



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 11:27 PM EDT

PDB ID : 4I1D / pdb_00004i1d
Title : The crystal structure of an ABC transporter substrate-binding protein from *Bradyrhizobium japonicum* USDA 110
Authors : Fan, Y.; Tan, K.; Mack, J.; Zerbs, S.; Collart, F.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2012-11-20
Resolution : 2.20 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

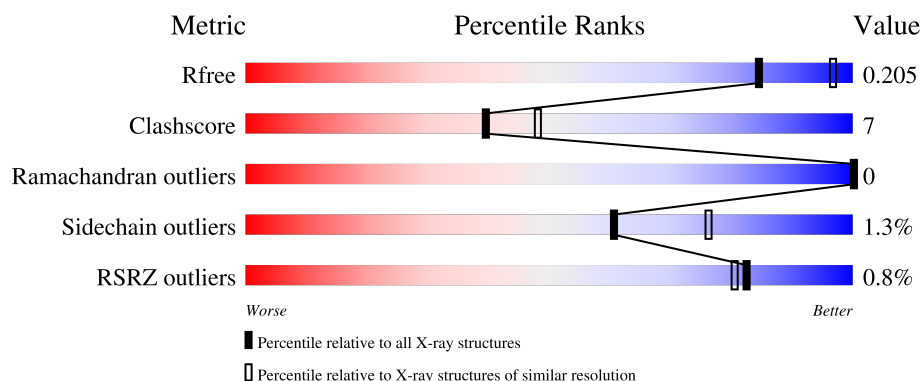
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	324	88% 11% ..
1	B	324	86% 13% ..
1	C	324	87% 12% .
1	D	324	87% 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
3	GOL	C	412	-	-	X	-
3	GOL	D	413	-	-	X	-
4	ACT	A	413	-	-	X	-
5	MLT	C	422	-	-	X	-
5	MLT	D	417	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10574 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 ABC transporter substrate-binding protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	Se	0	0	0
			2483	1590	419	466	2	6			
1	B	322	Total	C	N	O	S	Se	0	3	0
			2489	1594	418	469	2	6			
1	C	322	Total	C	N	O	S	Se	0	1	0
			2487	1594	418	467	2	6			
1	D	322	Total	C	N	O	S	Se	0	4	0
			2500	1603	419	470	2	6			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	41	SER	-	expression tag	UNP Q89F76
A	42	ASN	-	expression tag	UNP Q89F76
A	43	ALA	-	expression tag	UNP Q89F76
B	41	SER	-	expression tag	UNP Q89F76
B	42	ASN	-	expression tag	UNP Q89F76
B	43	ALA	-	expression tag	UNP Q89F76
C	41	SER	-	expression tag	UNP Q89F76
C	42	ASN	-	expression tag	UNP Q89F76
C	43	ALA	-	expression tag	UNP Q89F76
D	41	SER	-	expression tag	UNP Q89F76
D	42	ASN	-	expression tag	UNP Q89F76
D	43	ALA	-	expression tag	UNP Q89F76

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



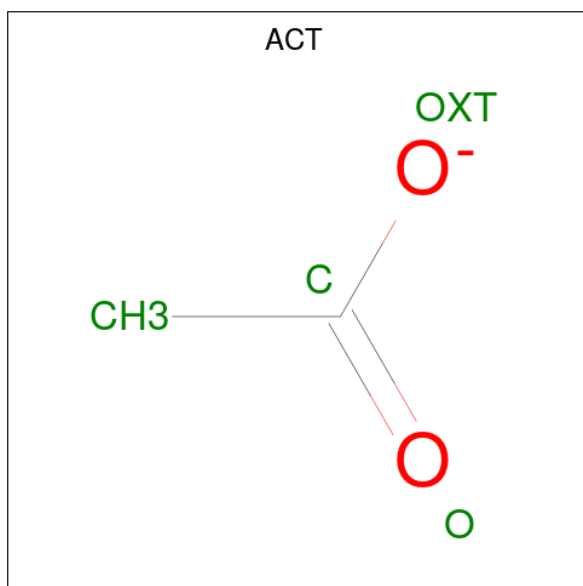
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total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

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

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



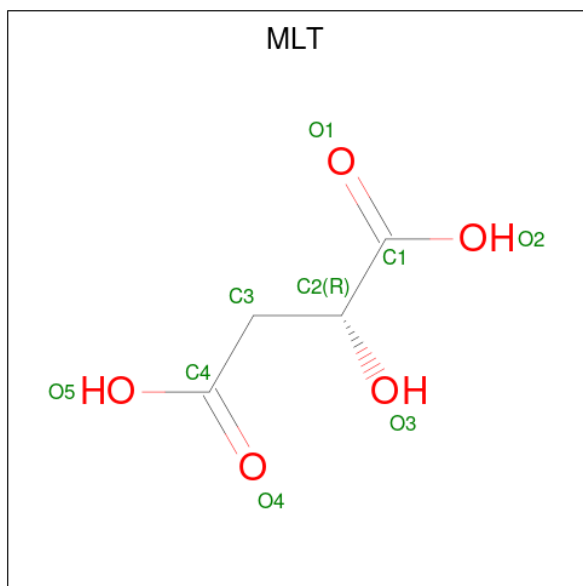
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total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

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

- Molecule 5 is D-MALATE (CCD ID: MLT) (formula: C₄H₆O₅).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			9	4	5		
5	D	1	Total	C	O	0	0
			9	4	5		

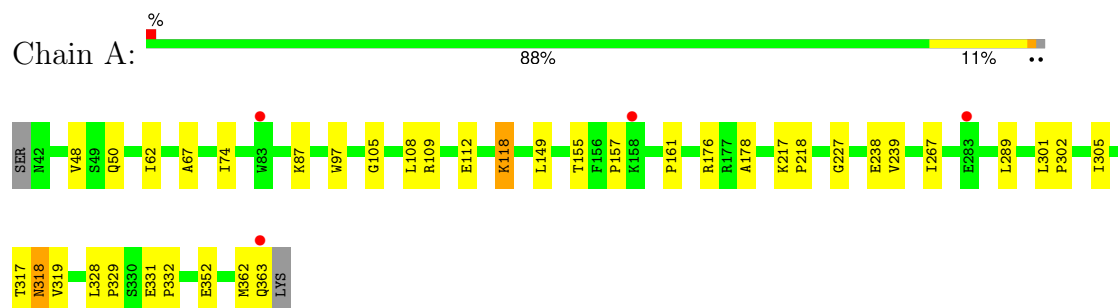
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	42	Total	O	0	0
			42	42		
6	B	65	Total	O	0	0
			65	65		
6	C	61	Total	O	0	0
			61	61		
6	D	75	Total	O	0	0
			75	75		

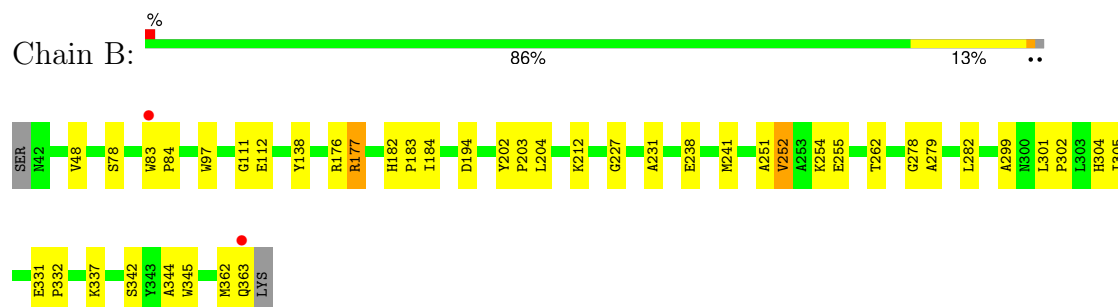
3 Residue-property plots

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.

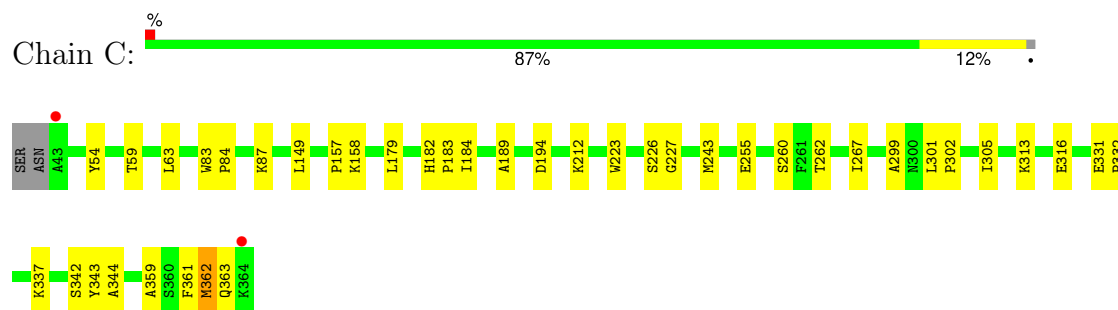
- Molecule 1: ABC transporter substrate-binding protein



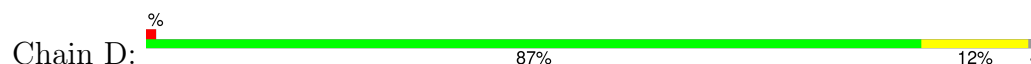
- Molecule 1: ABC transporter substrate-binding protein

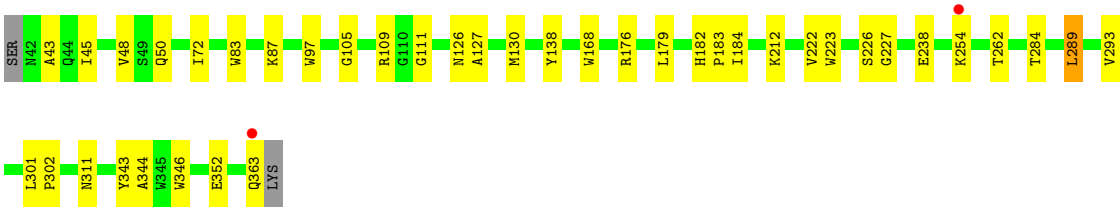


- Molecule 1: ABC transporter substrate-binding protein



- Molecule 1: ABC transporter substrate-binding protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	136.64Å 136.64Å 185.63Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.89 – 2.20 49.89 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.8 (49.89-2.20) 98.8 (49.89-2.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.16 (at 2.20Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, R_{free}	0.172 , 0.209 0.170 , 0.205	Depositor DCC
R_{free} test set	5027 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	46.3	Xtriage
Anisotropy	0.390	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.021 for -h,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10574	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% 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, SO4, MLT, ACT

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	0.43	0/2546	0.78	1/3462 (0.0%)
1	B	0.49	0/2560	0.74	0/3481
1	C	0.46	0/2553	0.74	0/3471
1	D	0.51	0/2575	0.78	0/3502
All	All	0.47	0/10234	0.76	1/13916 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	62	ILE	N-CA-C	6.43	116.70	111.62

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	2483	0	2437	27	0
1	B	2489	0	2444	36	0
1	C	2487	0	2441	39	0
1	D	2500	0	2467	35	0
2	A	45	0	0	2	0
2	B	35	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	45	0	0	0	0
2	D	45	0	0	1	0
3	A	6	0	8	0	0
3	B	30	0	40	7	0
3	C	30	0	40	8	0
3	D	24	0	32	5	0
4	A	24	0	18	3	0
4	B	12	0	9	0	0
4	C	28	0	21	1	0
4	D	12	0	9	1	0
5	A	9	0	4	3	0
5	B	9	0	4	3	0
5	C	9	0	4	5	0
5	D	9	0	4	4	0
6	A	42	0	0	0	0
6	B	65	0	0	3	0
6	C	61	0	0	1	0
6	D	75	0	0	0	0
All	All	10574	0	9982	136	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 7.

All (136) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:227:GLY:H	5:B:416:MLT:H31	1.06	1.08
1:C:227:GLY:H	5:C:422:MLT:H31	1.20	1.02
1:A:227:GLY:H	5:A:417:MLT:H2	1.26	0.99
1:B:282:LEU:HB3	3:B:412:GOL:H2	1.57	0.87
1:D:130:MSE:HE1	1:D:293:VAL:CG2	2.03	0.87
1:C:362:MSE:HA	1:C:362:MSE:HE2	1.56	0.86
1:B:227:GLY:N	5:B:416:MLT:H31	1.91	0.85
1:C:343:TYR:H	3:C:412:GOL:H32	1.41	0.84
1:C:227:GLY:H	5:C:422:MLT:C3	1.92	0.82
1:C:227:GLY:N	5:C:422:MLT:H31	1.97	0.80
1:D:130:MSE:HE2	1:D:311:ASN:ND2	1.96	0.80
1:B:182:HIS:HE1	1:B:184:ILE:HD12	1.47	0.79
1:A:362:MSE:O	1:A:363:GLN:HB2	1.83	0.79
1:B:342:SER:HA	3:B:409:GOL:H11	1.66	0.78
1:B:362:MSE:O	1:B:363:GLN:HB2	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:130:MSE:HE1	1:D:293:VAL:HB	1.69	0.73
1:D:227:GLY:H	5:D:417:MLT:C3	2.03	0.70
1:D:130:MSE:HE1	1:D:293:VAL:CB	2.21	0.69
1:D:127:ALA:HA	1:D:130:MSE:HE3	1.75	0.68
1:B:251:ALA:O	1:B:255[B]:GLU:HG2	1.93	0.68
1:D:126:ASN:O	1:D:130:MSE:HE3	1.96	0.65
1:D:344:ALA:N	2:D:407:SO4:O1	2.28	0.65
1:C:362:MSE:HA	1:C:362:MSE:CE	2.27	0.65
1:B:337:LYS:NZ	3:B:408:GOL:H11	2.12	0.65
1:D:301:LEU:HB3	1:D:302:PRO:HD3	1.80	0.64
1:C:343:TYR:H	3:C:412:GOL:C3	2.10	0.64
1:D:130:MSE:HE1	1:D:293:VAL:HG21	1.78	0.64
1:A:227:GLY:N	5:A:417:MLT:H2	2.06	0.62
1:B:78:SER:HB3	4:D:416:ACT:H3	1.82	0.62
1:B:304:HIS:HE1	6:B:530:HOH:O	1.82	0.62
1:C:359:ALA:O	1:C:363:GLN:HG2	1.99	0.62
1:B:337:LYS:HZ3	3:B:408:GOL:H11	1.65	0.61
1:B:182:HIS:CE1	1:B:184:ILE:HD12	2.33	0.61
1:B:344:ALA:H	3:B:409:GOL:H2	1.66	0.61
1:D:343:TYR:HE1	3:D:413:GOL:H12	1.66	0.61
1:C:344:ALA:H	3:C:412:GOL:H11	1.66	0.60
1:C:301:LEU:HB3	1:C:302:PRO:HD3	1.84	0.59
1:A:87:LYS:NZ	2:A:408:SO4:O3	2.37	0.58
1:C:262:THR:OG1	3:C:410:GOL:H2	2.05	0.57
1:B:83:TRP:HB3	1:B:84:PRO:HD3	1.86	0.57
1:D:262:THR:OG1	3:D:410:GOL:H2	2.05	0.57
1:D:184:ILE:HD13	3:D:413:GOL:H11	1.87	0.57
1:C:337:LYS:HZ2	3:C:410:GOL:H11	1.69	0.55
1:B:176:ARG:HD2	6:B:549:HOH:O	2.06	0.55
1:D:182:HIS:ND1	1:D:183:PRO:HD2	2.22	0.55
1:C:344:ALA:H	3:C:412:GOL:C1	2.19	0.55
1:D:48[B]:VAL:HG21	1:D:97:TRP:CD2	2.42	0.54
1:C:337:LYS:NZ	3:C:410:GOL:H11	2.23	0.54
1:B:282:LEU:CB	3:B:412:GOL:H2	2.34	0.54
1:C:194:ASP:HB2	1:C:212:LYS:HG3	1.90	0.54
1:C:361:PHE:CE2	1:C:362:MSE:HE3	2.43	0.54
1:D:226:SER:HA	5:D:417:MLT:H31	1.90	0.54
1:C:313:LYS:O	1:C:316[A]:GLU:HG2	2.09	0.53
1:C:149:LEU:C	1:C:149:LEU:HD23	2.34	0.52
1:D:130:MSE:HE2	1:D:311:ASN:HD22	1.72	0.52
1:D:105:GLY:HA3	3:D:413:GOL:O2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:182:HIS:CE1	1:C:184:ILE:HG12	2.45	0.52
1:C:157:PRO:HG2	1:C:158:LYS:HD2	1.92	0.52
1:A:87:LYS:HE3	4:A:412:ACT:H3	1.91	0.51
1:D:363:GLN:OE1	1:D:363:GLN:N	2.42	0.51
1:A:108:LEU:O	1:A:112:GLU:HG3	2.09	0.51
1:B:278:GLY:O	1:B:279:ALA:C	2.54	0.51
1:D:227:GLY:H	5:D:417:MLT:H32	1.76	0.50
1:B:301:LEU:HB3	1:B:302:PRO:HD3	1.91	0.50
1:B:48:VAL:HG21	1:B:97:TRP:CD2	2.47	0.50
1:D:48[B]:VAL:HG21	1:D:97:TRP:CG	2.46	0.50
1:B:182:HIS:CG	1:B:183:PRO:HD2	2.46	0.50
1:A:178:ALA:HB2	1:A:239:VAL:HG11	1.93	0.50
1:A:301:LEU:N	1:A:302:PRO:HD2	2.27	0.50
1:B:194:ASP:HB2	1:B:212:LYS:HG3	1.93	0.49
1:C:299:ALA:O	1:C:302:PRO:HD2	2.13	0.49
1:A:267:ILE:HD12	1:A:267:ILE:N	2.28	0.49
1:D:227:GLY:H	5:D:417:MLT:H31	1.75	0.49
1:A:331:GLU:HG2	1:A:332:PRO:HD2	1.95	0.48
1:B:48:VAL:HG21	1:B:97:TRP:CG	2.47	0.48
1:D:43:ALA:HB3	1:D:72:ILE:HG22	1.95	0.48
1:A:227:GLY:H	5:A:417:MLT:C2	2.12	0.48
1:B:362:MSE:O	1:B:363:GLN:CB	2.60	0.48
1:C:226:SER:HA	5:C:422:MLT:H31	1.96	0.48
1:B:331:GLU:HG2	1:B:332:PRO:HD2	1.96	0.48
1:C:83:TRP:HB3	1:C:84:PRO:HD3	1.96	0.48
1:A:105:GLY:HA3	4:A:413:ACT:H1	1.96	0.47
1:A:149:LEU:HD21	1:A:161:PRO:HB2	1.97	0.47
1:D:111:GLY:HA3	1:D:138:TYR:CD2	2.49	0.47
1:C:361:PHE:CD2	1:C:362:MSE:HE3	2.50	0.47
1:B:111:GLY:HA3	1:B:138:TYR:CD2	2.50	0.47
1:A:302:PRO:HA	1:A:305:ILE:O	2.15	0.46
1:C:316[A]:GLU:HG3	6:C:558:HOH:O	2.15	0.46
1:C:331:GLU:HG2	1:C:332:PRO:HD2	1.97	0.46
1:B:299:ALA:O	1:B:302:PRO:HD2	2.15	0.46
1:C:342:SER:HA	3:C:412:GOL:H31	1.97	0.46
1:D:346:TRP:HB3	3:D:413:GOL:H32	1.98	0.46
1:D:182:HIS:CG	1:D:183:PRO:HD2	2.51	0.46
1:A:67:ALA:HB2	1:A:74:ILE:HG13	1.98	0.45
1:A:118:LYS:HE3	1:A:118:LYS:HB2	1.67	0.45
1:B:227:GLY:H	5:B:416:MLT:C3	1.99	0.45
1:B:204:LEU:HD12	1:B:345:TRP:CZ3	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:302:PRO:HA	1:B:305:ILE:O	2.17	0.44
1:C:255:GLU:O	1:C:255:GLU:HG2	2.18	0.44
1:A:176:ARG:NH2	2:A:409:SO4:O2	2.51	0.44
1:C:179:LEU:O	1:C:223:TRP:HA	2.19	0.43
1:C:189:ALA:HB2	1:C:243:MSE:HE1	1.99	0.43
1:C:83:TRP:CE2	1:C:87:LYS:HD2	2.53	0.43
1:A:105:GLY:HA3	4:A:413:ACT:OXT	2.18	0.43
1:C:59:THR:HA	1:C:63:LEU:HB2	2.01	0.43
1:A:112:GLU:O	1:B:112:GLU:HB3	2.19	0.43
1:A:317:THR:O	1:A:318:ASN:HB2	2.18	0.43
1:B:231:ALA:HB1	1:B:252:VAL:CG2	2.48	0.43
1:C:149:LEU:C	1:C:149:LEU:CD2	2.92	0.43
1:D:83:TRP:CE2	1:D:87:LYS:HD2	2.54	0.43
1:D:168:TRP:CD1	1:D:212:LYS:HE3	2.53	0.43
1:A:48:VAL:HG21	1:A:97:TRP:CD2	2.54	0.43
1:C:54:TYR:HA	1:C:305:ILE:HD13	2.00	0.43
1:A:109:ARG:NH2	1:A:352:GLU:OE2	2.52	0.42
1:A:328:LEU:HA	1:A:329:PRO:HD3	1.88	0.42
1:A:48:VAL:HG21	1:A:97:TRP:CG	2.53	0.42
1:A:217:LYS:HB3	1:A:218:PRO:HD3	2.01	0.42
1:B:176:ARG:NH1	1:B:238:GLU:OE2	2.49	0.42
1:C:226:SER:CA	5:C:422:MLT:H31	2.50	0.42
1:D:48[B]:VAL:HG21	1:D:97:TRP:CE2	2.55	0.42
1:D:130:MSE:CE	1:D:293:VAL:HB	2.44	0.42
1:D:176:ARG:HD2	1:D:238:GLU:O	2.20	0.42
1:C:182:HIS:ND1	1:C:183:PRO:HD2	2.34	0.42
1:C:260:SER:OG	4:C:420:ACT:H2	2.20	0.42
1:D:109:ARG:NH1	1:D:352:GLU:HG3	2.35	0.42
1:B:202:TYR:HA	1:B:203:PRO:C	2.45	0.41
1:C:267:ILE:N	1:C:267:ILE:HD12	2.35	0.41
1:B:262:THR:OG1	3:B:408:GOL:H12	2.20	0.41
1:D:289:LEU:HD12	1:D:289:LEU:HA	1.81	0.41
1:A:176:ARG:NH1	1:A:238:GLU:OE2	2.54	0.41
1:D:45:ILE:HG23	1:D:284:THR:HG22	2.03	0.41
1:A:155:THR:C	1:A:157:PRO:HD3	2.46	0.41
1:D:179:LEU:O	1:D:223:TRP:HA	2.21	0.41
1:B:177:ARG:HD2	6:B:537:HOH:O	2.21	0.40
1:B:177:ARG:HA	1:B:241:MSE:O	2.21	0.40
1:C:189:ALA:CB	1:C:243:MSE:HE1	2.51	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	320/324 (99%)	311 (97%)	9 (3%)	0	100	100
1	B	323/324 (100%)	315 (98%)	8 (2%)	0	100	100
1	C	321/324 (99%)	314 (98%)	7 (2%)	0	100	100
1	D	324/324 (100%)	316 (98%)	8 (2%)	0	100	100
All	All	1288/1296 (99%)	1256 (98%)	32 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	258/255 (101%)	253 (98%)	5 (2%)	50	66
1	B	260/255 (102%)	257 (99%)	3 (1%)	63	78
1	C	258/255 (101%)	257 (100%)	1 (0%)	84	92
1	D	263/255 (103%)	259 (98%)	4 (2%)	57	73
All	All	1039/1020 (102%)	1026 (99%)	13 (1%)	61	76

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	GLN
1	A	118	LYS

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Mol	Chain	Res	Type
1	A	289	LEU
1	A	318	ASN
1	A	319	VAL
1	B	177	ARG
1	B	252	VAL
1	B	254	LYS
1	C	362	MSE
1	D	50	GLN
1	D	222	VAL
1	D	254	LYS
1	D	289	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	219	HIS
1	A	269	GLN
1	B	304	HIS
1	C	50	GLN
1	C	75	ASN
1	C	281	ASN
1	C	304	HIS
1	D	44	GLN
1	D	58	GLN
1	D	281	ASN
1	D	300	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

72 ligands are modelled in this entry.

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	GOL	B	409	-	5,5,5	0.34	0	5,5,5	0.39	0
4	ACT	A	411	-	3,3,3	0.84	0	3,3,3	1.25	0
2	SO4	D	406	-	4,4,4	0.25	0	6,6,6	0.09	0
4	ACT	A	415	-	3,3,3	0.79	0	3,3,3	1.23	0
2	SO4	C	407	-	4,4,4	0.25	0	6,6,6	0.15	0
5	MLT	D	417	-	8,8,8	1.15	0	10,10,10	1.63	2 (20%)
4	ACT	C	418	-	3,3,3	0.79	0	3,3,3	1.19	0
2	SO4	D	408	-	4,4,4	0.22	0	6,6,6	0.09	0
4	ACT	C	420	-	3,3,3	0.84	0	3,3,3	1.27	0
2	SO4	A	402	-	4,4,4	0.27	0	6,6,6	0.11	0
4	ACT	C	415	-	3,3,3	0.81	0	3,3,3	1.31	0
3	GOL	B	408	-	5,5,5	0.49	0	5,5,5	0.59	0
3	GOL	C	413	-	5,5,5	0.39	0	5,5,5	0.37	0
5	MLT	A	417	-	8,8,8	1.01	0	10,10,10	1.89	3 (30%)
3	GOL	C	412	-	5,5,5	0.34	0	5,5,5	0.26	0
2	SO4	C	408	-	4,4,4	0.25	0	6,6,6	0.06	0
2	SO4	C	409	-	4,4,4	0.26	0	6,6,6	0.26	0
2	SO4	B	407	-	4,4,4	0.22	0	6,6,6	0.13	0
3	GOL	B	412	-	5,5,5	0.36	0	5,5,5	0.21	0
4	ACT	D	414	-	3,3,3	0.80	0	3,3,3	1.28	0
3	GOL	A	410	-	5,5,5	0.28	0	5,5,5	0.48	0
3	GOL	C	411	-	5,5,5	0.36	0	5,5,5	0.25	0
2	SO4	D	403	-	4,4,4	0.25	0	6,6,6	0.22	0
2	SO4	D	405	-	4,4,4	0.24	0	6,6,6	0.13	0
2	SO4	A	401	-	4,4,4	0.26	0	6,6,6	0.18	0
4	ACT	A	413	-	3,3,3	0.82	0	3,3,3	1.55	0
4	ACT	A	416	-	3,3,3	0.79	0	3,3,3	1.27	0
4	ACT	C	419	-	3,3,3	0.83	0	3,3,3	1.27	0
2	SO4	A	407	-	4,4,4	0.26	0	6,6,6	0.20	0
2	SO4	C	401	-	4,4,4	0.27	0	6,6,6	0.20	0
2	SO4	D	407	-	4,4,4	0.20	0	6,6,6	0.19	0
4	ACT	A	414	-	3,3,3	0.81	0	3,3,3	1.34	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	A	403	-	4,4,4	0.25	0	6,6,6	0.06	0
2	SO4	D	409	-	4,4,4	0.25	0	6,6,6	0.05	0
2	SO4	C	403	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	D	404	-	4,4,4	0.28	0	6,6,6	0.11	0
5	MLT	B	416	-	8,8,8	1.18	0	10,10,10	1.71	1 (10%)
2	SO4	A	404	-	4,4,4	0.26	0	6,6,6	0.13	0
2	SO4	B	402	-	4,4,4	0.31	0	6,6,6	0.26	0
2	SO4	B	403	-	4,4,4	0.25	0	6,6,6	0.21	0
2	SO4	A	409	-	4,4,4	0.25	0	6,6,6	0.09	0
2	SO4	C	406	-	4,4,4	0.25	0	6,6,6	0.07	0
3	GOL	B	410	-	5,5,5	0.30	0	5,5,5	0.38	0
4	ACT	B	413	-	3,3,3	0.84	0	3,3,3	1.43	0
3	GOL	D	413	-	5,5,5	0.46	0	5,5,5	0.14	0
2	SO4	B	404	-	4,4,4	0.26	0	6,6,6	0.15	0
3	GOL	C	410	-	5,5,5	0.39	0	5,5,5	0.28	0
2	SO4	C	402	-	4,4,4	0.22	0	6,6,6	0.13	0
3	GOL	D	412	-	5,5,5	0.37	0	5,5,5	0.37	0
5	MLT	C	422	-	8,8,8	1.25	1 (12%)	10,10,10	1.62	3 (30%)
3	GOL	C	414	-	5,5,5	0.38	0	5,5,5	0.34	0
2	SO4	C	404	-	4,4,4	0.25	0	6,6,6	0.11	0
4	ACT	C	421	-	3,3,3	0.83	0	3,3,3	1.32	0
4	ACT	C	416	-	3,3,3	0.81	0	3,3,3	1.26	0
4	ACT	A	412	-	3,3,3	0.84	0	3,3,3	1.41	0
2	SO4	A	408	-	4,4,4	0.42	0	6,6,6	0.08	0
4	ACT	C	417	-	3,3,3	0.80	0	3,3,3	1.29	0
4	ACT	B	414	-	3,3,3	0.83	0	3,3,3	1.27	0
2	SO4	D	402	-	4,4,4	0.24	0	6,6,6	0.20	0
3	GOL	B	411	-	5,5,5	0.39	0	5,5,5	0.38	0
4	ACT	D	415	-	3,3,3	0.75	0	3,3,3	1.40	0
2	SO4	D	401	-	4,4,4	0.24	0	6,6,6	0.31	0
2	SO4	A	405	-	4,4,4	0.25	0	6,6,6	0.13	0
4	ACT	B	415	-	3,3,3	0.84	0	3,3,3	1.36	0
4	ACT	D	416	-	3,3,3	0.81	0	3,3,3	1.30	0
3	GOL	D	411	-	5,5,5	0.33	0	5,5,5	0.44	0
2	SO4	B	401	-	4,4,4	0.30	0	6,6,6	0.18	0
2	SO4	B	405	-	4,4,4	0.24	0	6,6,6	0.12	0
2	SO4	C	405	-	4,4,4	0.25	0	6,6,6	0.11	0
2	SO4	A	406	-	4,4,4	0.23	0	6,6,6	0.12	0
2	SO4	B	406	-	4,4,4	0.23	0	6,6,6	0.07	0
3	GOL	D	410	-	5,5,5	0.31	0	5,5,5	0.48	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	409	-	-	2/4/4/4	-
5	MLT	D	417	-	-	4/8/8/8	-
3	GOL	C	413	-	-	2/4/4/4	-
5	MLT	A	417	-	-	5/8/8/8	-
3	GOL	B	408	-	-	2/4/4/4	-
3	GOL	C	412	-	-	2/4/4/4	-
3	GOL	B	412	-	-	0/4/4/4	-
3	GOL	A	410	-	-	2/4/4/4	-
3	GOL	C	411	-	-	0/4/4/4	-
5	MLT	B	416	-	-	8/8/8/8	-
3	GOL	B	410	-	-	2/4/4/4	-
3	GOL	D	413	-	-	2/4/4/4	-
3	GOL	C	410	-	-	2/4/4/4	-
5	MLT	C	422	-	-	4/8/8/8	-
3	GOL	D	412	-	-	2/4/4/4	-
3	GOL	C	414	-	-	1/4/4/4	-
3	GOL	B	411	-	-	2/4/4/4	-
3	GOL	D	411	-	-	2/4/4/4	-
3	GOL	D	410	-	-	3/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	422	MLT	C2-C1	-2.19	1.49	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	416	MLT	O2-C1-C2	3.88	120.94	112.74
5	A	417	MLT	O2-C1-C2	3.72	120.61	112.74
5	D	417	MLT	O2-C1-C2	3.46	120.06	112.74
5	C	422	MLT	O2-C1-C2	2.94	118.95	112.74
5	A	417	MLT	O2-C1-O1	-2.36	118.73	124.08
5	C	422	MLT	O5-C4-C3	2.27	121.06	114.00
5	A	417	MLT	O5-C4-C3	2.24	120.98	114.00
5	D	417	MLT	O5-C4-O4	-2.15	117.81	123.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	422	MLT	C3-C2-C1	-2.08	105.42	110.53

There are no chirality outliers.

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	410	GOL	O1-C1-C2-C3
3	B	409	GOL	O1-C1-C2-C3
3	B	410	GOL	O1-C1-C2-C3
3	B	411	GOL	O1-C1-C2-O2
3	B	411	GOL	O1-C1-C2-C3
3	C	413	GOL	O1-C1-C2-C3
3	D	410	GOL	O1-C1-C2-C3
3	D	411	GOL	O1-C1-C2-O2
3	D	412	GOL	O1-C1-C2-C3
3	D	413	GOL	O1-C1-C2-O2
3	D	413	GOL	O1-C1-C2-C3
5	B	416	MLT	O1-C1-C2-O3
5	B	416	MLT	O2-C1-C2-O3
5	C	422	MLT	C1-C2-C3-C4
5	C	422	MLT	O3-C2-C3-C4
3	D	412	GOL	O1-C1-C2-O2
5	B	416	MLT	O3-C2-C3-C4
5	B	416	MLT	C1-C2-C3-C4
3	C	410	GOL	C1-C2-C3-O3
3	C	412	GOL	O1-C1-C2-C3
3	D	411	GOL	O1-C1-C2-C3
3	A	410	GOL	O1-C1-C2-O2
3	B	409	GOL	O1-C1-C2-O2
3	B	410	GOL	O1-C1-C2-O2
3	C	412	GOL	O1-C1-C2-O2
3	D	410	GOL	O1-C1-C2-O2
3	C	413	GOL	O1-C1-C2-O2
5	D	417	MLT	O2-C1-C2-O3
5	B	416	MLT	O1-C1-C2-C3
5	B	416	MLT	O2-C1-C2-C3
3	C	414	GOL	O1-C1-C2-O2
3	B	408	GOL	O1-C1-C2-O2
5	D	417	MLT	O1-C1-C2-O3
5	B	416	MLT	C2-C3-C4-O5
5	A	417	MLT	O2-C1-C2-C3
5	C	422	MLT	O2-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
5	D	417	MLT	O3-C2-C3-C4
5	B	416	MLT	C2-C3-C4-O4
3	B	408	GOL	O1-C1-C2-C3
3	C	410	GOL	O2-C2-C3-O3
5	A	417	MLT	C2-C3-C4-O5
5	D	417	MLT	C1-C2-C3-C4
5	A	417	MLT	O1-C1-C2-C3
5	C	422	MLT	O1-C1-C2-C3
3	D	410	GOL	O2-C2-C3-O3
5	A	417	MLT	C2-C3-C4-O4
5	A	417	MLT	O2-C1-C2-O3

There are no ring outliers.

18 monomers are involved in 43 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	409	GOL	2	0
5	D	417	MLT	4	0
4	C	420	ACT	1	0
3	B	408	GOL	3	0
5	A	417	MLT	3	0
3	C	412	GOL	5	0
3	B	412	GOL	2	0
4	A	413	ACT	2	0
2	D	407	SO4	1	0
5	B	416	MLT	3	0
2	A	409	SO4	1	0
3	D	413	GOL	4	0
3	C	410	GOL	3	0
5	C	422	MLT	5	0
4	A	412	ACT	1	0
2	A	408	SO4	1	0
4	D	416	ACT	1	0
3	D	410	GOL	1	0

5.7 Other polymers ⓘ

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 [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	316/324 (97%)	-0.19	4 (1%) 75 73	42, 56, 82, 101	0
1	B	316/324 (97%)	-0.34	2 (0%) 85 83	27, 49, 73, 100	3 (0%)
1	C	316/324 (97%)	-0.30	2 (0%) 85 83	33, 51, 75, 125	1 (0%)
1	D	316/324 (97%)	-0.38	2 (0%) 85 83	25, 48, 73, 98	4 (1%)
All	All	1264/1296 (97%)	-0.30	10 (0%) 82 80	25, 51, 77, 125	8 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	364	LYS	5.3
1	B	83	TRP	3.9
1	A	83	TRP	3.7
1	B	363	GLN	3.0
1	C	43	ALA	2.8
1	D	363	GLN	2.6
1	A	158	LYS	2.5
1	A	363	GLN	2.2
1	A	283	GLU	2.1
1	D	254	LYS	2.1

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 ⓘ

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	SO4	B	407	5/5	0.52	0.15	140,141,144,146	5
2	SO4	C	409	5/5	0.54	0.13	121,124,126,126	5
2	SO4	C	408	5/5	0.56	0.11	134,135,139,142	0
2	SO4	A	405	5/5	0.60	0.11	136,136,138,139	0
4	ACT	C	421	4/4	0.62	0.24	91,93,95,96	0
2	SO4	A	408	5/5	0.65	0.10	151,151,152,153	0
2	SO4	A	407	5/5	0.66	0.14	113,121,123,126	0
2	SO4	A	404	5/5	0.68	0.14	115,116,120,126	0
3	GOL	B	412	6/6	0.69	0.23	100,108,109,110	0
4	ACT	C	420	4/4	0.71	0.21	93,95,98,100	0
4	ACT	C	416	4/4	0.71	0.26	95,98,99,100	0
2	SO4	C	407	5/5	0.73	0.17	106,107,110,118	5
2	SO4	D	409	5/5	0.73	0.11	139,142,143,143	0
2	SO4	B	404	5/5	0.76	0.11	126,127,129,129	0
2	SO4	A	403	5/5	0.76	0.11	130,132,137,137	0
4	ACT	B	415	4/4	0.77	0.29	105,111,111,112	0
3	GOL	B	408	6/6	0.77	0.22	56,78,80,84	0
4	ACT	A	413	4/4	0.77	0.27	84,84,91,95	0
4	ACT	A	415	4/4	0.77	0.31	84,93,94,95	0
4	ACT	A	414	4/4	0.78	0.33	93,95,100,101	0
2	SO4	A	409	5/5	0.78	0.10	106,107,110,110	5
2	SO4	D	407	5/5	0.78	0.17	93,98,105,113	0
4	ACT	D	414	4/4	0.78	0.33	86,91,93,96	0
2	SO4	A	402	5/5	0.80	0.11	92,93,108,108	0
2	SO4	B	402	5/5	0.80	0.16	86,96,102,104	0
2	SO4	C	401	5/5	0.80	0.18	94,107,113,116	0
4	ACT	D	416	4/4	0.80	0.29	90,95,96,96	0
2	SO4	C	406	5/5	0.81	0.13	130,132,135,137	0
3	GOL	D	410	6/6	0.81	0.17	53,81,88,92	0
4	ACT	A	412	4/4	0.81	0.23	71,79,80,89	0
2	SO4	D	404	5/5	0.82	0.10	91,91,100,105	5
2	SO4	B	403	5/5	0.83	0.12	104,110,113,113	0
2	SO4	C	403	5/5	0.83	0.13	101,104,107,107	0
3	GOL	C	411	6/6	0.84	0.19	91,103,105,107	0
4	ACT	B	414	4/4	0.84	0.21	79,85,86,87	0
2	SO4	B	406	5/5	0.84	0.07	138,138,140,141	0
3	GOL	D	413	6/6	0.84	0.22	59,76,79,82	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MLT	A	417	9/9	0.84	0.14	68,88,95,101	0
5	MLT	B	416	9/9	0.84	0.13	58,73,85,90	0
5	MLT	D	417	9/9	0.84	0.14	66,78,86,89	0
5	MLT	C	422	9/9	0.85	0.16	71,75,86,92	0
2	SO4	A	406	5/5	0.85	0.12	123,124,125,126	0
2	SO4	D	402	5/5	0.86	0.16	71,80,87,100	5
3	GOL	D	412	6/6	0.86	0.25	77,89,92,92	0
3	GOL	B	411	6/6	0.86	0.15	60,69,74,76	0
2	SO4	D	405	5/5	0.87	0.09	83,85,96,97	5
2	SO4	C	402	5/5	0.87	0.10	91,99,101,104	0
2	SO4	D	408	5/5	0.87	0.18	116,118,121,126	5
4	ACT	B	413	4/4	0.87	0.18	72,77,79,83	0
2	SO4	D	401	5/5	0.87	0.16	85,89,99,100	0
2	SO4	C	405	5/5	0.88	0.09	100,101,106,113	5
2	SO4	D	406	5/5	0.88	0.16	142,143,145,147	0
4	ACT	A	411	4/4	0.88	0.24	80,81,83,86	0
3	GOL	C	410	6/6	0.88	0.14	60,79,81,82	0
2	SO4	B	405	5/5	0.88	0.14	123,125,127,130	0
4	ACT	C	419	4/4	0.88	0.21	87,89,92,94	0
3	GOL	B	409	6/6	0.88	0.14	70,74,77,80	0
3	GOL	B	410	6/6	0.89	0.19	82,85,87,92	0
2	SO4	D	403	5/5	0.89	0.10	90,90,98,99	5
4	ACT	C	417	4/4	0.89	0.17	78,86,87,88	0
3	GOL	A	410	6/6	0.89	0.19	68,71,79,83	0
2	SO4	A	401	5/5	0.89	0.16	102,103,110,111	0
2	SO4	B	401	5/5	0.89	0.12	59,59,71,81	5
3	GOL	D	411	6/6	0.90	0.18	77,79,80,82	0
3	GOL	C	414	6/6	0.90	0.18	90,91,93,94	0
4	ACT	C	418	4/4	0.92	0.21	73,75,81,84	0
4	ACT	A	416	4/4	0.92	0.14	86,92,93,97	0
4	ACT	D	415	4/4	0.92	0.18	77,79,81,82	0
3	GOL	C	412	6/6	0.92	0.16	58,69,80,92	0
4	ACT	C	415	4/4	0.93	0.18	92,98,99,100	0
2	SO4	C	404	5/5	0.95	0.13	115,116,121,127	5
3	GOL	C	413	6/6	0.95	0.19	88,92,95,101	0

6.5 Other polymers

There are no such residues in this entry.