



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 08:18 AM UTC

PDB ID : 1EUA / pdb_00001eua
Title : SCHIFF BASE INTERMEDIATE IN KDPG ALDOLASE FROM ES-
CHERICHIA COLI
Authors : Allard, J.; Grochulski, P.; Sygusch, J.
Deposited on : 2000-04-14
Resolution : 1.95 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

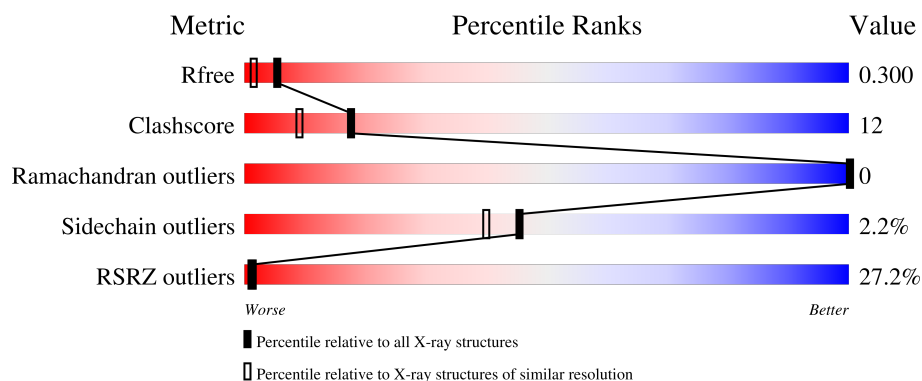
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	213	12% 77% 23%
1	B	213	35% 79% 21%
1	C	213	35% 74% 23% .

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
2	ACT	A	2002	-	-	X	-
4	PYR	A	2101	-	X	-	-
4	PYR	B	2102	-	X	-	-
4	PYR	C	2103	-	X	-	-

2 Entry composition [i](#)

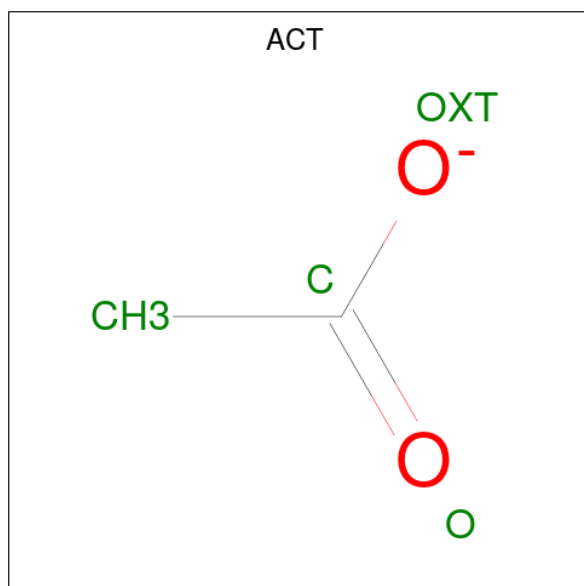
There are 5 unique types of molecules in this entry. The entry contains 5355 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 KDPG ALDOLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	213	Total	C	N	O	S	0	0	0
			1566	1006	260	293	7			
1	B	213	Total	C	N	O	S	0	0	0
			1566	1006	260	293	7			
1	C	213	Total	C	N	O	S	0	0	0
			1566	1006	260	293	7			

- Molecule 2 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



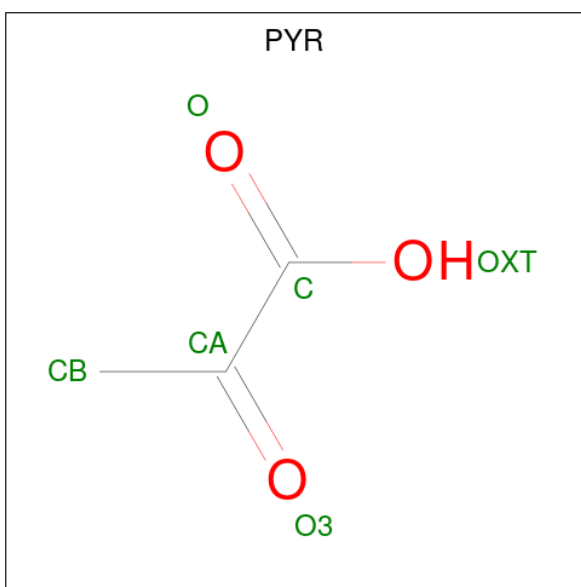
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



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

- Molecule 4 is PYRUVIC ACID (CCD ID: PYR) (formula: C₃H₄O₃).



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

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

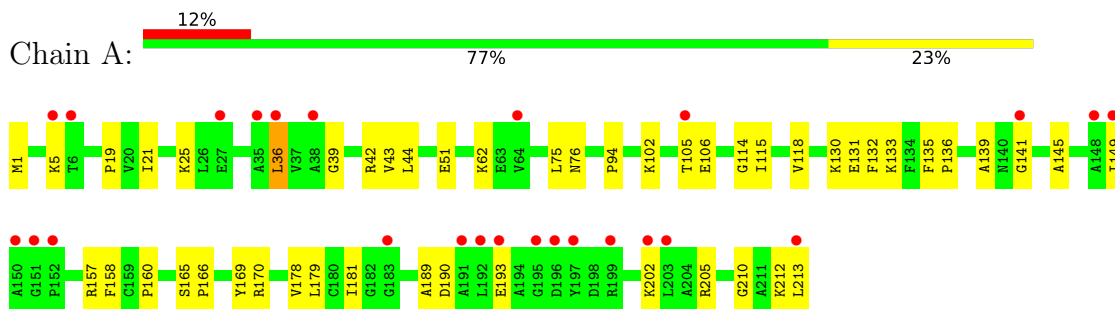
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	193	Total	O	0	0
			193	193		
5	B	248	Total	O	0	0
			248	248		
5	C	175	Total	O	0	0
			175	175		

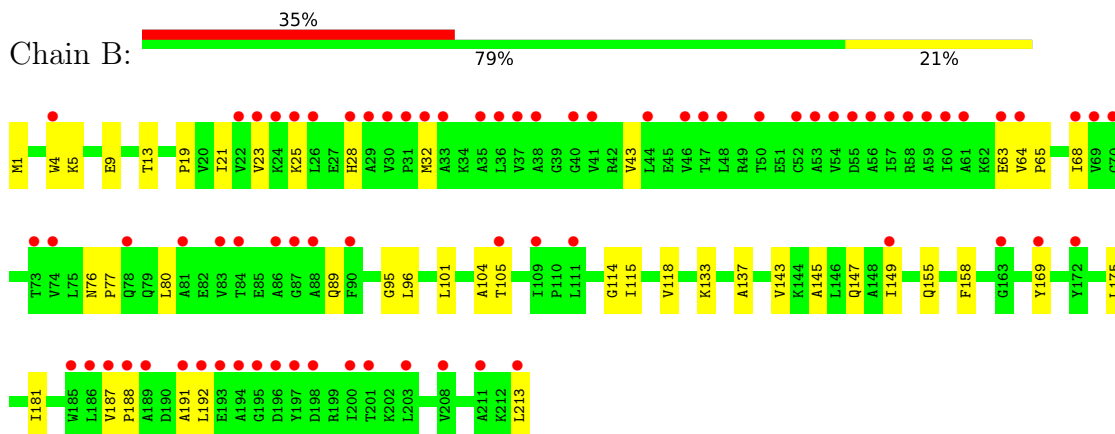
3 Residue-property plots [i](#)

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.

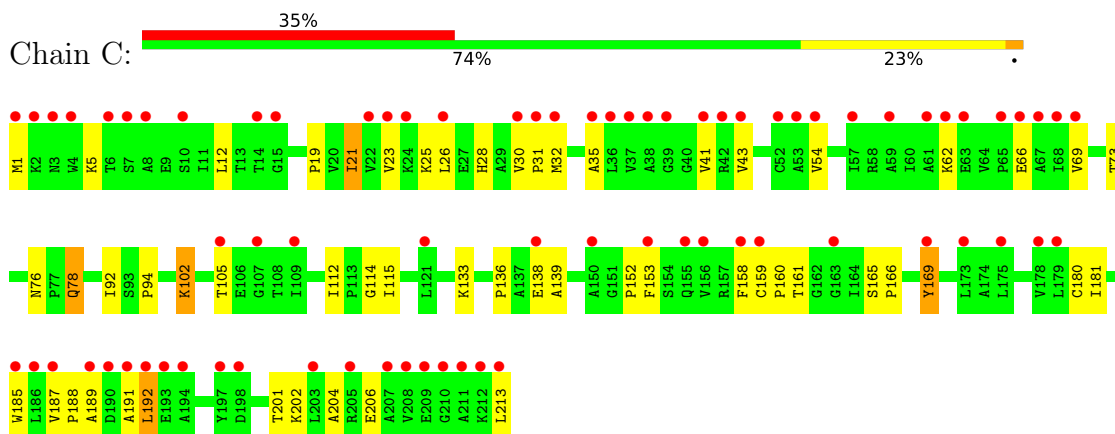
• Molecule 1: KDPG ALDOLASE



• Molecule 1: KDPG ALDOLASE



• Molecule 1: KDPG ALDOLASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.41Å 85.72Å 133.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.37 – 1.95 50.37 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.1 (50.37-1.95) 96.1 (50.37-1.95)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.70 (at 1.95Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.201 , 0.243 0.263 , 0.300	Depositor DCC
R_{free} test set	2280 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	25.7	Xtriage
Anisotropy	0.527	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 53.1	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	0.92	EDS
Total number of atoms	5355	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.11% 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: PYR, SO4, 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	0.39	0/1594	0.91	2/2168 (0.1%)
1	B	0.41	0/1594	0.94	4/2168 (0.2%)
1	C	0.38	0/1594	0.91	3/2168 (0.1%)
All	All	0.40	0/4782	0.92	9/6504 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	21	ILE	N-CA-C	7.51	119.39	108.58
1	A	21	ILE	N-CA-C	7.02	119.54	108.87
1	C	21	ILE	N-CA-C	6.08	118.11	108.87
1	B	158	PHE	N-CA-C	5.72	118.39	109.52
1	A	76	ASN	N-CA-C	5.63	113.08	108.07
1	B	137	ALA	N-CA-C	5.45	117.67	111.02
1	B	133	LYS	N-CA-C	-5.34	101.57	110.17
1	C	158	PHE	N-CA-C	5.26	117.41	109.41
1	C	69	VAL	N-CA-C	5.09	115.91	108.58

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	1566	0	1630	38	0
1	B	1566	0	1630	33	0
1	C	1566	0	1629	51	0
2	A	8	0	6	3	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	5	0	0	0	0
4	A	6	0	0	0	0
4	B	6	0	0	0	0
4	C	6	0	0	0	0
5	A	193	0	0	6	1
5	B	248	0	0	3	2
5	C	175	0	0	10	0
All	All	5355	0	4895	119	2

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 12.

All (119) 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:A:36:LEU:HD23	1:A:44:LEU:HD22	1.57	0.87
1:C:187:VAL:HG12	1:C:192:LEU:HD21	1.61	0.83
1:A:160:PRO:HG2	1:A:181:ILE:HG22	1.64	0.79
1:C:35:ALA:HB1	5:C:2368:HOH:O	1.82	0.78
1:A:51:GLU:H	1:A:51:GLU:CD	1.91	0.78
1:A:212:LYS:O	1:A:213:LEU:HB2	1.86	0.74
1:A:136:PRO:HG2	1:A:139:ALA:HB3	1.69	0.74
1:B:1:MET:HE1	1:B:105:THR:HG22	1.68	0.73
1:A:189:ALA:O	1:A:193:GLU:HG3	1.87	0.72
1:A:5:LYS:HD3	1:A:5:LYS:C	2.15	0.71
1:C:76:ASN:OD1	1:C:78:GLN:HG3	1.90	0.71
1:C:136:PRO:HG2	1:C:139:ALA:HB3	1.72	0.71
1:B:1:MET:CE	1:B:105:THR:HG22	2.20	0.71
1:B:155:GLN:HB2	5:B:2323:HOH:O	1.92	0.68
1:C:187:VAL:CG1	1:C:192:LEU:HD21	2.23	0.68
1:A:165:SER:HB2	1:A:166:PRO:HD2	1.75	0.68
1:C:21:ILE:HG21	1:C:32:MET:HE1	1.76	0.66
1:C:192:LEU:HD22	1:C:192:LEU:H	1.61	0.65
1:A:102:LYS:O	1:A:106:GLU:HG3	1.97	0.64
1:A:190:ASP:HA	1:A:193:GLU:OE1	1.99	0.62
1:C:165:SER:HB2	1:C:166:PRO:HD2	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:201:THR:N	5:C:2368:HOH:O	2.32	0.62
1:C:189:ALA:HB3	5:C:2311:HOH:O	1.98	0.61
1:C:160:PRO:HG2	1:C:181:ILE:HG22	1.83	0.60
1:A:19:PRO:HD2	1:A:43:VAL:O	2.02	0.59
1:B:145:ALA:O	1:B:149:ILE:HD13	2.01	0.59
1:C:1:MET:HE2	1:C:105:THR:HG22	1.83	0.59
1:C:1:MET:CE	1:C:105:THR:HG22	2.33	0.59
1:C:152:PRO:HG2	1:C:153:PHE:CD1	2.38	0.59
1:C:192:LEU:HD22	1:C:192:LEU:N	2.19	0.57
1:B:143:VAL:O	1:B:147:GLN:HG3	2.05	0.56
1:C:62:LYS:HG3	5:C:2278:HOH:O	2.06	0.55
1:C:12:LEU:HD21	1:C:112:ILE:HD13	1.87	0.55
1:C:181:ILE:HD12	1:C:181:ILE:C	2.31	0.55
1:B:188:PRO:HG2	1:B:191:ALA:HB2	1.88	0.55
1:C:114:GLY:HA2	1:C:133:LYS:HB3	1.89	0.55
1:C:202:LYS:HG2	1:C:206:GLU:OE2	2.07	0.54
1:A:136:PRO:HG3	1:B:149:ILE:HD12	1.91	0.52
1:B:181:ILE:C	1:B:181:ILE:HD12	2.34	0.52
1:A:114:GLY:HA2	1:A:133:LYS:O	2.09	0.52
1:B:188:PRO:HG2	1:B:191:ALA:CB	2.39	0.52
1:A:1:MET:HB3	5:A:2335:HOH:O	2.08	0.52
1:B:19:PRO:HD2	1:B:43:VAL:O	2.10	0.52
1:C:138:GLU:HB2	5:C:2258:HOH:O	2.09	0.52
1:B:76:ASN:HB2	1:B:77:PRO:HD2	1.92	0.52
1:C:26:LEU:O	1:C:26:LEU:HD23	2.09	0.52
1:B:32:MET:HE2	1:B:192:LEU:HD21	1.91	0.51
1:A:25:LYS:H	2:A:2002:ACT:CH3	2.24	0.50
1:C:188:PRO:O	1:C:191:ALA:HB3	2.11	0.50
1:A:130:LYS:HG3	1:A:131:GLU:OE2	2.12	0.49
1:C:192:LEU:H	1:C:192:LEU:CD2	2.23	0.49
1:B:68:ILE:HA	1:B:89:GLN:NE2	2.28	0.49
1:B:143:VAL:HG13	1:B:175:LEU:HD11	1.94	0.49
1:C:114:GLY:HA2	1:C:133:LYS:O	2.13	0.49
1:B:9:GLU:O	1:B:13:THR:HG23	2.13	0.49
1:C:26:LEU:HD23	1:C:26:LEU:C	2.38	0.49
1:A:181:ILE:C	1:A:181:ILE:HD12	2.38	0.48
1:C:161:THR:HG21	5:C:2378:HOH:O	2.13	0.48
1:A:213:LEU:HD22	5:A:2377:HOH:O	2.13	0.48
1:B:25:LYS:HB2	1:B:28:HIS:HD2	1.79	0.48
1:B:80:LEU:HD21	1:B:104:ALA:HA	1.96	0.47
1:C:169:TYR:HB2	1:C:181:ILE:HG12	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:MET:HE1	1:B:105:THR:CG2	2.40	0.47
1:C:102:LYS:HD3	5:C:2340:HOH:O	2.15	0.47
1:C:115:ILE:C	1:C:115:ILE:HD12	2.40	0.47
1:C:19:PRO:HD2	1:C:43:VAL:O	2.15	0.47
1:C:23:VAL:HG23	1:C:23:VAL:O	2.15	0.46
1:A:115:ILE:HD12	1:A:115:ILE:C	2.40	0.46
1:C:54:VAL:HG12	5:C:2275:HOH:O	2.14	0.46
1:A:25:LYS:H	2:A:2002:ACT:H3	1.79	0.46
1:C:102:LYS:HB2	1:C:102:LYS:NZ	2.31	0.46
1:A:105:THR:HG21	5:A:2315:HOH:O	2.15	0.46
1:A:131:GLU:HG3	1:A:157:ARG:HB2	1.98	0.46
1:B:95:GLY:HA2	1:B:114:GLY:O	2.16	0.45
1:B:76:ASN:HB2	1:B:77:PRO:CD	2.46	0.45
1:C:92:ILE:CG2	1:C:133:LYS:HD2	2.47	0.45
1:C:41:VAL:HG23	1:C:204:ALA:HB1	1.98	0.45
1:C:166:PRO:HD3	1:C:185:TRP:CE2	2.52	0.45
1:B:1:MET:HG3	1:B:4:TRP:CZ3	2.52	0.45
1:B:96:LEU:HD12	1:B:96:LEU:O	2.16	0.45
1:A:135:PHE:CD1	1:A:136:PRO:HA	2.52	0.45
1:A:202:LYS:HD2	5:A:2249:HOH:O	2.17	0.45
1:A:94:PRO:O	1:B:118:VAL:HG21	2.17	0.44
1:C:114:GLY:CA	1:C:133:LYS:HB3	2.47	0.44
1:B:23:VAL:HB	1:B:32:MET:HE1	1.99	0.44
1:A:39:GLY:HA2	1:A:205:ARG:HG3	2.00	0.44
1:C:202:LYS:HE2	1:C:206:GLU:OE2	2.17	0.44
1:A:118:VAL:HG21	1:C:94:PRO:O	2.17	0.44
1:B:5:LYS:HB2	5:B:2291:HOH:O	2.18	0.44
1:B:64:VAL:N	1:B:65:PRO:HD3	2.32	0.44
1:A:160:PRO:HD2	1:A:178:VAL:HG13	2.00	0.43
1:A:1:MET:HE3	1:A:105:THR:HG23	2.00	0.43
1:B:63:GLU:C	1:B:65:PRO:HD3	2.43	0.43
1:C:25:LYS:HB2	1:C:28:HIS:CD2	2.54	0.43
1:A:132:PHE:O	1:A:158:PHE:HA	2.19	0.43
1:C:25:LYS:HB2	1:C:28:HIS:HD2	1.84	0.43
1:C:30:VAL:N	1:C:31:PRO:HD2	2.34	0.43
1:C:213:LEU:HD23	1:C:213:LEU:H	1.83	0.43
1:B:25:LYS:HB2	1:B:28:HIS:CD2	2.54	0.42
1:A:157:ARG:HB3	1:A:179:LEU:HG	2.01	0.42
1:A:51:GLU:CD	1:A:51:GLU:N	2.69	0.42
1:A:170:ARG:NH1	1:A:210:GLY:O	2.51	0.42
1:A:5:LYS:C	1:A:5:LYS:CD	2.88	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:GLY:HA3	1:A:145:ALA:HB2	2.01	0.42
1:C:73:THR:HG22	1:C:73:THR:O	2.18	0.42
1:A:149:ILE:HD11	5:A:2395:HOH:O	2.18	0.42
1:B:32:MET:HE2	1:B:192:LEU:CD2	2.50	0.42
1:B:187:VAL:HB	5:B:2278:HOH:O	2.19	0.42
1:C:161:THR:CG2	5:C:2378:HOH:O	2.67	0.42
1:B:5:LYS:HB2	1:B:5:LYS:HE2	1.94	0.42
1:C:73:THR:HA	1:C:94:PRO:HD3	2.00	0.42
1:A:25:LYS:HB2	2:A:2002:ACT:H3	2.01	0.41
1:C:159:CYS:HA	1:C:180:CYS:O	2.19	0.41
1:A:62:LYS:HG3	5:A:2260:HOH:O	2.20	0.41
1:C:5:LYS:HB2	5:C:2297:HOH:O	2.20	0.41
1:B:101:LEU:O	1:B:105:THR:HG23	2.21	0.41
1:B:115:ILE:HD12	1:B:115:ILE:C	2.45	0.41
1:C:166:PRO:HA	1:C:169:TYR:CD2	2.56	0.41
1:C:213:LEU:HD23	1:C:213:LEU:N	2.36	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:2239:HOH:O	5:B:2430:HOH:O[4_446]	2.18	0.02
5:B:2399:HOH:O	5:B:2434:HOH:O[4_446]	2.19	0.01

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	211/213 (99%)	207 (98%)	4 (2%)	0	100	100
1	B	211/213 (99%)	210 (100%)	1 (0%)	0	100	100
1	C	211/213 (99%)	205 (97%)	6 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	633/639 (99%)	622 (98%)	11 (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	163/163 (100%)	159 (98%)	4 (2%)	42	34
1	B	163/163 (100%)	161 (99%)	2 (1%)	63	61
1	C	163/163 (100%)	158 (97%)	5 (3%)	35	26
All	All	489/489 (100%)	478 (98%)	11 (2%)	45	40

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	LEU
1	A	42	ARG
1	A	75	LEU
1	A	169	TYR
1	B	169	TYR
1	B	213	LEU
1	C	66	GLU
1	C	78	GLN
1	C	102	LYS
1	C	169	TYR
1	C	192	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	89	GLN
1	C	78	GLN

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Mol	Chain	Res	Type
1	C	89	GLN
1	C	147	GLN

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 ⓘ

8 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	ACT	A	2002	-	3,3,3	1.10	0	3,3,3	0.36	0
4	PYR	C	2103	1	5,5,5	4.31	2 (40%)	3,6,6	3.18	2 (66%)
3	SO4	B	2201	-	4,4,4	0.34	0	6,6,6	0.09	0
4	PYR	A	2101	1	5,5,5	4.40	3 (60%)	3,6,6	3.10	1 (33%)
4	PYR	B	2102	1	5,5,5	4.47	2 (40%)	3,6,6	3.05	2 (66%)
3	SO4	A	2202	-	4,4,4	0.36	0	6,6,6	0.10	0
3	SO4	C	2203	-	4,4,4	0.33	0	6,6,6	0.12	0
2	ACT	A	2001	-	3,3,3	1.11	0	3,3,3	0.39	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	PYR	C	2103	1	-	3/4/4/4	-
4	PYR	A	2101	1	-	4/4/4/4	-
4	PYR	B	2102	1	-	2/4/4/4	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	2102	PYR	O3-CA	9.27	1.44	1.23
4	A	2101	PYR	O3-CA	8.96	1.43	1.23
4	C	2103	PYR	O3-CA	8.89	1.43	1.23
4	B	2102	PYR	CA-C	-2.65	1.45	1.54
4	A	2101	PYR	CA-C	-2.60	1.45	1.54
4	C	2103	PYR	CA-C	-2.48	1.46	1.54
4	A	2101	PYR	CB-CA	-2.08	1.46	1.50

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	2101	PYR	O3-CA-CB	-4.82	108.97	119.77
4	C	2103	PYR	O3-CA-CB	-4.72	109.19	119.77
4	B	2102	PYR	O3-CA-CB	-4.52	109.65	119.77
4	B	2102	PYR	OXT-C-CA	2.34	120.07	113.59
4	C	2103	PYR	OXT-C-CA	2.32	120.03	113.59

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	2101	PYR	O-C-CA-O3
4	A	2101	PYR	OXT-C-CA-O3
4	A	2101	PYR	OXT-C-CA-CB
4	B	2102	PYR	O-C-CA-CB
4	B	2102	PYR	OXT-C-CA-CB
4	C	2103	PYR	O-C-CA-CB
4	C	2103	PYR	OXT-C-CA-CB
4	A	2101	PYR	O-C-CA-CB
4	C	2103	PYR	O-C-CA-O3

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2002	ACT	3	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	213/213 (100%)	1.01	25 (11%)	9 11	17, 31, 50, 62	0
1	B	213/213 (100%)	1.77	74 (34%)	1 0	14, 23, 38, 55	0
1	C	213/213 (100%)	1.76	75 (35%)	1 0	17, 33, 51, 65	0
All	All	639/639 (100%)	1.51	174 (27%)	1 1	14, 29, 48, 65	0

All (174) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	213	LEU	4.9
1	B	26	LEU	4.8
1	C	213	LEU	4.7
1	B	30	VAL	4.6
1	B	29	ALA	4.6
1	C	105	THR	4.6
1	B	52	CYS	4.6
1	B	105	THR	4.3
1	C	211	ALA	4.2
1	C	8	ALA	4.1
1	B	192	LEU	4.1
1	B	56	ALA	4.0
1	B	54	VAL	3.9
1	B	35	ALA	3.9
1	C	23	VAL	3.9
1	B	187	VAL	3.9
1	C	6	THR	3.8
1	B	59	ALA	3.8
1	B	37	VAL	3.8
1	A	27	GLU	3.8
1	B	36	LEU	3.8
1	B	60	ILE	3.8
1	C	38	ALA	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	64	VAL	3.8
1	B	186	LEU	3.7
1	C	54	VAL	3.6
1	C	62	LYS	3.6
1	C	212	LYS	3.6
1	B	70	GLY	3.5
1	C	156	VAL	3.5
1	B	189	ALA	3.4
1	B	22	VAL	3.4
1	B	48	LEU	3.4
1	C	197	TYR	3.3
1	B	83	VAL	3.3
1	B	68	ILE	3.3
1	A	149	ILE	3.3
1	B	193	GLU	3.2
1	B	38	ALA	3.2
1	B	149	ILE	3.2
1	B	188	PRO	3.2
1	B	69	VAL	3.2
1	A	192	LEU	3.1
1	A	150	ALA	3.1
1	A	203	LEU	3.1
1	B	213	LEU	3.1
1	B	197	TYR	3.0
1	C	178	VAL	3.0
1	B	163	GLY	3.0
1	B	61	ALA	3.0
1	C	68	ILE	3.0
1	C	42	ARG	3.0
1	A	197	TYR	3.0
1	B	196	ASP	3.0
1	C	194	ALA	3.0
1	B	63	GLU	3.0
1	A	64	VAL	2.9
1	B	55	ASP	2.9
1	C	107	GLY	2.9
1	C	186	LEU	2.9
1	C	10	SER	2.9
1	C	187	VAL	2.9
1	B	32	MET	2.9
1	B	195	GLY	2.9
1	B	88	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	194	ALA	2.9
1	C	4	TRP	2.9
1	B	57	ILE	2.9
1	C	36	LEU	2.9
1	B	25	LYS	2.9
1	C	190	ASP	2.9
1	B	23	VAL	2.8
1	C	191	ALA	2.8
1	C	37	VAL	2.8
1	C	203	LEU	2.8
1	A	105	THR	2.8
1	B	24	LYS	2.8
1	B	208	VAL	2.8
1	C	41	VAL	2.8
1	B	53	ALA	2.8
1	C	53	ALA	2.8
1	C	192	LEU	2.8
1	C	198	ASP	2.7
1	C	185	TRP	2.7
1	B	33	ALA	2.7
1	A	151	GLY	2.7
1	B	40	GLY	2.7
1	B	4	TRP	2.7
1	B	81	ALA	2.7
1	C	31	PRO	2.7
1	C	1	MET	2.6
1	B	201	THR	2.6
1	C	65	PRO	2.6
1	B	74	VAL	2.6
1	A	196	ASP	2.6
1	B	191	ALA	2.6
1	C	61	ALA	2.6
1	C	14	THR	2.6
1	A	36	LEU	2.6
1	B	86	ALA	2.6
1	C	35	ALA	2.6
1	B	87	GLY	2.6
1	C	15	GLY	2.6
1	A	141	GLY	2.5
1	B	211	ALA	2.5
1	C	22	VAL	2.5
1	B	31	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	90	PHE	2.5
1	A	191	ALA	2.5
1	B	41	VAL	2.5
1	B	84	THR	2.5
1	C	208	VAL	2.5
1	C	210	GLY	2.5
1	C	3	ASN	2.5
1	B	172	TYR	2.4
1	B	44	LEU	2.4
1	B	46	VAL	2.4
1	C	209	GLU	2.4
1	C	32	MET	2.4
1	C	205	ARG	2.4
1	B	185	TRP	2.4
1	C	24	LYS	2.4
1	C	121	LEU	2.4
1	C	179	LEU	2.4
1	C	109	ILE	2.4
1	C	30	VAL	2.4
1	B	50	THR	2.3
1	C	138	GLU	2.3
1	B	78	GLN	2.3
1	B	58	ARG	2.3
1	C	158	PHE	2.3
1	A	5	LYS	2.3
1	A	183	GLY	2.3
1	B	47	THR	2.3
1	A	148	ALA	2.3
1	C	67	ALA	2.3
1	C	163	GLY	2.3
1	C	155	GLN	2.3
1	C	69	VAL	2.3
1	C	150	ALA	2.2
1	C	189	ALA	2.2
1	A	202	LYS	2.2
1	A	6	THR	2.2
1	A	35	ALA	2.2
1	C	59	ALA	2.2
1	B	169	TYR	2.2
1	B	73	THR	2.2
1	C	52	CYS	2.2
1	C	26	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	198	ASP	2.2
1	C	39	GLY	2.2
1	C	153	PHE	2.2
1	C	173	LEU	2.2
1	A	38	ALA	2.2
1	C	207	ALA	2.2
1	C	57	ILE	2.1
1	A	152	PRO	2.1
1	C	175	LEU	2.1
1	C	43	VAL	2.1
1	C	63	GLU	2.1
1	B	28	HIS	2.1
1	B	111	LEU	2.1
1	B	200	ILE	2.1
1	C	169	TYR	2.1
1	A	199	ARG	2.1
1	A	193	GLU	2.1
1	B	203	LEU	2.1
1	C	66	GLU	2.0
1	C	193	GLU	2.0
1	C	7	SER	2.0
1	C	159	CYS	2.0
1	C	2	LYS	2.0
1	B	109	ILE	2.0
1	A	195	GLY	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ACT	A	2001	4/4	0.68	0.20	41,43,44,44	0
2	ACT	A	2002	4/4	0.72	0.18	46,48,49,49	0
4	PYR	A	2101	6/6	0.76	0.15	30,33,36,36	0
3	SO4	A	2202	5/5	0.79	0.15	64,64,65,66	5
4	PYR	B	2102	6/6	0.80	0.14	25,27,31,32	0
4	PYR	C	2103	6/6	0.81	0.14	32,36,39,42	0
3	SO4	B	2201	5/5	0.84	0.18	47,49,50,50	5
3	SO4	C	2203	5/5	0.89	0.13	45,46,47,47	5

6.5 Other polymers [i](#)

There are no such residues in this entry.