Full wwPDB X-ray Structure Validation Report ⓘ

Oct 26, 2026 – 10:24 PM JST

PDB ID	:	9XD8 / pdb_00009xd8
Title	:	Open-Sate Cleavage complex of human Topoisomerase IIA-Beta with oligonucleotide
Authors	:	Ali, M.S.; Zhao, W.; Tang, Y.J.
Deposited on	:	2025-10-27
Resolution	:	3.23 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:

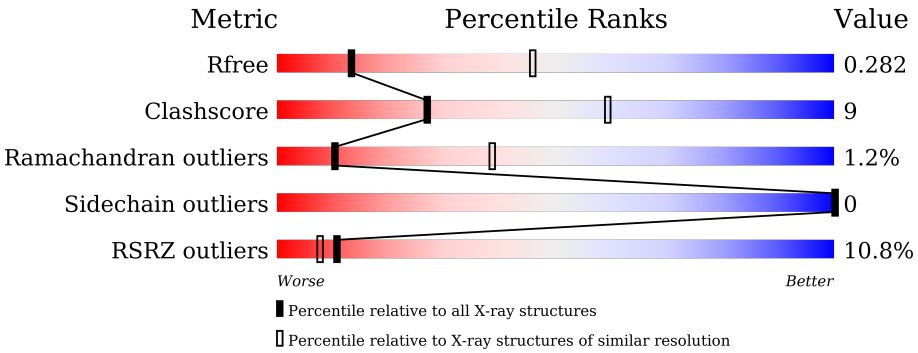
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.23 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	2153 (3.28-3.20)
Clashscore	190562	2275 (3.28-3.20)
Ramachandran outliers	187476	2233 (3.28-3.20)
Sidechain outliers	187428	2232 (3.28-3.20)
RSRZ outliers	180081	2153 (3.28-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	B	757	10% 72% 20% 8%
2	D	10	20% 30% 50% 20%
3	E	8	12% 38% 50% 12%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5701 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 DNA topoisomerase 2-beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	694	Total	C	N	O	S	0	1	0
			5343	3401	910	1009	23			

- Molecule 2 is a DNA chain called DNA (5'-D(P*CP*AP*GP*CP*TP*CP*GP*GP*CP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	10	Total	C	N	O	P	0	0	0
			203	96	36	61	10			

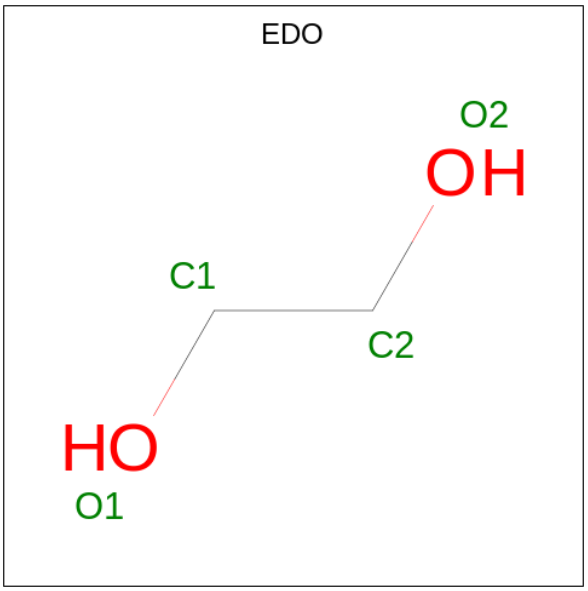
- Molecule 3 is a DNA chain called DNA (5'-D(*AP*GP*CP*CP*GP*AP*GP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	7	Total	C	N	O	P	0	0	0
			144	67	29	41	7			

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	Mg	0	0
			2	2		
4	D	1	Total	Mg	0	0
			1	1		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



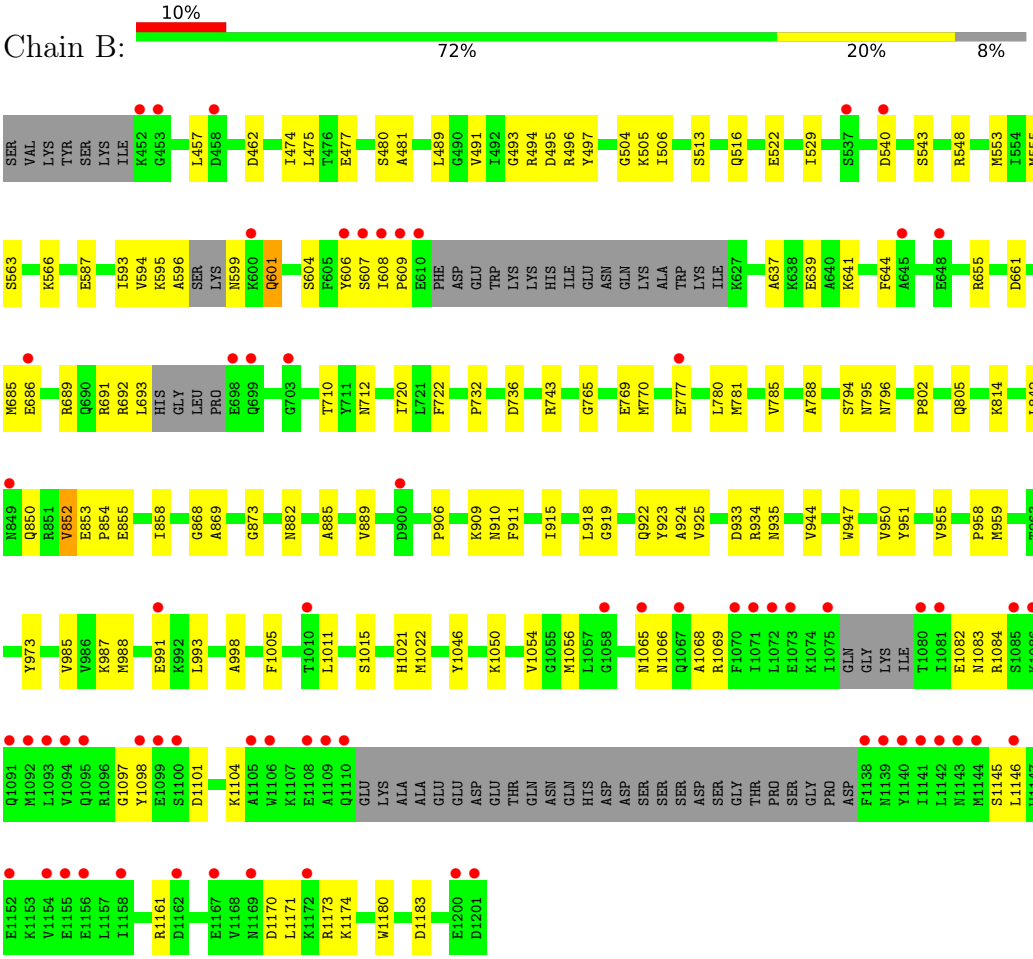
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		

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: DNA topoisomerase 2-beta

Chain B:



• Molecule 2: DNA (5'-D(P*CP*AP*GP*CP*TP*CP*GP*GP*CP*T)-3')

Chain D:



• Molecule 3: DNA (5'-D(*AP*GP*CP*CP*GP*AP*GP*C)-3')

Chain E:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	96.15Å 96.15Å 232.71Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.11 – 3.23 20.11 – 3.23	Depositor EDS
% Data completeness (in resolution range)	99.0 (20.11-3.23) 98.5 (20.11-3.23)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.59 (at 3.22Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487, PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.233 , 0.282 0.233 , 0.282	Depositor DCC
R_{free} test set	1026 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å^2)	-1.6	Xtriage
Anisotropy	-30.681	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 55.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.52$, $< L^2 > = 0.36$	Xtriage
Estimated twinning fraction	0.017 for -h,-k,l	Xtriage
F_o,F_c correlation	0.87	EDS
Total number of atoms	5701	wwPDB-VP
Average B, all atoms (Å^2)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.50% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: EDO, MG

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	B	0.46	0/5447	0.73	0/7375
2	D	1.61	3/226 (1.3%)	1.22	0/346
3	E	1.53	1/161 (0.6%)	1.17	1/246 (0.4%)
All	All	0.60	4/5834 (0.1%)	0.78	1/7967 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	5	DG	C3'-O3'	-7.82	1.33	1.44
2	D	11	DC	C1'-N1	5.79	1.55	1.47
2	D	11	DC	C3'-O3'	5.06	1.50	1.44
2	D	15	DT	C1'-N1	5.02	1.54	1.47

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	E	4	DC	O4'-C4'-C3'	-6.39	102.17	106.00

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	B	5343	0	5104	97	0
2	D	203	0	113	4	0
3	E	144	0	78	2	0
4	B	2	0	0	0	0
4	D	1	0	0	0	0
5	B	8	0	12	0	0
All	All	5701	0	5307	100	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 9.

All (100) 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:852:VAL:HG12	1:B:853:GLU:H	1.38	0.86
1:B:925:VAL:HG11	1:B:944:VAL:HG22	1.69	0.74
1:B:959:MET:HE1	1:B:1005:PHE:HA	1.68	0.74
1:B:1101:ASP:HB3	1:B:1104:LYS:HB2	1.70	0.73
1:B:732:PRO:HG2	1:B:869:ALA:HB1	1.70	0.72
1:B:1021:HIS:CE1	1:B:1022:MET:HG2	2.32	0.65
1:B:457:LEU:HD22	1:B:529:ILE:HG12	1.78	0.64
1:B:555:MET:HE1	1:B:644:PHE:HZ	1.62	0.63
1:B:680:TRP:CE2	1:B:720:ILE:HD12	2.33	0.63
1:B:852:VAL:HG12	1:B:853:GLU:N	2.12	0.63
1:B:593:ILE:HG23	1:B:594:VAL:HG23	1.81	0.62
1:B:686:GLU:OE1	1:B:689:ARG:NH1	2.33	0.61
1:B:873:GLY:O	2:D:16:DC:H2"	2.01	0.60
1:B:1054:VAL:C	1:B:1056:MET:H	2.09	0.60
1:B:991:GLU:H	1:B:991:GLU:CD	2.11	0.58
1:B:1082:GLU:O	1:B:1084:ARG:N	2.38	0.57
1:B:595:LYS:NZ	1:B:639:GLU:OE2	2.38	0.57
1:B:915:ILE:HG12	1:B:925:VAL:HG22	1.86	0.57
1:B:1069:ARG:NH1	1:B:1097:GLY:O	2.38	0.56
1:B:918:LEU:HD11	1:B:924:ALA:HB2	1.87	0.56
1:B:462:ASP:OD2	1:B:496:ARG:NE	2.28	0.56
1:B:845:LEU:HD21	1:B:855:GLU:HB3	1.86	0.55
1:B:474:ILE:HG21	1:B:555:MET:HE3	1.89	0.55
1:B:777:GLU:HB3	1:B:781:MET:HE3	1.89	0.54
1:B:923:TYR:HB2	1:B:1011:LEU:HB3	1.88	0.54
1:B:506:ILE:HG22	1:B:522:GLU:OE1	2.06	0.54
1:B:1065:ASN:OD1	1:B:1161:ARG:HB2	2.08	0.54
1:B:655:ARG:O	1:B:710:THR:HA	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:955:VAL:C	1:B:958:PRO:HD2	2.34	0.53
1:B:988:MET:HE3	1:B:993:LEU:HD13	1.90	0.53
1:B:540:ASP:HB2	1:B:543:SER:H	1.74	0.53
1:B:993:LEU:HD12	1:B:993:LEU:O	2.09	0.52
1:B:604:SER:O	1:B:608:ILE:HG12	2.10	0.52
1:B:765:GLY:HA2	1:B:777:GLU:OE2	2.09	0.52
1:B:844:PHE:HA	1:B:854:PRO:HA	1.92	0.52
1:B:736:ASP:OD1	1:B:743:ARG:NH1	2.43	0.51
1:B:491:VAL:HG13	1:B:494:ARG:HH21	1.74	0.51
1:B:691:ARG:C	1:B:693:LEU:H	2.19	0.51
1:B:1046:TYR:OH	1:B:1183:ASP:OD2	2.21	0.50
1:B:935:ASN:HA	1:B:988:MET:O	2.12	0.50
1:B:1069:ARG:HB2	1:B:1098:TYR:CE2	2.47	0.49
1:B:788:ALA:O	1:B:802:PRO:HG3	2.13	0.49
1:B:637:ALA:O	1:B:641:LYS:HG2	2.13	0.49
1:B:906:PRO:HB2	1:B:915:ILE:HG21	1.95	0.49
1:B:1170:ASP:OD1	1:B:1173:ARG:NH2	2.46	0.48
1:B:796:ASN:ND2	1:B:882:ASN:HB3	2.28	0.48
1:B:477:GLU:HB2	1:B:556:THR:HG22	1.95	0.48
1:B:769:GLU:HG3	1:B:770:MET:N	2.27	0.48
1:B:951:TYR:OH	1:B:1005:PHE:O	2.29	0.48
1:B:852:VAL:HG12	1:B:853:GLU:OE1	2.13	0.48
1:B:601:GLN:HG2	1:B:607:SER:HB2	1.95	0.47
1:B:462:ASP:HB2	1:B:495:ASP:O	2.15	0.47
1:B:850:GLN:O	1:B:852:VAL:HG23	2.15	0.46
1:B:1082:GLU:C	1:B:1084:ARG:H	2.22	0.46
1:B:529:ILE:O	1:B:548:ARG:HD3	2.14	0.46
1:B:560:GLN:HB3	1:B:722:PHE:HA	1.97	0.46
1:B:885:ALA:O	1:B:889:VAL:HG23	2.16	0.46
1:B:493:GLY:O	1:B:497:TYR:HB2	2.16	0.46
1:B:553:MET:HA	1:B:587:GLU:O	2.15	0.46
1:B:933:ASP:C	1:B:934:ARG:O	2.58	0.45
1:B:909:LYS:O	1:B:911:PHE:N	2.50	0.45
1:B:935:ASN:O	1:B:987:LYS:HA	2.16	0.45
2:D:14:DC:H2"	2:D:15:DT:H5'	1.98	0.45
1:B:563:SER:O	1:B:566:LYS:HB2	2.17	0.45
1:B:596:ALA:H	1:B:599:ASN:HA	1.82	0.44
1:B:973:TYR:HA	1:B:985:VAL:O	2.17	0.44
1:B:504:GLY:HA2	3:E:8:DC:H2"	1.99	0.44
1:B:814:LYS:HA	1:B:814:LYS:HD3	1.59	0.44
1:B:947:TRP:HB2	1:B:950:VAL:HG23	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:491:VAL:HA	1:B:494:ARG:NE	2.32	0.44
1:B:604:SER:C	1:B:606:TYR:H	2.25	0.44
1:B:1066:ASN:C	1:B:1068:ALA:H	2.25	0.44
1:B:843:LYS:HD2	1:B:855:GLU:CD	2.43	0.43
1:B:606:TYR:O	1:B:609:PRO:HD2	2.19	0.43
1:B:1054:VAL:O	1:B:1056:MET:N	2.48	0.43
1:B:1145:SER:O	1:B:1148:SER:OG	2.32	0.43
1:B:1050:LYS:HE3	1:B:1171:LEU:O	2.19	0.42
1:B:842:LEU:HD21	1:B:858:ILE:HG22	2.02	0.42
1:B:1046:TYR:CZ	1:B:1180:TRP:HA	2.54	0.42
1:B:685:MET:O	1:B:689:ARG:HG3	2.20	0.42
1:B:1054:VAL:C	1:B:1056:MET:N	2.77	0.42
1:B:780:LEU:HA	1:B:780:LEU:HD12	1.76	0.42
1:B:489:LEU:HD11	1:B:497:TYR:HB3	2.02	0.42
1:B:781:MET:O	1:B:785:VAL:HG23	2.20	0.42
1:B:691:ARG:O	1:B:693:LEU:N	2.53	0.41
2:D:19:DC:H2"	2:D:20:DT:H72	2.01	0.41
1:B:805:GLN:O	1:B:805:GLN:HG3	2.20	0.41
1:B:475:LEU:HD23	1:B:475:LEU:HA	1.72	0.41
1:B:736:ASP:CG	1:B:743:ARG:HH11	2.28	0.41
1:B:794:SER:OG	1:B:795:ASN:N	2.53	0.41
1:B:951:TYR:CD1	1:B:955:VAL:HG21	2.55	0.41
1:B:1170:ASP:O	1:B:1174:LYS:HG3	2.21	0.41
1:B:505:LYS:HG3	3:E:7:DG:H21	1.86	0.40
1:B:513:SER:CB	1:B:516:GLN:HB2	2.50	0.40
1:B:677:ARG:O	1:B:680:TRP:HB3	2.20	0.40
1:B:919:GLY:HA3	1:B:922:GLN:OE1	2.21	0.40
2:D:11:DC:H4'	2:D:12:DA:OP1	2.21	0.40
1:B:661:ASP:OD1	1:B:712:ASN:HB2	2.21	0.40
1:B:480:SER:O	1:B:481:ALA:C	2.65	0.40
1:B:1146:LEU:C	1:B:1148:SER:N	2.79	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	B	681/757 (90%)	608 (89%)	65 (10%)	8 (1%)	10 38

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	601	GLN
1	B	1015	SER
1	B	910	ASN
1	B	692	ARG
1	B	868	GLY
1	B	998	ALA
1	B	1083	ASN
1	B	852	VAL

5.3.2 Protein sidechains (i)

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	B	542/665 (82%)	542 (100%)	0	100 100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	576	ASN
1	B	650	HIS
1	B	712	ASN
1	B	742	GLN
1	B	774	HIS
1	B	840	ASN
1	B	935	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 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	$\#\|Z\| > 2$
5	EDO	B	1303	-	3,3,3	0.53	0	2,2,2	0.28	0
5	EDO	B	1304	-	3,3,3	0.48	0	2,2,2	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
5	EDO	B	1303	-	-	1/1/1/1	-
5	EDO	B	1304	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	1303	EDO	O1-C1-C2-O2

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, 95^{th} 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	B	694/757 (91%)	0.61	74 (10%)	11	8	2, 13, 51, 69	1 (0%)
2	D	10/10 (100%)	0.98	2 (20%)	3	2	4, 8, 45, 70	0
3	E	7/8 (87%)	0.54	1 (14%)	6	5	4, 5, 18, 32	0
All	All	711/775 (91%)	0.61	77 (10%)	11	8	2, 13, 51, 70	1 (0%)

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1091	GLN	4.7
1	B	452	LYS	4.4
1	B	1073	GLU	4.4
1	B	1095	GLN	4.3
1	B	1162	ASP	4.0
1	B	1154	VAL	4.0
1	B	1109	ALA	3.9
1	B	1100	SER	3.8
1	B	1080	THR	3.8
1	B	1141	ILE	3.8
1	B	609	PRO	3.7
1	B	900	ASP	3.7
1	B	1142	LEU	3.6
1	B	1169	ASN	3.6
1	B	1093	LEU	3.5
1	B	1139	ASN	3.4
1	B	1148	SER	3.3
1	B	1152	GLU	3.3
1	B	1088	ASP	3.3
1	B	1065	ASN	3.2
1	B	1072	LEU	3.2
1	B	1151	LYS	3.2
1	B	1087	LYS	3.2

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Mol	Chain	Res	Type	RSRZ
2	D	11	DC	3.1
1	B	1143	ASN	3.0
1	B	991	GLU	3.0
1	B	1089	LEU	2.9
1	B	1099	GLU	2.8
1	B	606	TYR	2.7
1	B	1140	TYR	2.7
1	B	1146	LEU	2.7
1	B	1155	GLU	2.7
1	B	1075	ILE	2.6
1	B	1105	ALA	2.6
1	B	1085	SER	2.6
1	B	1092	MET	2.6
1	B	1150	THR	2.6
1	B	1098	TYR	2.6
1	B	1144	MET	2.6
1	B	610	GLU	2.5
1	B	1086	LYS	2.5
1	B	1071	ILE	2.5
2	D	12	DA	2.5
1	B	849	ASN	2.5
1	B	645	ALA	2.4
1	B	1081	ILE	2.4
1	B	1200	GLU	2.4
1	B	703	GLY	2.4
1	B	1167	GLU	2.4
1	B	1106	TRP	2.4
1	B	1149	LEU	2.4
1	B	607	SER	2.3
1	B	1158	ILE	2.3
1	B	686	GLU	2.3
1	B	1156	GLU	2.3
1	B	1058	GLY	2.3
1	B	699	GLN	2.3
1	B	1110	GLN	2.2
1	B	1108	GLU	2.2
1	B	600	LYS	2.2
1	B	1070	PHE	2.2
1	B	1138	PHE	2.2
1	B	648	GLU	2.2
1	B	1172	LYS	2.1
1	B	967	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	537	SER	2.1
1	B	453	GLY	2.1
1	B	777	GLU	2.1
1	B	540	ASP	2.1
1	B	1201	ASP	2.1
1	B	698	GLU	2.1
1	B	608	ILE	2.0
1	B	458	ASP	2.0
1	B	1067	GLN	2.0
1	B	1094	VAL	2.0
3	E	2	DG	2.0
1	B	1010	THR	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, 95^{th} 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	MG	D	101	1/1	0.89	0.10	17,17,17,17	0
5	EDO	B	1304	4/4	0.92	0.13	13,16,18,23	0
4	MG	B	1302	1/1	0.93	0.10	11,11,11,11	0
5	EDO	B	1303	4/4	0.94	0.10	4,7,8,13	0
4	MG	B	1301	1/1	0.95	0.13	3,3,3,3	0

6.5 Other polymers (i)

There are no such residues in this entry.