



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:48 AM UTC

PDB ID : 6EUO / pdb_00006euo
Title : Crystal structure of APO Fe(II)/alpha-ketoglutarate dependent dioxygenase KDO5
Authors : Isabet, T.; Stura, E.; Legrand, P.; Zaparucha, A.; Bastard, K.
Deposited on : 2017-10-30
Resolution : 2.30 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

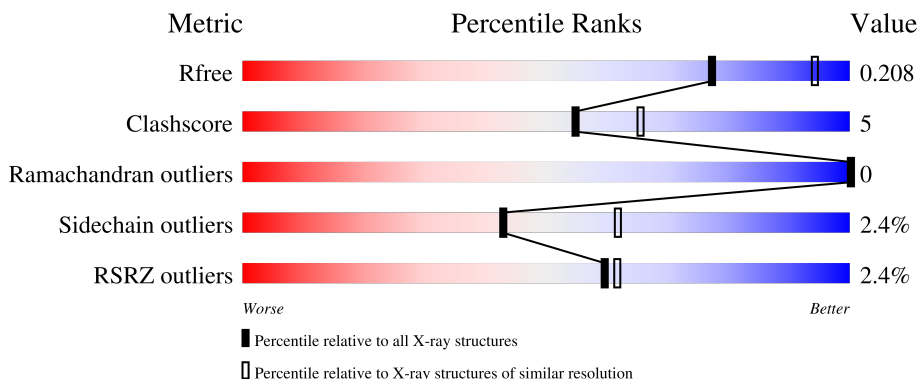
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	372	% 82% 11% 7%
1	B	372	5% 77% 12% 10%
1	C	372	3% 83% 12% 5%
1	D	372	% 84% 9% 8%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11693 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.

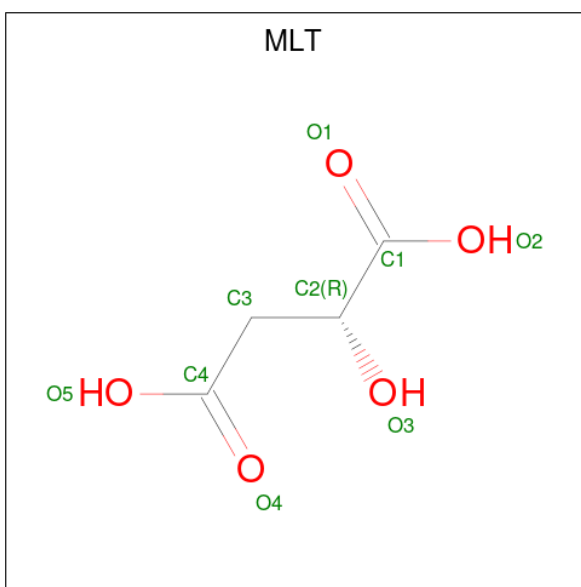
- Molecule 1 is a protein called L-lysine 4-hydroxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	346	Total	C	N	O	S	0	8	0
			2814	1787	482	532	13			
1	B	335	Total	C	N	O	S	0	8	0
			2706	1719	459	513	15			
1	C	354	Total	C	N	O	S	0	3	0
			2817	1787	482	536	12			
1	D	344	Total	C	N	O	S	0	1	0
			2734	1738	466	518	12			

- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe).

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

- Molecule 3 is D-MALATE (CCD ID: MLT) (formula: C₄H₆O₅).



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

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

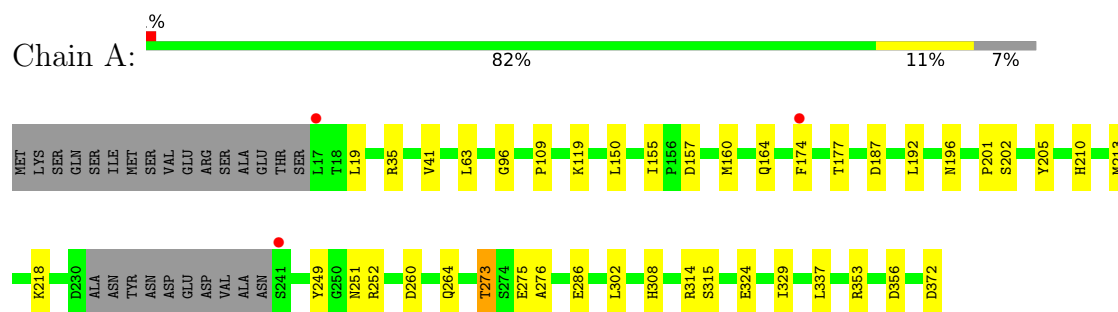
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	172	Total 172	O 172	0	0
6	B	105	Total 105	O 105	0	0
6	C	117	Total 117	O 117	0	0
6	D	131	Total 131	O 131	0	0

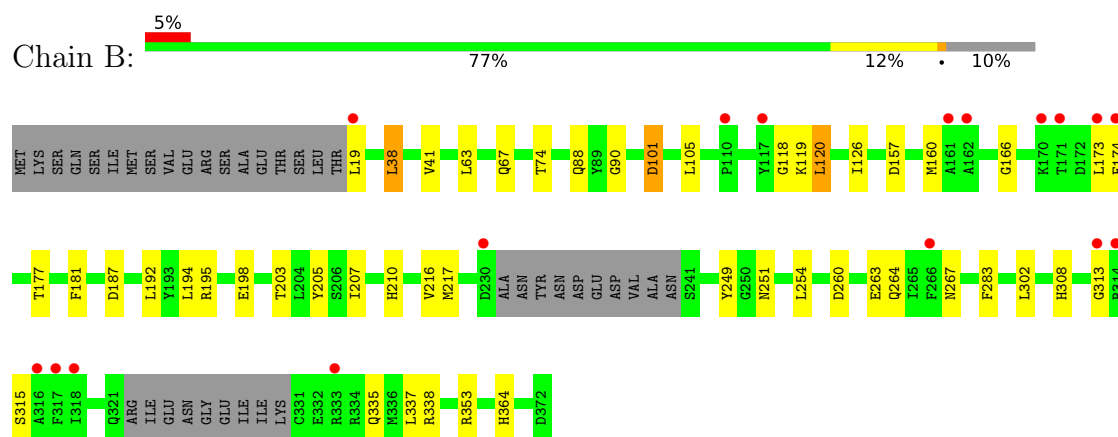
3 Residue-property plots [i](#)

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.

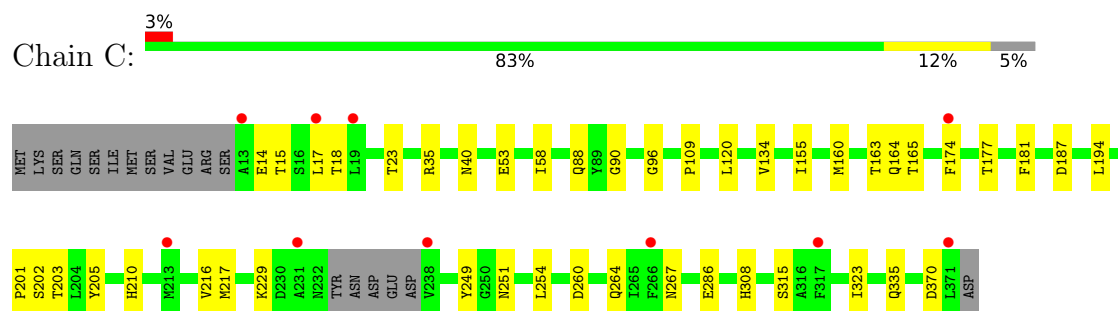
- Molecule 1: L-lysine 4-hydroxylase



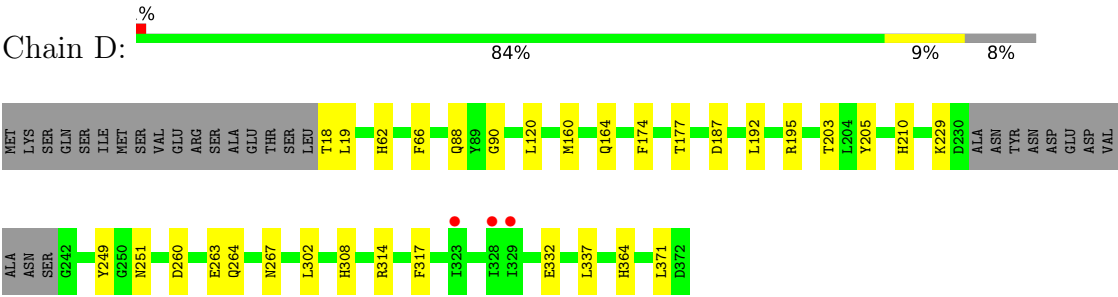
- Molecule 1: L-lysine 4-hydroxylase



- Molecule 1: L-lysine 4-hydroxylase



- Molecule 1: L-lysine 4-hydroxylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	91.28Å 98.93Å 166.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	62.20 – 2.30 62.20 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (62.20-2.30) 100.0 (62.20-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 2.29Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.169 , 0.205 0.173 , 0.208	Depositor DCC
R_{free} test set	3359 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	49.4	Xtriage
Anisotropy	0.505	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 75.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	11693	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.66% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FE, SO4, GOL, MLT

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	1/2878 (0.0%)	1.19	2/3892 (0.1%)
1	B	0.81	0/2768	1.23	6/3747 (0.2%)
1	C	0.80	1/2879 (0.0%)	1.25	7/3898 (0.2%)
1	D	0.83	0/2796	1.23	8/3784 (0.2%)
All	All	0.83	2/11321 (0.0%)	1.22	23/15321 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	134	VAL	CA-C	5.64	1.57	1.52
1	A	196	ASN	CA-CB	5.03	1.57	1.53

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	101	ASP	CA-CB-CG	7.35	119.95	112.60
1	C	14	GLU	CA-C-N	6.89	130.44	120.38
1	C	14	GLU	C-N-CA	6.89	130.44	120.38
1	B	267	ASN	CA-CB-CG	6.51	119.11	112.60
1	D	177	THR	N-CA-C	-6.35	100.22	110.32
1	A	177	THR	N-CA-C	-6.16	100.53	110.32
1	C	177	THR	N-CA-C	-6.16	100.53	110.32
1	C	323	ILE	N-CA-CB	6.08	118.33	111.21
1	D	18	THR	CA-C-N	5.83	132.67	121.54
1	D	18	THR	C-N-CA	5.83	132.67	121.54
1	B	120	LEU	N-CA-C	-5.75	105.15	111.82
1	A	157	ASP	CA-CB-CG	5.42	118.02	112.60
1	D	267	ASN	CA-CB-CG	5.29	117.89	112.60
1	C	267	ASN	CA-CB-CG	5.28	117.88	112.60
1	B	41	VAL	CA-C-N	5.22	125.73	119.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	41	VAL	C-N-CA	5.22	125.73	119.94
1	B	177	THR	N-CA-C	-5.18	102.09	110.32
1	C	120	LEU	CA-C-N	5.17	127.52	120.54
1	C	120	LEU	C-N-CA	5.17	127.52	120.54
1	D	120	LEU	CA-C-N	5.16	127.15	120.44
1	D	120	LEU	C-N-CA	5.16	127.15	120.44
1	D	160	MET	CA-C-N	5.16	127.92	120.38
1	D	160	MET	C-N-CA	5.16	127.92	120.38

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2814	0	2744	33	0
1	B	2706	0	2622	29	0
1	C	2817	0	2757	23	0
1	D	2734	0	2671	22	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	9	0	4	0	0
3	B	9	0	4	2	0
3	C	9	0	4	0	0
3	D	9	0	4	0	0
4	A	12	0	16	4	0
4	B	12	0	16	1	0
4	C	18	0	24	0	0
5	A	5	0	0	0	0
5	C	5	0	0	0	0
5	D	5	0	0	0	0
6	A	172	0	0	1	0
6	B	105	0	0	0	0
6	C	117	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	D	131	0	0	0	0
All	All	11693	0	10866	105	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 5.

All (105) 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:88:GLN:HE21	1:B:90:GLY:H	1.23	0.85
1:C:251:ASN:HD22	1:C:254:LEU:H	1.25	0.84
1:A:273:THR:HG22	1:A:276:ALA:H	1.40	0.84
1:D:88:GLN:HE21	1:D:90:GLY:H	1.25	0.84
1:B:251:ASN:HD22	1:B:254:LEU:H	1.26	0.84
1:C:88:GLN:HE21	1:C:90:GLY:H	1.24	0.83
1:B:67:GLN:HB2	1:C:40:ASN:HD22	1.55	0.71
1:B:166:GLY:H	3:B:402:MLT:H32	1.55	0.71
1:A:353:ARG:HA	4:A:403:GOL:H32	1.74	0.68
1:B:353:ARG:HA	4:B:404:GOL:H2	1.76	0.68
1:A:260:ASP:H	1:A:264:GLN:HE21	1.44	0.66
1:D:249:TYR:OH	1:D:308:HIS:HD2	1.79	0.65
1:D:260:ASP:H	1:D:264:GLN:HE21	1.45	0.65
1:C:260:ASP:H	1:C:264:GLN:HE21	1.45	0.64
1:B:249:TYR:OH	1:B:308:HIS:HD2	1.82	0.63
1:A:249:TYR:OH	1:A:308:HIS:HD2	1.81	0.63
1:B:260:ASP:H	1:B:264:GLN:HE21	1.45	0.63
1:C:249:TYR:OH	1:C:308:HIS:HD2	1.82	0.62
1:D:205:TYR:OH	1:D:210:HIS:HE1	1.84	0.60
1:A:205:TYR:OH	1:A:210:HIS:HE1	1.85	0.60
1:A:273:THR:CG2	1:A:276:ALA:H	2.13	0.58
1:B:205:TYR:OH	1:B:210:HIS:HE1	1.87	0.57
1:A:356:ASP:OD1	4:A:404:GOL:H12	2.05	0.57
1:C:205:TYR:OH	1:C:210:HIS:HE1	1.87	0.56
1:C:88:GLN:HE21	1:C:90:GLY:N	2.01	0.56
1:A:353:ARG:HG2	4:A:403:GOL:H11	1.88	0.55
1:B:216:VAL:HG12	1:B:217[B]:MET:HE2	1.90	0.54
1:B:88:GLN:HE21	1:B:90:GLY:N	2.01	0.54
1:A:205:TYR:OH	1:A:210:HIS:CE1	2.61	0.53
1:B:88:GLN:NE2	1:B:90:GLY:H	2.02	0.53
1:B:217[B]:MET:HE1	1:B:283:PHE:HD1	1.71	0.53
1:A:164:GLN:HE22	1:A:174[B]:PHE:H	1.57	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:PRO:HG2	1:A:155:ILE:HD13	1.91	0.52
1:D:88:GLN:HE21	1:D:90:GLY:N	2.00	0.52
1:D:192:LEU:HD23	1:D:337:LEU:HD12	1.93	0.51
1:C:205:TYR:OH	1:C:210:HIS:CE1	2.63	0.51
1:D:205:TYR:OH	1:D:210:HIS:CE1	2.63	0.51
1:B:205:TYR:OH	1:B:210:HIS:CE1	2.64	0.51
1:C:109:PRO:HG2	1:C:155:ILE:HD13	1.93	0.51
1:C:88:GLN:NE2	1:C:90:GLY:H	2.02	0.51
1:B:338:ARG:HH12	3:B:402:MLT:H2	1.76	0.50
1:B:38:LEU:HD22	1:B:126:ILE:HG12	1.93	0.50
1:A:252:ARG:HD2	6:A:538:HOH:O	2.12	0.50
1:C:163:THR:OG1	1:C:165:THR:HG22	2.11	0.50
1:D:88:GLN:NE2	1:D:90:GLY:H	2.02	0.50
1:B:249:TYR:OH	1:B:308:HIS:CD2	2.65	0.49
1:A:324:GLU:HB3	1:A:329:ILE:CD1	2.42	0.49
1:D:249:TYR:OH	1:D:308:HIS:CD2	2.63	0.49
1:C:164:GLN:HE22	1:C:174:PHE:H	1.59	0.49
1:D:164:GLN:HE22	1:D:174:PHE:H	1.59	0.49
1:C:249:TYR:OH	1:C:308:HIS:CD2	2.65	0.48
1:A:192:LEU:HD23	1:A:337:LEU:HD12	1.95	0.48
1:A:249:TYR:OH	1:A:308:HIS:CD2	2.64	0.48
1:A:314:ARG:HH11	1:A:314:ARG:HG2	1.78	0.48
1:B:63:LEU:O	1:B:67:GLN:HG2	2.14	0.48
1:C:251:ASN:ND2	1:C:254:LEU:H	2.03	0.47
1:A:164:GLN:HE22	1:A:174[A]:PHE:H	1.60	0.47
1:D:195:ARG:HD3	1:D:332:GLU:HB3	1.95	0.47
1:B:173:LEU:HD23	1:B:313:GLY:HA2	1.96	0.47
1:C:53:GLU:HA	1:C:58:ILE:HD11	1.95	0.47
1:B:192:LEU:HD23	1:B:337:LEU:HD12	1.97	0.46
1:A:187:ASP:OD1	1:A:308:HIS:HE1	1.98	0.46
1:A:210:HIS:HD2	1:A:286:GLU:OE2	1.99	0.46
1:D:187:ASP:OD1	1:D:308:HIS:HE1	2.00	0.45
1:C:201:PRO:HD2	1:C:315:SER:HB2	2.00	0.44
1:C:187:ASP:OD1	1:C:308:HIS:HE1	1.99	0.44
1:A:213[A]:MET:HG3	1:A:218:LYS:HE2	1.99	0.44
1:A:164:GLN:NE2	1:A:174[B]:PHE:HD1	2.15	0.44
1:B:187:ASP:OD1	1:B:308:HIS:HE1	2.00	0.44
1:B:251:ASN:ND2	1:B:254:LEU:H	2.05	0.44
1:C:217:MET:HE3	1:C:286:GLU:OE2	2.17	0.44
1:D:19:LEU:HD12	1:D:19:LEU:H	1.83	0.44
1:C:216:VAL:HG12	1:C:217:MET:HE2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:LEU:HD21	4:A:404:GOL:H32	2.01	0.43
1:B:194:LEU:HD12	1:B:335:GLN:HG2	1.99	0.43
1:D:314:ARG:HD2	1:D:317:PHE:HE1	1.84	0.43
1:A:164:GLN:HE22	1:A:174[B]:PHE:HD1	1.67	0.43
1:B:118:GLY:C	1:B:120:LEU:H	2.26	0.43
1:C:260:ASP:H	1:C:264:GLN:NE2	2.15	0.42
1:A:192:LEU:HB2	1:A:302:LEU:HD12	2.02	0.42
1:A:201:PRO:HD2	1:A:315:SER:HB2	2.02	0.42
1:C:194:LEU:HD12	1:C:335:GLN:HG2	2.01	0.41
1:D:229:LYS:HD3	1:D:364:HIS:HB2	2.01	0.41
1:B:157:ASP:HB3	1:B:160:MET:HB2	2.02	0.41
1:D:164:GLN:NE2	1:D:174:PHE:H	2.19	0.41
1:A:35:ARG:HD3	1:A:96:GLY:O	2.20	0.41
1:B:207:ILE:HD11	1:B:217[A]:MET:HE3	2.03	0.41
1:D:192:LEU:HB2	1:D:302:LEU:HD12	2.02	0.41
1:C:181:PHE:CG	1:C:229:LYS:HA	2.56	0.41
1:D:62:HIS:O	1:D:66:PHE:HD2	2.04	0.41
1:A:260:ASP:H	1:A:264:GLN:NE2	2.14	0.41
1:B:105:LEU:HD12	1:B:195:ARG:CZ	2.51	0.41
1:C:35:ARG:HD3	1:C:96:GLY:O	2.21	0.41
1:B:192:LEU:HB2	1:B:302:LEU:HD12	2.03	0.40
1:B:181:PHE:HB2	1:B:364:HIS:HD2	1.87	0.40
1:A:63:LEU:HD13	1:B:74:THR:HG21	2.03	0.40
1:A:324:GLU:HB3	1:A:329:ILE:HD11	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	350/372 (94%)	345 (99%)	5 (1%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	337/372 (91%)	329 (98%)	8 (2%)	0	100	100
1	C	353/372 (95%)	342 (97%)	11 (3%)	0	100	100
1	D	341/372 (92%)	331 (97%)	10 (3%)	0	100	100
All	All	1381/1488 (93%)	1347 (98%)	34 (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	307/322 (95%)	299 (97%)	8 (3%)	40	59
1	B	295/322 (92%)	286 (97%)	9 (3%)	35	52
1	C	307/322 (95%)	299 (97%)	8 (3%)	40	59
1	D	297/322 (92%)	294 (99%)	3 (1%)	68	82
All	All	1206/1288 (94%)	1178 (98%)	28 (2%)	43	63

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	41	VAL
1	A	119	LYS
1	A	160	MET
1	A	202	SER
1	A	273	THR
1	A	275	GLU
1	A	372	ASP
1	B	19	LEU
1	B	38	LEU
1	B	101	ASP
1	B	119	LYS
1	B	174	PHE

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Mol	Chain	Res	Type
1	B	198	GLU
1	B	203	THR
1	B	263	GLU
1	B	315	SER
1	C	15	THR
1	C	17	LEU
1	C	18	THR
1	C	23	THR
1	C	160	MET
1	C	202	SER
1	C	203	THR
1	C	370	ASP
1	D	203	THR
1	D	263	GLU
1	D	371	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (46) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	ASN
1	A	67	GLN
1	A	79	GLN
1	A	164	GLN
1	A	184	ASN
1	A	210	HIS
1	A	226	GLN
1	A	251	ASN
1	A	264	GLN
1	A	267	ASN
1	A	308	HIS
1	B	79	GLN
1	B	88	GLN
1	B	184	ASN
1	B	210	HIS
1	B	226	GLN
1	B	251	ASN
1	B	264	GLN
1	B	298	ASN
1	B	308	HIS
1	C	40	ASN
1	C	43	ASN
1	C	88	GLN

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Mol	Chain	Res	Type
1	C	113	GLN
1	C	164	GLN
1	C	184	ASN
1	C	210	HIS
1	C	226	GLN
1	C	251	ASN
1	C	264	GLN
1	C	267	ASN
1	C	298	ASN
1	C	308	HIS
1	C	347	HIS
1	D	40	ASN
1	D	43	ASN
1	D	51	ASN
1	D	79	GLN
1	D	88	GLN
1	D	164	GLN
1	D	184	ASN
1	D	210	HIS
1	D	226	GLN
1	D	251	ASN
1	D	264	GLN
1	D	308	HIS

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 18 ligands modelled in this entry, 4 are monoatomic - leaving 14 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$
4	GOL	B	404	-	5,5,5	0.08	0	5,5,5	0.16	0
5	SO4	D	403	-	4,4,4	0.28	0	6,6,6	0.20	0
3	MLT	B	402	-	8,8,8	1.05	0	10,10,10	1.49	2 (20%)
3	MLT	A	402	2	8,8,8	1.17	0	10,10,10	1.54	1 (10%)
3	MLT	D	402	2	8,8,8	1.08	0	10,10,10	1.59	3 (30%)
4	GOL	A	404	-	5,5,5	0.07	0	5,5,5	0.33	0
4	GOL	A	403	-	5,5,5	0.11	0	5,5,5	0.29	0
5	SO4	A	405	-	4,4,4	0.47	0	6,6,6	0.34	0
4	GOL	B	403	-	5,5,5	0.09	0	5,5,5	0.14	0
4	GOL	C	404	-	5,5,5	0.05	0	5,5,5	0.15	0
4	GOL	C	403	-	5,5,5	0.10	0	5,5,5	0.42	0
4	GOL	C	405	-	5,5,5	0.07	0	5,5,5	0.31	0
3	MLT	C	402	-	8,8,8	1.14	0	10,10,10	1.51	3 (30%)
5	SO4	C	406	-	4,4,4	0.37	0	6,6,6	0.17	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
4	GOL	B	404	-	-	0/4/4/4	-
3	MLT	B	402	-	-	4/8/8/8	-
3	MLT	A	402	2	-	3/8/8/8	-
3	MLT	D	402	2	-	4/8/8/8	-
4	GOL	A	404	-	-	0/4/4/4	-
4	GOL	A	403	-	-	0/4/4/4	-
4	GOL	B	403	-	-	0/4/4/4	-
4	GOL	C	404	-	-	0/4/4/4	-
4	GOL	C	403	-	-	2/4/4/4	-
4	GOL	C	405	-	-	0/4/4/4	-
3	MLT	C	402	-	-	6/8/8/8	-

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	MLT	O2-C1-C2	3.33	119.79	112.74
3	B	402	MLT	O2-C1-C2	3.19	119.49	112.74
3	D	402	MLT	O5-C4-C3	2.59	122.05	114.00
3	C	402	MLT	O2-C1-C2	2.38	117.78	112.74
3	D	402	MLT	O4-C4-C3	-2.18	116.14	122.84
3	D	402	MLT	O2-C1-C2	2.14	117.27	112.74
3	C	402	MLT	O3-C2-C1	-2.12	104.98	110.36
3	B	402	MLT	O1-C1-C2	-2.10	118.40	122.60
3	C	402	MLT	O5-C4-C3	2.04	120.34	114.00

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	402	MLT	C1-C2-C3-C4
3	B	402	MLT	O3-C2-C3-C4
3	C	402	MLT	O1-C1-C2-O3
3	C	402	MLT	O2-C1-C2-O3
3	C	402	MLT	O2-C1-C2-C3
3	D	402	MLT	O1-C1-C2-O3
3	D	402	MLT	O2-C1-C2-O3
4	C	403	GOL	O1-C1-C2-C3
3	B	402	MLT	O1-C1-C2-O3
3	C	402	MLT	O1-C1-C2-C3
4	C	403	GOL	O1-C1-C2-O2
3	C	402	MLT	O3-C2-C3-C4
3	B	402	MLT	O2-C1-C2-O3
3	D	402	MLT	O2-C1-C2-C3
3	C	402	MLT	C1-C2-C3-C4
3	A	402	MLT	O2-C1-C2-O3
3	A	402	MLT	O2-C1-C2-C3
3	D	402	MLT	O1-C1-C2-C3
3	A	402	MLT	O1-C1-C2-O3

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	404	GOL	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	MLT	2	0
4	A	404	GOL	2	0
4	A	403	GOL	2	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	346/372 (93%)	-0.31	3 (0%) 81 82	24, 52, 89, 131	8 (2%)
1	B	335/372 (90%)	0.08	17 (5%) 33 35	23, 65, 140, 177	8 (2%)
1	C	354/372 (95%)	0.03	10 (2%) 55 57	26, 71, 132, 170	3 (0%)
1	D	344/372 (92%)	-0.10	3 (0%) 81 82	29, 65, 130, 190	1 (0%)
All	All	1379/1488 (92%)	-0.08	33 (2%) 59 62	23, 62, 130, 190	20 (1%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	19	LEU	4.3
1	D	323	ILE	4.1
1	A	174[A]	PHE	3.1
1	B	162	ALA	3.1
1	C	266	PHE	3.1
1	B	173	LEU	3.0
1	B	117	TYR	3.0
1	C	17	LEU	2.9
1	B	313	GLY	2.9
1	B	174	PHE	2.9
1	B	316	ALA	2.8
1	B	318	ILE	2.8
1	B	333	ARG	2.6
1	C	317	PHE	2.6
1	C	213	MET	2.5
1	B	170	LYS	2.4
1	B	110	PRO	2.3
1	B	161	ALA	2.3
1	B	314	ARG	2.3
1	B	230	ASP	2.3
1	C	13	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	371	LEU	2.3
1	B	19	LEU	2.2
1	C	231	ALA	2.2
1	D	328	ILE	2.2
1	C	174	PHE	2.1
1	B	266	PHE	2.1
1	B	171	THR	2.1
1	A	17	LEU	2.1
1	D	329	ILE	2.1
1	C	238	VAL	2.0
1	B	317	PHE	2.0
1	A	241	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	C	405	6/6	0.54	0.15	134,136,137,137	0
3	MLT	B	402	9/9	0.76	0.13	123,125,128,128	0
4	GOL	C	404	6/6	0.81	0.09	89,91,93,95	0
4	GOL	B	404	6/6	0.81	0.12	92,94,95,96	0
4	GOL	C	403	6/6	0.83	0.17	69,80,83,83	0
3	MLT	D	402	9/9	0.85	0.16	90,92,98,101	0
4	GOL	A	403	6/6	0.85	0.14	81,89,91,93	0
4	GOL	B	403	6/6	0.87	0.12	81,86,87,90	0
5	SO4	C	406	5/5	0.87	0.10	130,130,131,131	0
4	GOL	A	404	6/6	0.89	0.13	68,81,86,88	0
5	SO4	A	405	5/5	0.91	0.07	80,80,83,85	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MLT	A	402	9/9	0.92	0.12	72,79,83,88	0
5	SO4	D	403	5/5	0.93	0.08	108,110,111,112	0
3	MLT	C	402	9/9	0.94	0.13	109,112,114,115	0
2	FE	D	401	1/1	0.98	0.06	65,65,65,65	0
2	FE	B	401	1/1	0.99	0.08	60,60,60,60	0
2	FE	C	401	1/1	1.00	0.05	68,68,68,68	0
2	FE	A	401	1/1	1.00	0.03	50,50,50,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.