	OXL	A	502	2	-	4/4/4/4	-
3	OXL	C	502	2	-	4/4/4/4	-
3	OXL	D	502	2	-	4/4/4/4	-
3	OXL	B	502	2	-	4/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	502	OXL	C2-C1	-4.23	1.47	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	502	OXL	C2-C1	-4.05	1.47	1.54
3	D	502	OXL	C2-C1	-3.99	1.47	1.54
3	A	502	OXL	C2-C1	-3.69	1.48	1.54
3	C	502	OXL	O3-C1	-2.06	1.25	1.30

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	502	OXL	O4-C2-C1	4.31	121.19	112.83
3	A	502	OXL	O4-C2-C1	4.29	121.16	112.83
3	A	502	OXL	O3-C1-C2	4.20	120.97	112.83
3	C	502	OXL	O3-C1-C2	4.14	120.86	112.83
3	B	502	OXL	O3-C1-C2	2.89	118.44	112.83
3	D	502	OXL	O3-C1-C2	2.87	118.39	112.83
3	D	502	OXL	O4-C2-C1	2.46	117.61	112.83
3	B	502	OXL	O4-C2-C1	2.19	117.08	112.83

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	OXL	O3-C1-C2-O4
3	C	502	OXL	O3-C1-C2-O4
3	A	502	OXL	O1-C1-C2-O2
3	B	502	OXL	O3-C1-C2-O4
3	D	502	OXL	O3-C1-C2-O4
3	C	502	OXL	O1-C1-C2-O2
3	B	502	OXL	O1-C1-C2-O2
3	D	502	OXL	O1-C1-C2-O2
3	A	502	OXL	O1-C1-C2-O4
3	A	502	OXL	O3-C1-C2-O2
3	C	502	OXL	O3-C1-C2-O2
3	C	502	OXL	O1-C1-C2-O4
3	B	502	OXL	O1-C1-C2-O4
3	B	502	OXL	O3-C1-C2-O2
3	D	502	OXL	O1-C1-C2-O4
3	D	502	OXL	O3-C1-C2-O2

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	464/475 (97%)	-1.24	0 100 100	41, 65, 112, 139	1 (0%)
1	B	464/475 (97%)	-1.24	0 100 100	41, 65, 114, 140	1 (0%)
1	C	464/475 (97%)	-1.23	0 100 100	40, 65, 113, 143	1 (0%)
1	D	464/475 (97%)	-1.22	0 100 100	39, 65, 113, 137	1 (0%)
All	All	1856/1900 (97%)	-1.23	0 100 100	39, 65, 114, 143	4 (0%)

There are no RSRZ outliers to report.

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
3	OXL	A	502	6/6	0.99	0.04	53,62,74,78	0
3	OXL	B	502	6/6	0.99	0.04	56,61,73,80	0
3	OXL	D	502	6/6	0.99	0.04	54,64,83,86	0
2	MG	D	501	1/1	1.00	0.02	53,53,53,53	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MG	A	501	1/1	1.00	0.01	53,53,53,53	0
2	MG	B	501	1/1	1.00	0.02	53,53,53,53	0
3	OXL	C	502	6/6	1.00	0.05	53,60,78,79	0
2	MG	C	501	1/1	1.00	0.01	55,55,55,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.