



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 02:44 AM UTC

PDB ID : 2O59 / pdb_00002o59
Title : Structure of E. coli topoisomerase III in complex with an 8-base single stranded oligonucleotide. Frozen in glycerol pH 8.0
Authors : Changela, A.; DiGate, R.J.; Mondragon, A.
Deposited on : 2006-12-05
Resolution : 2.50 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

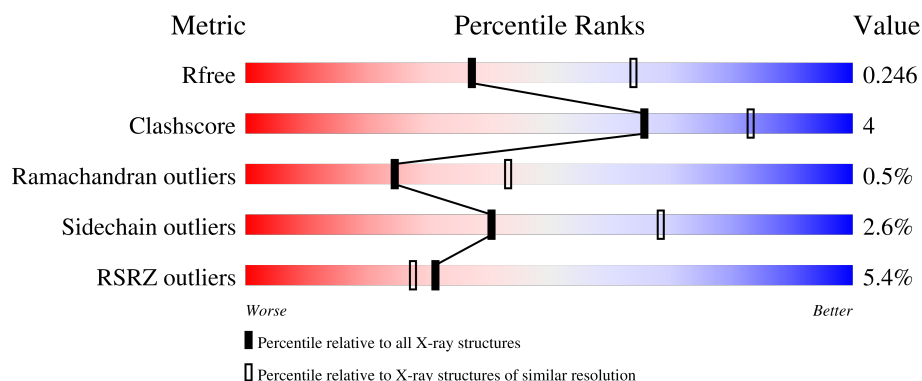
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	C	8	
1	D	8	
2	A	659	
2	B	659	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10353 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 DNA chain called 5'-D(*CP*GP*CP*AP*AP*CP*TP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	7	Total	C	N	O	P	0	0	0
			138	67	26	39	6			
1	D	8	Total	C	N	O	P	0	0	0
			158	77	28	46	7			

- Molecule 2 is a protein called DNA topoisomerase 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	619	Total	C	N	O	S	0	0	0
			4927	3117	898	892	20			
2	B	626	Total	C	N	O	S	0	0	0
			4982	3149	904	909	20			

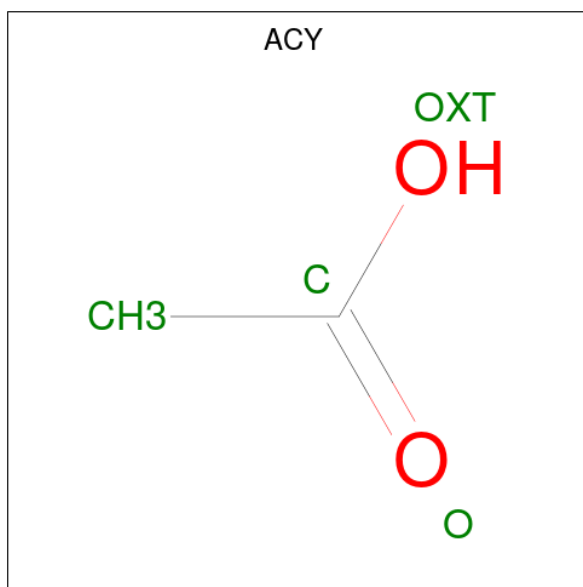
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	654	HIS	-	expression tag	UNP P14294
A	655	HIS	-	expression tag	UNP P14294
A	656	HIS	-	expression tag	UNP P14294
A	657	HIS	-	expression tag	UNP P14294
A	658	HIS	-	expression tag	UNP P14294
A	659	HIS	-	expression tag	UNP P14294
B	654	HIS	-	expression tag	UNP P14294
B	655	HIS	-	expression tag	UNP P14294
B	656	HIS	-	expression tag	UNP P14294
B	657	HIS	-	expression tag	UNP P14294
B	658	HIS	-	expression tag	UNP P14294
B	659	HIS	-	expression tag	UNP P14294

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Cl 1 1	0	0
3	B	1	Total Cl 1 1	0	0

- Molecule 4 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂).



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

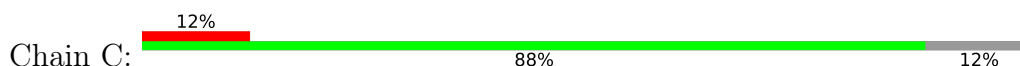
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	4	Total O 4 4	0	0
5	D	6	Total O 6 6	0	0
5	A	53	Total O 53 53	0	0
5	B	79	Total O 79 79	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: 5'-D(*CP*GP*CP*AP*AP*CP*TP*T)-3'

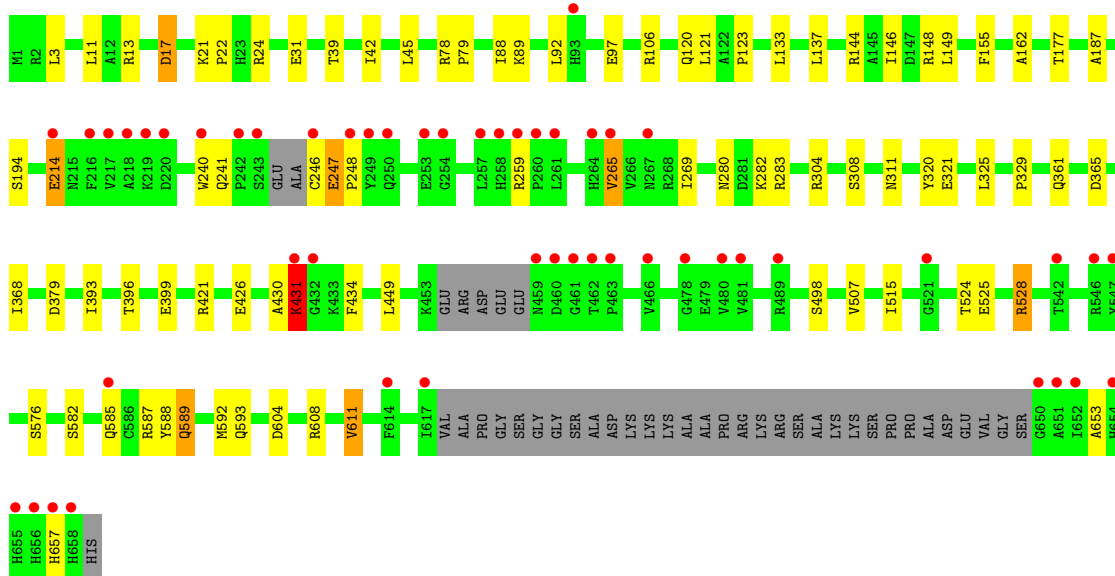
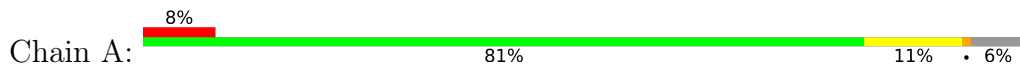


- Molecule 1: 5'-D(*CP*GP*CP*AP*AP*CP*TP*T)-3'

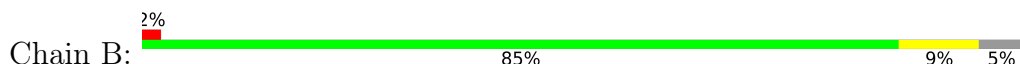


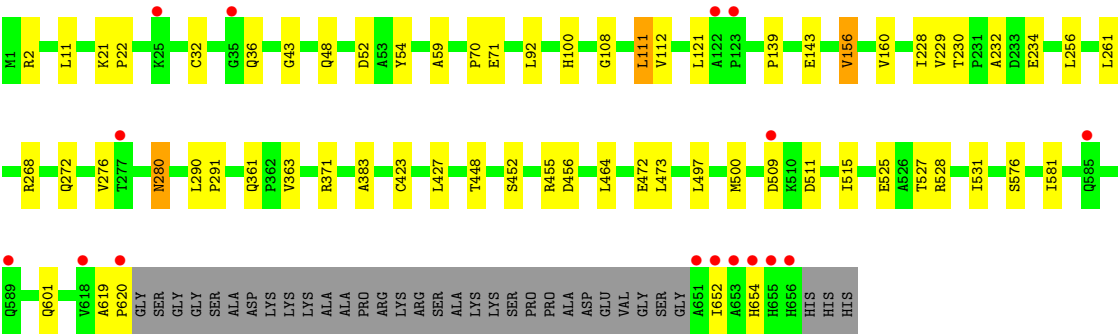
There are no outlier residues recorded for this chain.

- Molecule 2: DNA topoisomerase 3



- Molecule 2: DNA topoisomerase 3





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	101.59Å 101.59Å 453.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.53 – 2.50 29.53 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.2 (29.53-2.50) 97.2 (29.53-2.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.86 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.221 , 0.258 0.212 , 0.246	Depositor DCC
R_{free} test set	4038 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	55.9	Xtriage
Anisotropy	0.254	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10353	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.99% 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: ACY, CL

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	C	0.44	0/154	1.10	0/235
1	D	0.46	0/176	1.12	0/269
2	A	0.58	0/5035	0.84	3/6824 (0.0%)
2	B	0.62	1/5091 (0.0%)	0.86	4/6904 (0.1%)
All	All	0.59	1/10456 (0.0%)	0.86	7/14232 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	156	VAL	CA-CB	5.28	1.57	1.54

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	361	GLN	CA-C-N	6.41	125.91	119.24
2	B	361	GLN	C-N-CA	6.41	125.91	119.24
2	B	464	LEU	CA-C-N	5.81	125.78	120.21
2	B	464	LEU	C-N-CA	5.81	125.78	120.21
2	A	528	ARG	N-CA-C	5.26	117.01	111.28
2	A	241	GLN	CA-C-N	5.20	125.49	119.93
2	A	241	GLN	C-N-CA	5.20	125.49	119.93

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	C	138	0	80	0	0
1	D	158	0	92	0	0
2	A	4927	0	4943	46	0
2	B	4982	0	4995	31	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	B	4	0	3	0	0
5	A	53	0	0	2	0
5	B	79	0	0	1	0
5	C	4	0	0	0	0
5	D	6	0	0	0	0
All	All	10353	0	10113	77	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 4.

All (77) 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)
2:A:396:THR:HG23	2:A:399:GLU:H	1.43	0.83
2:B:230:THR:HG22	2:B:232:ALA:H	1.44	0.82
2:A:308:SER:OG	2:A:311:ASN:HB2	1.88	0.72
2:A:187:ALA:HB2	2:A:611:VAL:HG13	1.74	0.70
2:A:24:ARG:HB3	2:A:31:GLU:HG3	1.76	0.66
2:A:396:THR:HG22	2:A:399:GLU:CG	2.29	0.62
2:B:619:ALA:HB1	2:B:620:PRO:HD2	1.81	0.62
2:A:283:ARG:HE	2:A:421:ARG:HH21	1.47	0.61
2:A:430:ALA:O	2:A:431:LYS:HB3	1.99	0.60
2:B:43:GLY:C	2:B:111:LEU:HD12	2.27	0.60
2:B:525:GLU:OE2	2:B:525:GLU:HA	2.03	0.59
2:A:396:THR:HG22	2:A:399:GLU:HG3	1.86	0.57
2:B:290:LEU:HB3	2:B:291:PRO:HD2	1.87	0.56
2:A:42:ILE:H	2:A:45:LEU:HD21	1.71	0.56
2:B:256:LEU:HD11	2:B:261:LEU:HD23	1.87	0.56
2:B:32:CYS:HB2	2:B:36:GLN:HB2	1.88	0.55
2:A:524:THR:O	2:A:528:ARG:HG3	2.08	0.54
2:B:363:VAL:HG21	2:B:448:THR:HG21	1.90	0.53
2:A:88:ILE:O	2:A:92:LEU:HB2	2.09	0.52
2:A:21:LYS:HB3	2:A:22:PRO:HA	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:587:ARG:NH1	5:A:853:HOH:O	2.42	0.51
2:B:500:MET:HE2	2:B:528:ARG:HG2	1.92	0.51
2:A:361:GLN:HG2	2:A:449:LEU:HD21	1.91	0.51
2:A:265:VAL:O	2:A:269:ILE:HG12	2.11	0.50
2:A:280:ASN:HD21	2:A:282:LYS:HE3	1.77	0.49
2:A:92:LEU:HG	2:A:121:LEU:HD13	1.95	0.48
2:A:396:THR:CG2	2:A:399:GLU:HG3	2.43	0.48
2:B:228:ILE:HG21	2:B:473:LEU:HD23	1.96	0.48
2:A:11:LEU:HD12	2:A:137:LEU:HD21	1.96	0.47
2:A:515:ILE:HD13	2:A:576:SER:HB2	1.95	0.47
2:B:509:ASP:OD1	2:B:509:ASP:C	2.58	0.47
2:A:13:ARG:O	2:A:17:ASP:HB2	2.15	0.47
2:A:396:THR:HG22	2:A:399:GLU:CD	2.40	0.47
2:A:320:TYR:CG	2:A:329:PRO:HD3	2.51	0.46
2:B:455:ARG:HD2	2:B:456:ASP:OD2	2.16	0.45
2:A:106:ARG:HG2	2:A:162:ALA:HB2	1.99	0.45
2:A:246:CYS:SG	2:A:247:GLU:N	2.90	0.45
2:B:100:HIS:NE2	2:B:112:VAL:HB	2.32	0.45
2:B:276:VAL:HA	2:B:427:LEU:HD23	2.00	0.44
2:B:156:VAL:HG11	2:B:654:HIS:HB2	1.99	0.44
2:A:3:LEU:HD11	2:A:39:THR:HG22	1.99	0.44
2:B:160:VAL:HG12	2:B:581:ILE:HG21	1.99	0.44
2:A:146:ILE:O	2:A:149:LEU:HG	2.17	0.44
2:B:108:GLY:HA2	2:B:111:LEU:HD22	1.99	0.44
2:A:365:ASP:HB3	2:A:368:ILE:HD12	1.99	0.44
2:A:604:ASP:O	2:A:608:ARG:HD2	2.17	0.44
2:B:291:PRO:HB2	2:B:383:ALA:CB	2.48	0.44
2:B:515:ILE:HD13	2:B:576:SER:HB2	2.00	0.43
2:A:78:ARG:HA	2:A:79:PRO:HD3	1.89	0.43
2:B:92:LEU:HG	2:B:121:LEU:HD13	2.00	0.43
2:A:247:GLU:H	2:A:248:PRO:CD	2.31	0.43
2:A:588:TYR:CE1	2:A:592:MET:HE3	2.54	0.43
2:B:21:LYS:HE2	2:B:22:PRO:HA	2.01	0.43
2:B:527:THR:O	2:B:531:ILE:HG12	2.19	0.43
2:B:139:PRO:O	2:B:143:GLU:HB2	2.18	0.42
2:B:268:ARG:O	2:B:272:GLN:HG3	2.19	0.42
2:A:177:THR:HG23	2:A:194:SER:HA	2.00	0.42
2:A:133:LEU:HD11	2:A:155:PHE:CZ	2.54	0.42
2:B:54:TYR:CE2	2:B:70:PRO:HB3	2.54	0.42
2:B:228:ILE:CG2	2:B:229:VAL:N	2.82	0.42
2:A:133:LEU:HD11	2:A:155:PHE:HZ	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:653:ALA:O	2:A:657:HIS:ND1	2.49	0.42
2:B:509:ASP:OD1	2:B:511:ASP:N	2.52	0.42
2:B:52:ASP:HB3	2:B:59:ALA:HB2	2.02	0.41
2:B:371:ARG:HG2	5:B:805:HOH:O	2.19	0.41
2:A:240:TRP:HB2	2:A:434:PHE:CE2	2.55	0.41
2:A:308:SER:HG	2:A:311:ASN:HB2	1.85	0.41
2:A:396:THR:CG2	2:A:399:GLU:H	2.22	0.41
2:A:589:GLN:HG3	2:A:593:GLN:HG2	2.02	0.41
2:A:144:ARG:O	2:A:148:ARG:HG3	2.21	0.41
2:A:79:PRO:HD2	5:A:813:HOH:O	2.20	0.41
2:B:230:THR:HB	2:B:234:GLU:HG2	2.03	0.41
2:A:11:LEU:HD12	2:A:137:LEU:CD2	2.51	0.41
2:A:137:LEU:HD12	2:A:321:GLU:HA	2.03	0.41
2:A:325:LEU:HD21	2:A:393:ILE:HD12	2.03	0.41
2:A:78:ARG:HH11	2:A:78:ARG:HG2	1.87	0.40
2:B:280:ASN:O	2:B:423:CYS:HA	2.21	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	
2	A	611/659 (93%)	590 (97%)	15 (2%)	6 (1%)	12	24
2	B	622/659 (94%)	606 (97%)	16 (3%)	0	100	100
All	All	1233/1318 (94%)	1196 (97%)	31 (2%)	6 (0%)	24	43

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	120	GLN
2	A	247	GLU

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Mol	Chain	Res	Type
2	A	379	ASP
2	A	431	LYS
2	A	214	GLU
2	A	123	PRO

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	
2	A	526/555 (95%)	510 (97%)	16 (3%)	36	64
2	B	532/555 (96%)	521 (98%)	11 (2%)	47	74
All	All	1058/1110 (95%)	1031 (97%)	27 (3%)	40	68

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	17	ASP
2	A	89	LYS
2	A	97	GLU
2	A	214	GLU
2	A	259	ARG
2	A	265	VAL
2	A	304	ARG
2	A	426	GLU
2	A	431	LYS
2	A	498	SER
2	A	507	VAL
2	A	525	GLU
2	A	582	SER
2	A	585	GLN
2	A	589	GLN
2	A	611	VAL
2	B	2	ARG
2	B	11	LEU
2	B	48	GLN

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Mol	Chain	Res	Type
2	B	71	GLU
2	B	111	LEU
2	B	280	ASN
2	B	452	SER
2	B	472	GLU
2	B	497	LEU
2	B	601	GLN
2	B	652	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	23	HIS
2	A	127	GLN
2	A	280	ASN
2	A	317	GLN
2	A	370	ASN
2	A	490	HIS
2	A	585	GLN
2	A	589	GLN
2	A	593	GLN
2	A	613	GLN
2	B	74	GLN
2	B	323	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

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	ACY	B	803	-	3,3,3	1.04	0	3,3,3	0.73	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

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 ⓘ

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	C	7/8 (87%)	0.37	1 (14%) 6 5	41, 43, 74, 104	0
1	D	8/8 (100%)	-0.86	0 100 100	34, 39, 54, 63	0
2	A	619/659 (93%)	0.48	51 (8%) 17 15	33, 66, 120, 164	0
2	B	626/659 (94%)	0.05	16 (2%) 57 52	28, 52, 95, 198	0
All	All	1260/1334 (94%)	0.26	68 (5%) 31 27	28, 58, 103, 198	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	652	ILE	11.0
2	A	651	ALA	8.0
2	B	651	ALA	7.5
2	A	216	PHE	7.4
2	B	656	HIS	6.1
2	B	653	ALA	6.0
2	B	122	ALA	5.2
2	A	655	HIS	5.1
2	A	652	ILE	4.8
2	A	617	ILE	4.7
2	A	217	VAL	4.6
2	A	658	HIS	4.5
2	A	650	GLY	4.4
2	B	655	HIS	4.2
2	A	656	HIS	4.2
2	A	547	TYR	4.1
2	A	260	PRO	4.1
2	A	654	HIS	4.1
2	A	246	CYS	3.9
2	B	654	HIS	3.8
1	C	7	DT	3.7

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Mol	Chain	Res	Type	RSRZ
2	A	243	SER	3.6
2	A	461	GLY	3.5
2	A	248	PRO	3.5
2	B	620	PRO	3.5
2	A	93	HIS	3.4
2	A	478	GLY	3.3
2	A	585	GLN	3.2
2	A	257	LEU	3.2
2	A	431	LYS	3.1
2	A	214	GLU	3.1
2	A	261	LEU	3.0
2	A	259	ARG	3.0
2	A	459	ASN	3.0
2	A	462	THR	2.9
2	A	264	HIS	2.9
2	B	277	THR	2.8
2	A	254	GLY	2.7
2	A	460	ASP	2.7
2	A	265	VAL	2.6
2	A	220	ASP	2.6
2	A	657	HIS	2.6
2	A	219	LYS	2.6
2	A	542	THR	2.5
2	A	249	TYR	2.5
2	A	258	HIS	2.4
2	A	466	VAL	2.4
2	A	463	PRO	2.4
2	A	218	ALA	2.4
2	B	25	LYS	2.4
2	B	123	PRO	2.4
2	B	585	GLN	2.3
2	A	481	VAL	2.3
2	A	432	GLY	2.3
2	A	521	GLY	2.3
2	A	489	ARG	2.3
2	B	35	GLY	2.2
2	A	253	GLU	2.2
2	A	250	GLN	2.2
2	A	614	PHE	2.2
2	A	480	VAL	2.1
2	A	242	PRO	2.1
2	B	618	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
2	B	509	ASP	2.1
2	A	546	ARG	2.0
2	A	240	TRP	2.0
2	B	589	GLN	2.0
2	A	267	ASN	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
4	ACY	B	803	4/4	0.84	0.22	45,45,46,46	0
3	CL	B	801	1/1	0.98	0.06	39,39,39,39	0
3	CL	A	800	1/1	0.98	0.07	37,37,37,37	0

6.5 Other polymers [i](#)

There are no such residues in this entry.