



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:03 AM UTC

PDB ID : 5XDC / pdb_00005xdc
Title : Crystal structure of Indole-bound TdsC from Paenibacillus sp. A11-2
Authors : Hino, T.; Hamamoto, H.; Ohshiro, T.; Nagano, S.
Deposited on : 2017-03-28
Resolution : 1.58 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

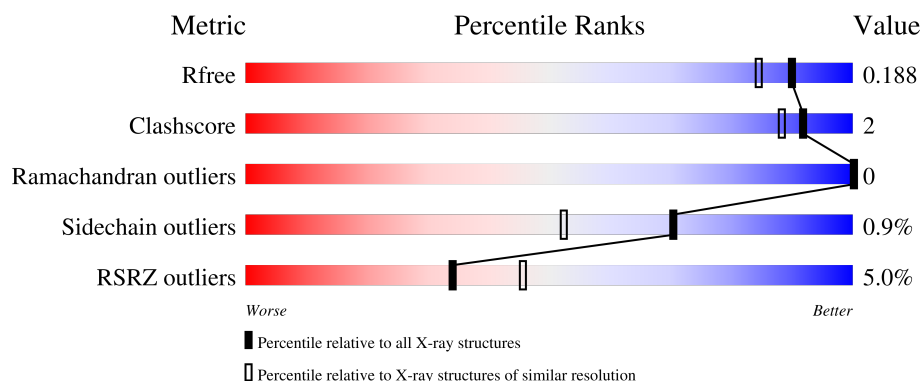
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.58 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	1094 (1.58-1.58)
Clashscore	190562	1105 (1.58-1.58)
Ramachandran outliers	187476	1082 (1.58-1.58)
Sidechain outliers	187428	1081 (1.58-1.58)
RSRZ outliers	180081	1094 (1.58-1.58)

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	414	7% 93% ...
1	B	414	5% 93% ..
1	C	414	4% 93% ..
1	D	414	4% 92% ..

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	IND	A	501	-	X	-	-
2	IND	B	501	-	X	-	-
2	IND	C	502	-	X	-	-
2	IND	D	501	-	X	-	-

2 Entry composition [i](#)

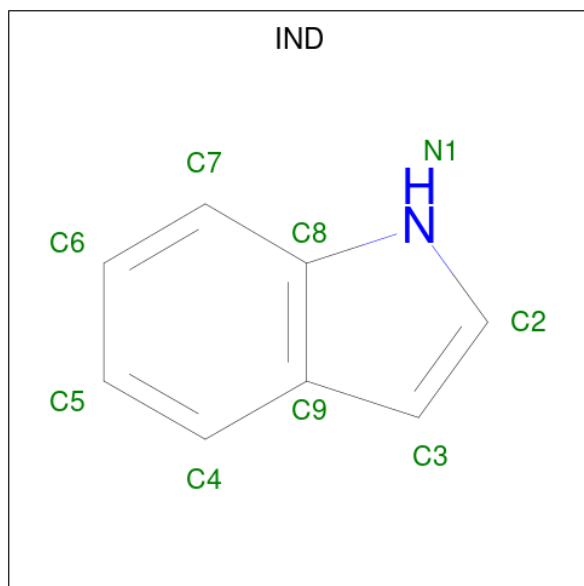
There are 5 unique types of molecules in this entry. The entry contains 26633 atoms, of which 12224 are hydrogens and 0 are deuteriums.

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 Thermophilic dibenzothiophene desulfurization enzyme C.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	401	Total	C	H	N	O	S	0	20	0
			6320	2023	3108	582	603	4			
1	B	402	Total	C	H	N	O	S	0	10	0
			6150	1983	3012	560	591	4			
1	C	401	Total	C	H	N	O	S	0	7	0
			6149	1980	3016	559	590	4			
1	D	396	Total	C	H	N	O	S	0	11	0
			6171	1982	3036	559	590	4			

- Molecule 2 is INDOLE (CCD ID: IND) (formula: C₈H₇N).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	N	0	0
			16	8	7	1		
2	B	1	Total	C	H	N	0	0
			16	8	7	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	H	N	0	0
			16	8	7	1		
2	D	1	Total	C	H	N	0	0
			16	8	7	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	H	O	0	0
			14	3	8	3		
4	C	1	Total	C	H	O	0	0
			14	3	8	3		
4	D	1	Total	C	H	O	0	0
			14	3	8	3		

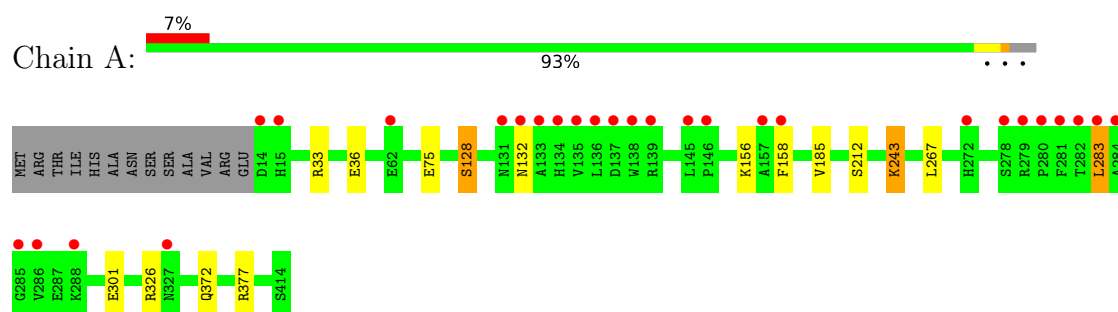
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	373	Total 373	O 373	0	0
5	B	421	Total 421	O 421	0	0
5	C	406	Total 406	O 406	0	0
5	D	457	Total 457	O 457	0	0

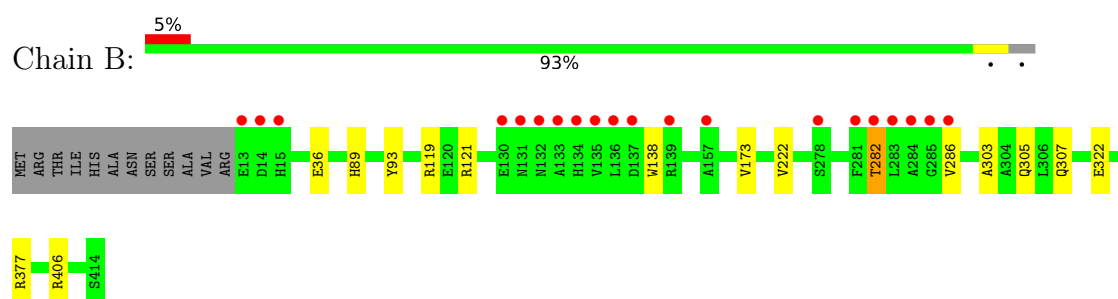
3 Residue-property plots [i](#)

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.

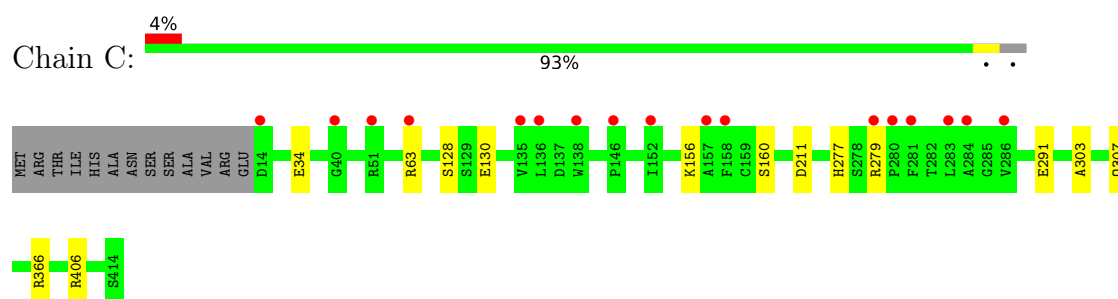
- Molecule 1: Thermophilic dibenzothiophene desulfurization enzyme C



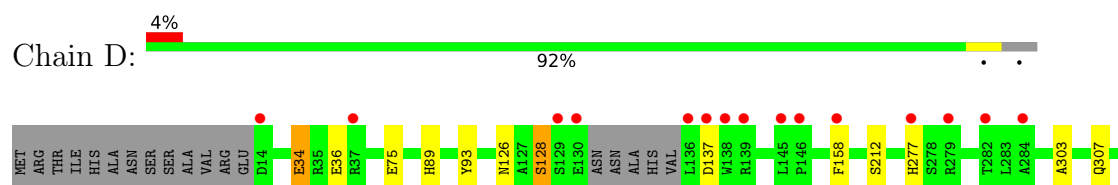
- Molecule 1: Thermophilic dibenzothiophene desulfurization enzyme C



- Molecule 1: Thermophilic dibenzothiophene desulfurization enzyme C



- Molecule 1: Thermophilic dibenzothiophene desulfurization enzyme C





4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	100.83Å 100.83Å 424.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.49 – 1.58 47.49 – 1.58	Depositor EDS
% Data completeness (in resolution range)	88.6 (47.49-1.58) 84.1 (47.49-1.58)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 1.58Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155)	Depositor
R, R_{free}	0.165 , 0.186 0.167 , 0.188	Depositor DCC
R_{free} test set	14152 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	18.6	Xtriage
Anisotropy	0.006	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 49.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	26633	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.28	0/3367	0.47	0/4576
1	B	0.29	0/3244	0.47	0/4417
1	C	0.28	0/3235	0.47	0/4403
1	D	0.28	0/3252	0.46	0/4420
All	All	0.28	0/13098	0.47	0/17816

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3212	3108	3035	11	0
1	B	3138	3012	2982	10	0
1	C	3133	3016	2993	11	0
1	D	3135	3036	2998	11	0
2	A	9	7	7	0	0
2	B	9	7	7	0	0
2	C	9	7	7	3	0
2	D	9	7	7	0	0
3	A	20	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	15	0	0	0	0
3	C	25	0	0	0	0
3	D	20	0	0	0	0
4	B	6	8	8	0	0
4	C	6	8	8	0	0
4	D	6	8	8	0	0
5	A	373	0	0	1	0
5	B	421	0	0	4	1
5	C	406	0	0	3	0
5	D	457	0	0	3	1
All	All	14409	12224	12060	40	1

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 2.

All (40) 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:75[B]:GLU:OE2	5:A:601:HOH:O	1.95	0.82
1:C:277:HIS:O	1:C:366:ARG:NH2	2.12	0.82
1:C:406:ARG:NH2	5:C:601:HOH:O	2.14	0.79
1:A:132[A]:ASN:OD1	1:B:282:THR:OG1	2.06	0.73
1:C:160:SER:OG	2:C:502:IND:H2	1.92	0.70
1:D:406:ARG:NE	5:D:607:HOH:O	2.33	0.61
1:D:137:ASP:OD1	5:D:601:HOH:O	2.18	0.54
1:B:406:ARG:NH2	1:C:291:GLU:OE1	2.37	0.53
1:C:303:ALA:O	1:C:307:GLN:HG2	2.09	0.52
1:A:36:GLU:HG2	1:A:377:ARG:HD3	1.93	0.50
1:B:119:ARG:NH2	5:B:603:HOH:O	2.26	0.49
1:D:36:GLU:HG2	1:D:377:ARG:HD3	1.94	0.48
1:C:130:GLU:OE1	2:C:502:IND:H3	2.13	0.48
1:C:211:ASP:OD2	5:C:602:HOH:O	2.20	0.47
1:C:34:GLU:HG2	5:C:825:HOH:O	2.15	0.47
1:C:160:SER:HG	2:C:502:IND:H2	1.80	0.46
1:B:36:GLU:HG2	1:B:377:ARG:HD3	1.98	0.46
1:B:89:HIS:CE1	1:B:93:TYR:CE1	3.05	0.46
1:D:34:GLU:H	1:D:34:GLU:CD	2.24	0.46
1:C:128:SER:HB3	1:C:156:LYS:HG2	1.99	0.45
1:B:322:GLU:HG2	5:B:876:HOH:O	2.17	0.44
1:A:158:PHE:HA	1:A:212:SER:O	2.18	0.44
1:A:267:LEU:HG	1:D:402:PHE:CE2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:121[A]:ARG:NE	5:B:616:HOH:O	2.50	0.44
1:D:75:GLU:OE2	5:D:602:HOH:O	2.21	0.44
1:B:303:ALA:O	1:B:307:GLN:HG2	2.18	0.43
1:B:138:TRP:CE3	1:B:173:VAL:HG13	2.53	0.43
1:D:277:HIS:O	1:D:366:ARG:NH2	2.51	0.42
1:A:128:SER:HB3	1:A:156:LYS:HG2	2.02	0.42
1:D:303:ALA:O	1:D:307:GLN:HG2	2.18	0.42
1:D:158:PHE:HA	1:D:212:SER:O	2.20	0.42
1:D:126:ASN:OD1	1:D:128:SER:HA	2.20	0.42
1:D:89:HIS:CE1	1:D:93[A]:TYR:CE1	3.08	0.41
1:C:279:ARG:HG3	1:C:366:ARG:NH2	2.35	0.41
1:A:243:LYS:HE3	1:A:326:ARG:O	2.21	0.41
1:B:305:GLN:NE2	5:B:605:HOH:O	2.38	0.41
1:A:33[A]:ARG:NH1	1:A:372[A]:GLN:HG3	2.36	0.41
1:A:283:LEU:HD23	1:A:283:LEU:N	2.37	0.40

All (1) 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:B:951:HOH:O	5:D:1005:HOH:O[1_455]	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	419/414 (101%)	412 (98%)	7 (2%)	0	100	100
1	B	410/414 (99%)	405 (99%)	5 (1%)	0	100	100
1	C	406/414 (98%)	402 (99%)	4 (1%)	0	100	100
1	D	403/414 (97%)	398 (99%)	5 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1638/1656 (99%)	1617 (99%)	21 (1%)	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	332/331 (100%)	327 (98%)	5 (2%)	57	31
1	B	316/331 (96%)	313 (99%)	3 (1%)	70	52
1	C	318/331 (96%)	317 (100%)	1 (0%)	86	79
1	D	321/331 (97%)	319 (99%)	2 (1%)	78	65
All	All	1287/1324 (97%)	1276 (99%)	11 (1%)	70	52

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

Mol	Chain	Res	Type
1	A	128	SER
1	A	185	VAL
1	A	243	LYS
1	A	283	LEU
1	A	301	GLU
1	B	222	VAL
1	B	282	THR
1	B	286	VAL
1	C	63	ARG
1	D	34	GLU
1	D	128	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	340	GLN

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Mol	Chain	Res	Type
1	A	405	HIS
1	B	328	HIS
1	C	134	HIS
1	C	340	GLN
1	D	89	HIS
1	D	328	HIS
1	D	340	GLN
1	D	405	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](#)

23 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	$\# Z > 2$
3	SO4	D	505	-	4,4,4	0.27	0	6,6,6	0.12	0
2	IND	D	501	-	10,10,10	2.61	5 (50%)	13,13,13	4.66	7 (53%)
4	GOL	B	504	-	5,5,5	0.53	0	5,5,5	0.39	0
3	SO4	C	507	-	4,4,4	0.33	0	6,6,6	0.21	0
3	SO4	C	503	-	4,4,4	0.25	0	6,6,6	0.26	0
3	SO4	B	503	-	4,4,4	0.23	0	6,6,6	0.24	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	D	502	-	4,4,4	0.23	0	6,6,6	0.36	0
3	SO4	A	503	-	4,4,4	0.23	0	6,6,6	0.22	0
3	SO4	B	502	-	4,4,4	0.25	0	6,6,6	0.27	0
3	SO4	C	501	-	4,4,4	0.24	0	6,6,6	0.13	0
3	SO4	D	504	-	4,4,4	0.22	0	6,6,6	0.05	0
3	SO4	A	504	-	4,4,4	0.24	0	6,6,6	0.16	0
3	SO4	A	502	-	4,4,4	0.23	0	6,6,6	0.17	0
4	GOL	C	506	-	5,5,5	0.56	0	5,5,5	0.24	0
3	SO4	A	505	-	4,4,4	0.23	0	6,6,6	0.19	0
3	SO4	B	505	-	4,4,4	0.24	0	6,6,6	0.17	0
2	IND	B	501	-	10,10,10	2.58	5 (50%)	13,13,13	4.71	7 (53%)
4	GOL	D	503	-	5,5,5	0.42	0	5,5,5	0.23	0
3	SO4	C	504	-	4,4,4	0.23	0	6,6,6	0.26	0
3	SO4	D	506	-	4,4,4	0.23	0	6,6,6	0.15	0
2	IND	A	501	-	10,10,10	2.60	5 (50%)	13,13,13	4.72	7 (53%)
2	IND	C	502	-	10,10,10	2.63	4 (40%)	13,13,13	4.55	7 (53%)
3	SO4	C	505	-	4,4,4	0.23	0	6,6,6	0.21	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	D	503	-	-	0/4/4/4	-
2	IND	D	501	-	-	-	0/2/2/2
2	IND	A	501	-	-	-	0/2/2/2
2	IND	C	502	-	-	-	0/2/2/2
4	GOL	B	504	-	-	0/4/4/4	-
4	GOL	C	506	-	-	0/4/4/4	-
2	IND	B	501	-	-	-	0/2/2/2

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	502	IND	C9-C3	-4.47	1.30	1.43
2	A	501	IND	C9-C3	-4.37	1.31	1.43
2	D	501	IND	C9-C3	-4.37	1.31	1.43
2	A	501	IND	C2-C3	-4.35	1.28	1.36
2	D	501	IND	C2-C3	-4.29	1.28	1.36
2	B	501	IND	C9-C3	-4.28	1.31	1.43
2	B	501	IND	C2-C3	-4.20	1.28	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	502	IND	C2-C3	-4.13	1.28	1.36
2	C	502	IND	C8-N1	3.64	1.43	1.37
2	B	501	IND	C8-N1	3.55	1.43	1.37
2	A	501	IND	C8-N1	3.44	1.43	1.37
2	D	501	IND	C8-N1	3.19	1.42	1.37
2	C	502	IND	C2-N1	-2.97	1.32	1.38
2	D	501	IND	C2-N1	-2.91	1.32	1.38
2	A	501	IND	C2-N1	-2.68	1.33	1.38
2	B	501	IND	C2-N1	-2.53	1.33	1.38
2	D	501	IND	C9-C8	2.20	1.44	1.41
2	A	501	IND	C9-C8	2.08	1.44	1.41
2	B	501	IND	C9-C8	2.08	1.44	1.41

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	IND	C9-C3-C2	11.01	120.48	107.05
2	B	501	IND	C9-C3-C2	10.90	120.34	107.05
2	C	502	IND	C9-C3-C2	10.84	120.27	107.05
2	D	501	IND	C9-C3-C2	10.82	120.25	107.05
2	C	502	IND	C8-C9-C3	-7.07	102.12	106.59
2	D	501	IND	C8-C9-C3	-6.87	102.24	106.59
2	A	501	IND	C8-C9-C3	-6.72	102.34	106.59
2	B	501	IND	C3-C2-N1	-6.62	102.61	110.25
2	A	501	IND	C3-C2-N1	-6.46	102.78	110.25
2	B	501	IND	C9-C8-N1	-6.35	102.44	107.53
2	D	501	IND	C3-C2-N1	-6.29	102.98	110.25
2	B	501	IND	C8-C9-C3	-6.28	102.62	106.59
2	A	501	IND	C9-C8-N1	-6.08	102.66	107.53
2	D	501	IND	C9-C8-N1	-5.99	102.73	107.53
2	C	502	IND	C3-C2-N1	-5.95	103.38	110.25
2	C	502	IND	C9-C8-N1	-5.30	103.28	107.53
2	B	501	IND	C7-C8-N1	4.36	137.73	130.74
2	A	501	IND	C7-C8-N1	4.34	137.69	130.74
2	B	501	IND	C8-N1-C2	4.22	112.00	108.91
2	C	502	IND	C7-C8-N1	4.15	137.39	130.74
2	D	501	IND	C7-C8-N1	4.10	137.31	130.74
2	D	501	IND	C8-N1-C2	3.94	111.79	108.91
2	A	501	IND	C8-N1-C2	3.86	111.73	108.91
2	C	502	IND	C7-C8-C9	-3.10	119.32	122.13
2	C	502	IND	C8-N1-C2	2.79	110.95	108.91
2	A	501	IND	C7-C8-C9	-2.74	119.65	122.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	IND	C7-C8-C9	-2.55	119.83	122.13
2	D	501	IND	C7-C8-C9	-2.40	119.96	122.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	502	IND	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	401/414 (96%)	0.13	28 (6%)	22 32	11, 22, 53, 80	11 (2%)
1	B	402/414 (97%)	-0.13	20 (4%)	34 46	8, 20, 46, 90	6 (1%)
1	C	401/414 (96%)	-0.04	17 (4%)	40 52	9, 23, 42, 75	4 (0%)
1	D	396/414 (95%)	-0.23	15 (3%)	44 55	8, 21, 40, 75	6 (1%)
All	All	1600/1656 (96%)	-0.07	80 (5%)	34 46	8, 21, 46, 90	27 (1%)

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	136	LEU	7.4
1	B	283	LEU	6.2
1	A	283	LEU	5.6
1	A	282	THR	5.5
1	D	138	TRP	5.2
1	A	281	PHE	5.0
1	C	283	LEU	4.7
1	A	286	VAL	4.6
1	A	158	PHE	4.5
1	B	135	VAL	4.4
1	B	132	ASN	4.4
1	B	282	THR	4.4
1	B	136	LEU	4.3
1	A	14	ASP	4.3
1	A	284	ALA	4.2
1	B	281	PHE	4.1
1	A	146	PRO	4.0
1	A	157	ALA	3.9
1	C	284	ALA	3.6
1	B	13	GLU	3.6
1	D	137	ASP	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	133	ALA	3.5
1	C	157	ALA	3.5
1	B	131	ASN	3.5
1	A	138	TRP	3.4
1	A	279	ARG	3.4
1	A	139	ARG	3.3
1	C	279	ARG	3.3
1	B	15	HIS	3.3
1	A	15	HIS	3.2
1	B	284	ALA	3.2
1	B	139	ARG	3.2
1	B	286	VAL	3.2
1	B	278	SER	3.2
1	A	133[A]	ALA	3.1
1	D	130	GLU	3.1
1	A	327[A]	ASN	3.0
1	A	280	PRO	3.0
1	A	134	HIS	2.9
1	C	136	LEU	2.9
1	C	281	PHE	2.8
1	A	62[A]	GLU	2.8
1	C	51[A]	ARG	2.8
1	D	146	PRO	2.7
1	B	134	HIS	2.7
1	A	136	LEU	2.6
1	A	272[A]	HIS	2.6
1	A	285	GLY	2.6
1	A	137	ASP	2.5
1	C	14	ASP	2.5
1	D	37	ARG	2.5
1	A	135	VAL	2.5
1	C	286	VAL	2.5
1	A	132[A]	ASN	2.5
1	A	131[A]	ASN	2.5
1	A	278	SER	2.4
1	C	280	PRO	2.4
1	C	40	GLY	2.4
1	D	145	LEU	2.4
1	C	158	PHE	2.4
1	C	138	TRP	2.3
1	B	285	GLY	2.3
1	B	137	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	282	THR	2.3
1	D	279	ARG	2.3
1	A	288	LYS	2.3
1	B	130	GLU	2.2
1	D	129	SER	2.2
1	A	145	LEU	2.2
1	D	139	ARG	2.2
1	D	284	ALA	2.2
1	B	14	ASP	2.2
1	D	158	PHE	2.2
1	D	14	ASP	2.1
1	D	277	HIS	2.1
1	C	135	VAL	2.1
1	B	157	ALA	2.1
1	C	152	ILE	2.1
1	C	63	ARG	2.0
1	C	146	PRO	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
3	SO4	C	507	5/5	0.74	0.17	34,45,69,71	0
3	SO4	D	506	5/5	0.84	0.14	37,57,71,72	0
3	SO4	A	504	5/5	0.85	0.13	36,58,61,72	0
3	SO4	C	501	5/5	0.86	0.11	52,54,97,106	0
3	SO4	C	505	5/5	0.88	0.10	41,41,56,57	0
2	IND	A	501	9/9	0.88	0.15	32,49,57,63	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	IND	D	501	9/9	0.88	0.13	32,38,48,52	0
2	IND	C	502	9/9	0.89	0.11	27,34,44,53	0
3	SO4	D	505	5/5	0.89	0.11	43,48,71,78	0
3	SO4	A	505	5/5	0.89	0.11	34,45,64,73	0
3	SO4	B	503	5/5	0.90	0.12	49,55,61,80	0
3	SO4	D	504	5/5	0.91	0.10	41,45,61,72	0
3	SO4	B	505	5/5	0.91	0.09	37,42,63,79	0
2	IND	B	501	9/9	0.91	0.11	26,34,42,50	0
4	GOL	C	506	6/6	0.94	0.08	25,30,36,41	0
3	SO4	B	502	5/5	0.95	0.08	35,35,44,49	0
3	SO4	A	503	5/5	0.95	0.09	35,40,43,46	0
3	SO4	C	503	5/5	0.96	0.07	31,32,34,49	0
4	GOL	B	504	6/6	0.96	0.07	19,29,35,35	0
3	SO4	D	502	5/5	0.96	0.08	29,31,39,39	0
3	SO4	A	502	5/5	0.97	0.07	35,37,38,47	0
3	SO4	C	504	5/5	0.97	0.07	27,32,39,53	0
4	GOL	D	503	6/6	0.98	0.04	20,24,26,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.