



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 03:09 AM UTC

PDB ID : 1LUR / pdb_00001lur
Title : Crystal Structure of the GalM/aldose Epimerase Homologue from *C. elegans*, Northeast Structural Genomics Target WR66
Authors : Keller, J.P.; Xiao, R.; MacDonald, L.; Shen, J.; Acton, T.; Montelione, G.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2002-05-23
Resolution : 1.85 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

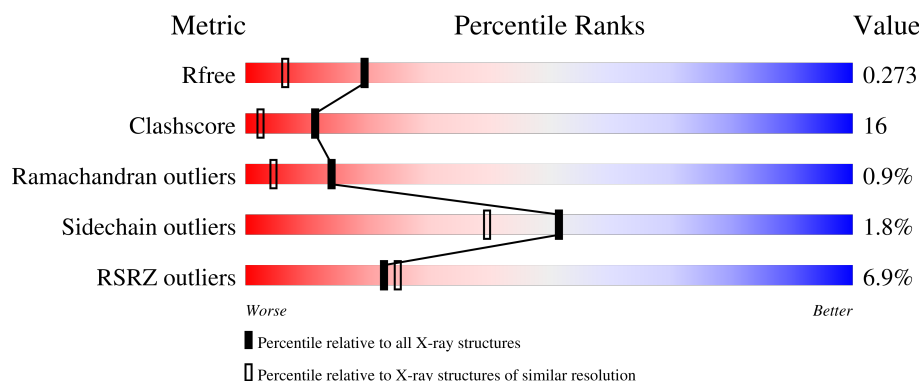
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	339	8% 69% 24% . .
1	B	339	5% 68% 24% . .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5504 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.

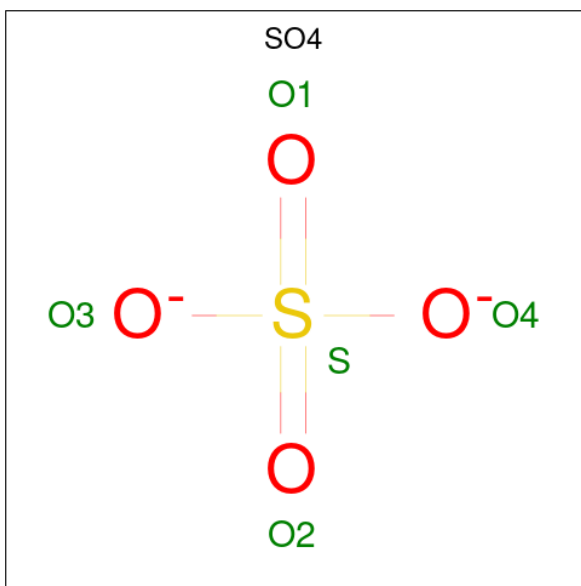
- Molecule 1 is a protein called aldose 1-epimerase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	Se	50	0	0
			2548	1610	433	499	4	2			
1	B	325	Total	C	N	O	S	Se	0	0	0
			2548	1610	433	499	4	2			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	992	MSE	-	expression tag	UNP Q966D4
A	993	SER	-	expression tag	UNP Q966D4
A	994	HIS	-	expression tag	UNP Q966D4
A	995	HIS	-	expression tag	UNP Q966D4
A	996	HIS	-	expression tag	UNP Q966D4
A	997	HIS	-	expression tag	UNP Q966D4
A	998	HIS	-	expression tag	UNP Q966D4
A	999	HIS	-	expression tag	UNP Q966D4
A	1000	SER	-	expression tag	UNP Q966D4
A	1001	MSE	MET	modified residue	UNP Q966D4
A	1082	MSE	MET	modified residue	UNP Q966D4
A	1181	MSE	MET	modified residue	UNP Q966D4
B	1992	MSE	-	expression tag	UNP Q966D4
B	1993	SER	-	expression tag	UNP Q966D4
B	1994	HIS	-	expression tag	UNP Q966D4
B	1995	HIS	-	expression tag	UNP Q966D4
B	1996	HIS	-	expression tag	UNP Q966D4
B	1997	HIS	-	expression tag	UNP Q966D4
B	1998	HIS	-	expression tag	UNP Q966D4
B	1999	HIS	-	expression tag	UNP Q966D4
B	2000	SER	-	expression tag	UNP Q966D4
B	2001	MSE	MET	modified residue	UNP Q966D4
B	2082	MSE	MET	modified residue	UNP Q966D4
B	2181	MSE	MET	modified residue	UNP Q966D4

- 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	B	1	Total	O	S	0	0
			5	4	1		

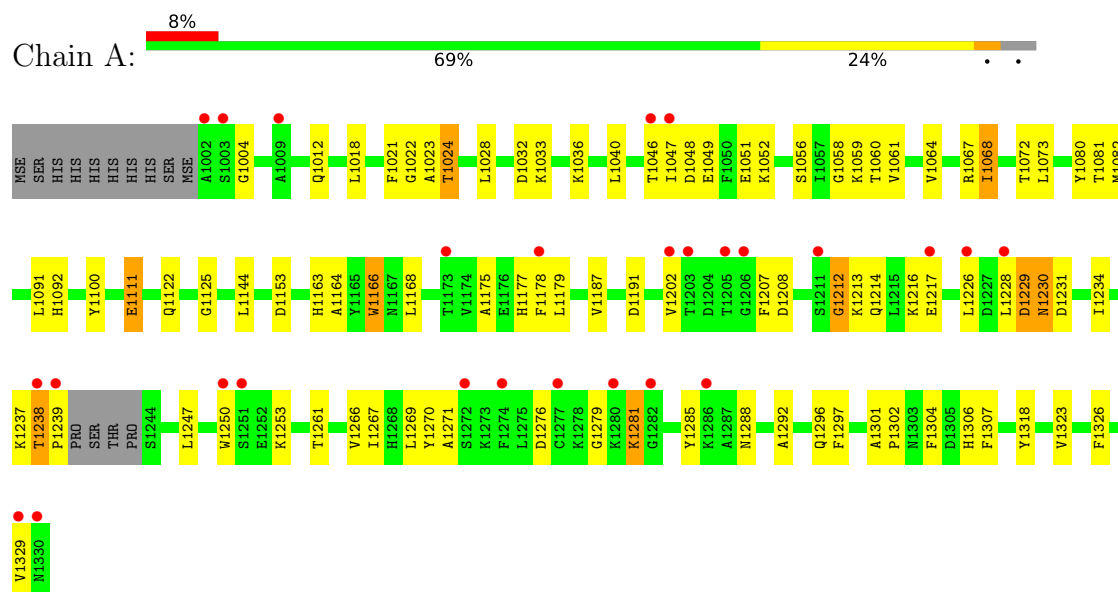
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	162	Total	O	0	0
			162	162		
3	B	236	Total	O	0	0
			236	236		

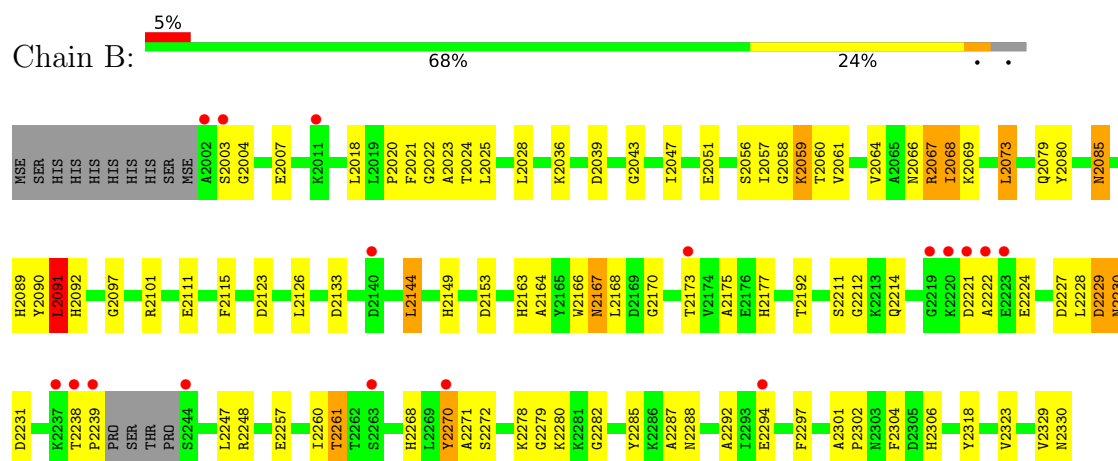
3 Residue-property plots [i](#)

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: aldose 1-epimerase



• Molecule 1: aldose 1-epimerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	50.81Å 91.01Å 83.70Å 90.00° 100.67° 90.00°	Depositor
Resolution (Å)	20.00 – 1.85 20.00 – 1.85	Depositor EDS
% Data completeness (in resolution range)	95.2 (20.00-1.85) 95.1 (20.00-1.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.40 (at 1.84Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.224 , 0.262 0.230 , 0.273	Depositor DCC
R_{free} test set	3097 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	27.2	Xtriage
Anisotropy	0.349	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 43.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5504	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% 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: SO4

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	0.35	0/2608	0.96	12/3535 (0.3%)
1	B	0.37	0/2608	1.03	19/3535 (0.5%)
All	All	0.36	0/5216	0.99	31/7070 (0.4%)

There are no bond length outliers.

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1067	ARG	N-CA-C	9.40	125.31	110.17
1	B	2060	THR	N-CA-C	-8.98	94.68	108.96
1	B	2067	ARG	N-CA-C	8.64	124.09	110.17
1	A	1060	THR	N-CA-C	-8.03	95.64	108.73
1	B	2068	ILE	N-CA-C	-7.74	95.92	107.37
1	A	1068	ILE	N-CA-C	-7.64	96.06	107.37
1	B	2144	LEU	N-CA-C	-7.17	96.72	108.41
1	A	1024	THR	N-CA-C	7.01	121.28	110.20
1	A	1307	PHE	N-CA-C	-6.92	102.18	110.13
1	B	2024	THR	N-CA-C	6.87	121.05	110.20
1	B	2089	HIS	N-CA-C	6.80	120.53	110.46
1	B	2297	PHE	N-CA-C	-6.78	100.48	110.52
1	B	2039	ASP	N-CA-C	-6.25	99.82	109.76
1	B	2153	ASP	N-CA-C	-6.12	106.35	113.88
1	A	1091	LEU	N-CA-C	6.08	119.15	109.24
1	A	1318	TYR	N-CA-C	-6.08	100.10	109.76
1	A	1297	PHE	N-CA-C	-6.07	101.54	110.52
1	B	2192	THR	N-CA-C	-6.00	104.61	113.61
1	A	1168	LEU	N-CA-C	5.98	120.42	113.18
1	A	1153	ASP	N-CA-C	-5.98	106.32	113.97
1	B	2261	THR	N-CA-C	-5.96	99.29	109.24
1	B	2059	LYS	N-CA-C	5.79	118.17	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	2091	LEU	N-CA-C	5.51	119.21	109.96
1	A	1144	LEU	N-CA-C	-5.42	99.69	108.52
1	B	2168	LEU	N-CA-C	5.31	119.74	113.16
1	B	2061	VAL	N-CA-C	5.23	116.50	108.71
1	B	2123	ASP	N-CA-C	5.23	116.79	111.14
1	B	2025	LEU	N-CA-C	-5.08	99.04	107.99
1	A	1166	TRP	N-CA-C	5.04	118.02	109.76
1	B	2272	SER	N-CA-C	5.03	118.54	111.90
1	B	2260	ILE	N-CA-C	5.00	115.92	108.46

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	2548	0	2440	64	0
1	B	2548	0	2440	96	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	162	0	0	4	0
3	B	236	0	0	6	0
All	All	5504	0	4880	159	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 16.

All (159) 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:2278:LYS:HD3	1:B:2282:GLY:HA2	1.39	1.02
1:B:2067:ARG:HH12	1:B:2091:LEU:HD12	1.24	1.00
1:B:2085:ASN:HD21	1:B:2091:LEU:H	0.99	0.93
1:B:2268:HIS:HB3	1:B:2294:GLU:HG3	1.54	0.89
1:A:1238:THR:H	1:A:1239:PRO:HD2	1.42	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2270:TYR:N	1:B:2270:TYR:HD2	1.75	0.85
1:A:1080:TYR:HB3	1:A:1082:MSE:HE1	1.61	0.83
1:A:1082:MSE:HE2	1:A:1125:GLY:O	1.83	0.78
1:B:2067:ARG:HG2	1:B:2067:ARG:HH11	1.47	0.78
1:B:2270:TYR:N	1:B:2270:TYR:CD2	2.50	0.77
1:B:2003:SER:HB3	1:B:2020:PRO:HG2	1.69	0.75
1:B:2066:ASN:HD22	1:B:2067:ARG:H	1.34	0.74
1:B:2091:LEU:HD23	1:B:2091:LEU:O	1.89	0.72
1:B:2085:ASN:ND2	1:B:2091:LEU:H	1.81	0.72
1:B:2066:ASN:H	1:B:2092:HIS:HD2	1.36	0.71
1:B:2270:TYR:CE2	1:B:2292:ALA:HB3	2.27	0.70
1:A:1279:GLY:HA3	1:A:1285:TYR:CE2	2.28	0.69
1:B:2101:ARG:HH11	1:B:2101:ARG:HB3	1.58	0.69
1:A:1081:THR:C	1:A:1082:MSE:HE3	2.18	0.69
1:A:1175:ALA:HB1	1:A:1216:LYS:HB3	1.74	0.69
1:B:2270:TYR:HH	1:B:2294:GLU:CD	2.01	0.69
1:B:2101:ARG:HB3	1:B:2101:ARG:NH1	2.08	0.68
1:B:2091:LEU:HD23	1:B:2091:LEU:C	2.21	0.66
1:A:1187:VAL:HB	1:A:1231:ASP:HB2	1.75	0.66
1:B:2067:ARG:NH1	1:B:2091:LEU:HD12	2.04	0.66
1:B:2066:ASN:H	1:B:2092:HIS:CD2	2.13	0.66
1:B:2085:ASN:HD21	1:B:2091:LEU:N	1.84	0.65
1:A:1228:LEU:HB3	1:A:1271:ALA:HB2	1.79	0.64
1:B:2229:ASP:H	1:B:2230:ASN:HD22	1.45	0.64
1:A:1072:THR:HG23	1:A:1081:THR:HG22	1.79	0.64
1:A:1046:THR:OG1	1:A:1049:GLU:HG3	1.98	0.63
1:A:1048:ASP:OD1	1:A:1052:LYS:HE3	2.00	0.62
1:B:2167:ASN:ND2	1:B:2170:GLY:H	1.97	0.62
1:A:1261:THR:HG22	1:A:1323:VAL:HB	1.80	0.61
1:B:2270:TYR:CZ	1:B:2294:GLU:CD	2.78	0.61
1:B:2270:TYR:HE2	1:B:2294:GLU:HG2	1.65	0.61
1:B:2069:LYS:HE2	3:B:3095:HOH:O	2.00	0.61
1:A:1270:TYR:HB3	1:A:1292:ALA:HB3	1.83	0.60
1:B:2066:ASN:ND2	1:B:2067:ARG:H	2.00	0.59
1:A:1238:THR:N	1:A:1239:PRO:HD2	2.17	0.59
1:A:1018:LEU:N	1:A:1018:LEU:HD22	2.18	0.58
1:A:1247:LEU:C	1:A:1247:LEU:HD23	2.29	0.57
1:B:2247:LEU:C	1:B:2247:LEU:HD23	2.30	0.57
1:A:1238:THR:H	1:A:1239:PRO:CD	2.16	0.57
1:B:2238:THR:N	1:B:2239:PRO:HD2	2.20	0.57
1:B:2270:TYR:CE2	1:B:2294:GLU:HG2	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1213:LYS:HD2	1:A:1217:GLU:HB2	1.87	0.56
1:A:1271:ALA:O	1:A:1288:ASN:HA	2.05	0.56
1:B:2073:LEU:HD21	1:B:2080:TYR:HB2	1.86	0.56
1:A:1021:PHE:CE2	1:A:1051:GLU:HA	2.41	0.56
1:B:2270:TYR:OH	1:B:2294:GLU:CD	2.49	0.56
1:A:1056:SER:O	1:A:1163:HIS:HB3	2.06	0.56
1:A:1266:VAL:HG22	1:A:1296:GLN:O	2.06	0.55
1:B:2270:TYR:CD2	1:B:2292:ALA:HB3	2.42	0.55
1:B:2163:HIS:CE1	1:B:2270:TYR:CE1	2.94	0.55
1:A:1081:THR:O	1:A:1082:MSE:HE3	2.07	0.54
1:B:2018:LEU:HD22	1:B:2018:LEU:N	2.23	0.54
1:B:2004:GLY:HA2	1:B:2047:ILE:CD1	2.36	0.54
1:B:2028:LEU:HD12	1:B:2144:LEU:HD22	1.90	0.54
1:A:1004:GLY:HA2	1:A:1047:ILE:HD11	1.90	0.54
1:B:2073:LEU:N	1:B:2073:LEU:HD23	2.24	0.53
1:B:2228:LEU:O	1:B:2270:TYR:HA	2.09	0.53
1:B:2073:LEU:N	1:B:2073:LEU:CD2	2.72	0.53
1:B:2268:HIS:CD2	1:B:2294:GLU:OE2	2.61	0.53
1:A:1046:THR:HB	3:A:3016:HOH:O	2.09	0.53
1:B:2043:GLY:HA3	1:B:2280:LYS:HE3	1.90	0.53
1:B:2004:GLY:HA2	1:B:2047:ILE:HD11	1.90	0.52
1:B:2067:ARG:HH11	1:B:2067:ARG:CG	2.15	0.52
1:B:2021:PHE:CE2	1:B:2051:GLU:HA	2.44	0.52
1:B:2085:ASN:HD22	1:B:2085:ASN:H	1.57	0.52
1:A:1032:ASP:OD2	1:A:1036:LYS:HB3	2.10	0.51
1:A:1229:ASP:H	1:A:1230:ASN:HD22	1.58	0.51
1:B:2231:ASP:HA	1:B:2268:HIS:HA	1.93	0.51
1:B:2064:VAL:HG11	1:B:2068:ILE:HD11	1.92	0.51
1:A:1281:LYS:HA	1:A:1281:LYS:HE3	1.93	0.51
1:B:2270:TYR:CE2	1:B:2292:ALA:CB	2.93	0.51
1:B:2073:LEU:CD2	1:B:2080:TYR:HB2	2.40	0.51
1:B:2270:TYR:HD2	1:B:2270:TYR:H	1.58	0.50
1:A:1028:LEU:HD22	1:A:1040:LEU:HD12	1.94	0.50
1:B:2271:ALA:O	1:B:2288:ASN:HA	2.11	0.50
1:B:2067:ARG:NH1	1:B:2067:ARG:CG	2.75	0.49
1:A:1279:GLY:HA3	1:A:1285:TYR:CZ	2.47	0.49
1:A:1304:PHE:HB3	1:A:1306:HIS:CE1	2.47	0.49
1:A:1028:LEU:HD13	1:A:1166:TRP:CH2	2.48	0.49
1:B:2133:ASP:OD1	1:B:2149:HIS:HB3	2.13	0.49
1:B:2227:ASP:HB2	3:B:3331:HOH:O	2.12	0.48
1:B:2163:HIS:CE1	1:B:2270:TYR:HE1	2.32	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2278:LYS:CD	1:B:2282:GLY:HA2	2.28	0.48
1:A:1028:LEU:HD11	1:A:1326:PHE:CZ	2.49	0.48
1:B:2028:LEU:HD13	1:B:2166:TRP:CH2	2.48	0.48
1:B:2090:TYR:HE2	1:B:2126:LEU:HD21	1.79	0.48
1:A:1226:LEU:H	1:A:1288:ASN:HD21	1.62	0.47
1:A:1179:LEU:O	1:A:1212:GLY:HA2	2.15	0.47
1:B:2097:GLY:O	1:B:2101:ARG:HG2	2.13	0.47
1:A:1033:LYS:HE2	1:A:1253:LYS:O	2.15	0.47
1:A:1073:LEU:C	1:A:1073:LEU:HD12	2.39	0.47
1:B:2043:GLY:CA	1:B:2280:LYS:HE3	2.43	0.47
1:A:1329:VAL:O	1:A:1329:VAL:HG13	2.15	0.47
1:A:1022:GLY:O	1:A:1023:ALA:C	2.58	0.46
1:B:2073:LEU:HD23	1:B:2073:LEU:H	1.80	0.46
1:A:1177:HIS:O	1:A:1214:GLN:HA	2.14	0.46
1:B:2028:LEU:HD13	1:B:2166:TRP:HH2	1.80	0.46
1:B:2007:GLU:HB3	3:B:3236:HOH:O	2.15	0.46
1:B:2167:ASN:C	1:B:2167:ASN:HD22	2.24	0.46
1:A:1024:THR:HG23	1:A:1058:GLY:C	2.41	0.46
1:A:1202:VAL:CG2	1:A:1207:PHE:C	2.89	0.45
1:B:2022:GLY:O	1:B:2023:ALA:C	2.59	0.45
1:B:2248:ARG:HD2	1:B:2257:GLU:OE2	2.17	0.45
1:A:1301:ALA:N	1:A:1302:PRO:CD	2.80	0.45
1:B:2173:THR:HG23	1:B:2175:ALA:H	1.80	0.45
1:B:2221:ASP:OD1	1:B:2224:GLU:HB3	2.17	0.45
1:B:2229:ASP:H	1:B:2230:ASN:ND2	2.13	0.45
1:A:1122:GLN:CD	3:A:3255:HOH:O	2.60	0.45
1:A:1072:THR:HG23	1:A:1081:THR:CG2	2.46	0.45
1:A:1064:VAL:HG11	1:A:1068:ILE:HD11	1.99	0.44
1:A:1056:SER:HA	1:A:1059:LYS:HD2	1.99	0.44
1:B:2301:ALA:N	1:B:2302:PRO:CD	2.80	0.44
1:A:1208:ASP:O	1:A:1213:LYS:NZ	2.50	0.44
1:A:1237:LYS:N	1:A:1237:LYS:HD2	2.32	0.44
1:A:1058:GLY:HA2	1:A:1164:ALA:HA	1.99	0.44
1:B:2173:THR:HG21	3:B:3274:HOH:O	2.17	0.44
1:B:2058:GLY:HA2	1:B:2164:ALA:HA	1.99	0.43
1:A:1028:LEU:HD13	1:A:1166:TRP:CZ3	2.53	0.43
1:B:2028:LEU:CD1	1:B:2144:LEU:HD22	2.49	0.43
1:B:2056:SER:HA	1:B:2059:LYS:HD2	1.99	0.43
1:B:2133:ASP:OD2	1:B:2149:HIS:HD2	2.02	0.43
1:B:2149:HIS:HA	1:B:2318:TYR:O	2.18	0.43
1:B:2304:PHE:HB3	1:B:2306:HIS:CE1	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2115:PHE:O	1:B:2133:ASP:HA	2.18	0.43
1:B:2261:THR:HG22	1:B:2323:VAL:HB	2.00	0.43
1:B:2279:GLY:HA3	1:B:2285:TYR:CZ	2.54	0.43
1:B:2079:GLN:HG3	3:B:3209:HOH:O	2.19	0.43
1:B:2085:ASN:HD22	1:B:2085:ASN:N	2.15	0.43
1:B:2279:GLY:HA3	1:B:2285:TYR:CE1	2.54	0.43
1:A:1100:TYR:HE2	3:A:3151:HOH:O	2.02	0.42
1:B:2287:ALA:O	1:B:2288:ASN:HB2	2.18	0.42
1:A:1080:TYR:CB	1:A:1082:MSE:HE1	2.43	0.42
1:A:1226:LEU:H	1:A:1288:ASN:ND2	2.18	0.42
1:A:1229:ASP:HA	1:A:1269:LEU:O	2.19	0.42
1:A:1261:THR:CG2	1:A:1323:VAL:HB	2.47	0.42
1:A:1028:LEU:C	1:A:1028:LEU:HD23	2.45	0.42
1:B:2057:ILE:HG22	3:B:3122:HOH:O	2.17	0.42
1:A:1111:GLU:CD	1:A:1111:GLU:H	2.27	0.42
1:A:1061:VAL:HG13	1:A:1092:HIS:HB3	2.01	0.42
1:B:2163:HIS:NE2	1:B:2270:TYR:HE1	2.18	0.42
1:A:1271:ALA:HB3	3:A:3054:HOH:O	2.20	0.41
1:B:2028:LEU:C	1:B:2028:LEU:HD23	2.46	0.41
1:B:2270:TYR:CE2	1:B:2294:GLU:CD	2.98	0.41
1:B:2270:TYR:CZ	1:B:2292:ALA:CB	3.04	0.41
1:B:2067:ARG:HG2	1:B:2067:ARG:NH1	2.25	0.41
1:B:2177:HIS:O	1:B:2214:GLN:HA	2.21	0.41
1:A:1012:GLN:HA	1:B:2036:LYS:HB2	2.03	0.40
1:B:2329:VAL:O	1:B:2330:ASN:HB2	2.20	0.40
1:A:1234:ILE:HD11	1:A:1267:ILE:HG13	2.03	0.40
1:B:2167:ASN:HD21	1:B:2170:GLY:H	1.70	0.40
1:A:1178:PHE:HB2	1:A:1250:TRP:CE2	2.57	0.40
1:A:1238:THR:N	1:A:1239:PRO:CD	2.80	0.40
1:B:2111:GLU:CD	1:B:2111:GLU:H	2.30	0.40
1:B:2211:SER:O	1:B:2212:GLY:C	2.65	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	321/339 (95%)	302 (94%)	15 (5%)	4 (1%)	10	2
1	B	321/339 (95%)	304 (95%)	15 (5%)	2 (1%)	21	10
All	All	642/678 (95%)	606 (94%)	30 (5%)	6 (1%)	14	4

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1229	ASP
1	B	2222	ALA
1	A	1212	GLY
1	A	1238	THR
1	A	1276	ASP
1	B	2229	ASP

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	276/286 (96%)	272 (99%)	4 (1%)	59	49
1	B	276/286 (96%)	270 (98%)	6 (2%)	45	32
All	All	552/572 (96%)	542 (98%)	10 (2%)	51	40

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1111	GLU
1	A	1191	ASP
1	A	1230	ASN
1	A	1281	LYS
1	B	2073	LEU
1	B	2085	ASN
1	B	2091	LEU

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Mol	Chain	Res	Type
1	B	2167	ASN
1	B	2230	ASN
1	B	2270	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1037	ASN
1	A	1079	GLN
1	A	1149	HIS
1	A	1230	ASN
1	A	1288	ASN
1	A	1317	HIS
1	B	2012	GLN
1	B	2038	GLN
1	B	2066	ASN
1	B	2070	ASN
1	B	2079	GLN
1	B	2085	ASN
1	B	2086	ASN
1	B	2092	HIS
1	B	2096	ASN
1	B	2122	GLN
1	B	2149	HIS
1	B	2167	ASN
1	B	2214	GLN
1	B	2230	ASN
1	B	2317	HIS

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 [i](#)

2 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	$\# Z > 2$
2	SO4	A	4001	-	4,4,4	0.39	0	6,6,6	0.08	0
2	SO4	B	4002	-	4,4,4	0.36	0	6,6,6	0.09	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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

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	316/339 (93%)	0.70	27 (8%) 16 17	20, 35, 57, 71	0
1	B	323/339 (95%)	0.29	17 (5%) 32 35	18, 28, 46, 68	0
All	All	639/678 (94%)	0.50	44 (6%) 23 25	18, 31, 56, 71	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	2270	TYR	10.4
1	A	1239	PRO	5.3
1	A	1002	ALA	5.0
1	A	1003	SER	4.6
1	B	2222	ALA	4.4
1	A	1238	THR	4.1
1	B	2220	LYS	3.8
1	B	2239	PRO	3.6
1	B	2221	ASP	3.4
1	A	1205	THR	3.3
1	B	2173	THR	3.2
1	B	2002	ALA	3.1
1	B	2219	GLY	3.0
1	A	1173	THR	2.9
1	A	1202	VAL	2.7
1	A	1046	THR	2.7
1	A	1282	GLY	2.6
1	A	1286	LYS	2.6
1	B	2237	LYS	2.6
1	A	1330	ASN	2.5
1	A	1211	SER	2.5
1	B	2238	THR	2.5
1	A	1251	SER	2.4
1	A	1329	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	2294	GLU	2.4
1	A	1277	CYS	2.4
1	B	2003	SER	2.4
1	A	1228	LEU	2.4
1	B	2223	GLU	2.3
1	A	1274	PHE	2.3
1	A	1217	GLU	2.3
1	A	1250	TRP	2.2
1	A	1206	GLY	2.2
1	A	1047	ILE	2.2
1	B	2263	SER	2.2
1	A	1009	ALA	2.2
1	A	1203	THR	2.1
1	A	1272	SER	2.1
1	A	1280	LYS	2.1
1	B	2011	LYS	2.0
1	A	1226	LEU	2.0
1	B	2140	ASP	2.0
1	B	2244	SER	2.0
1	A	1178	PHE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	A	4001	5/5	0.83	0.09	71,71,71,71	0
2	SO4	B	4002	5/5	0.95	0.10	44,45,46,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.