



wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 16, 2024 – 09:20 AM EDT

PDB ID : 5AI4
Title : ligand complex structure of soluble epoxide hydrolase
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Deposited on : 2015-02-12
Resolution : 1.93 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

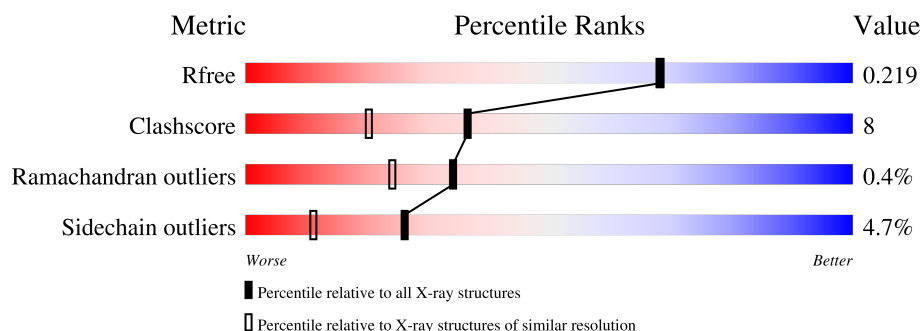
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 1.93 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}	130704	4310 (1.96-1.92)
Clashscore	141614	1023 (1.94-1.94)
Ramachandran outliers	138981	1007 (1.94-1.94)
Sidechain outliers	138945	1007 (1.94-1.94)

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	549	

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	DMS	A	1552	-	-	X	-
4	4VY	A	1555	-	X	-	-
4	4VY	A	1556	-	X	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4887 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.

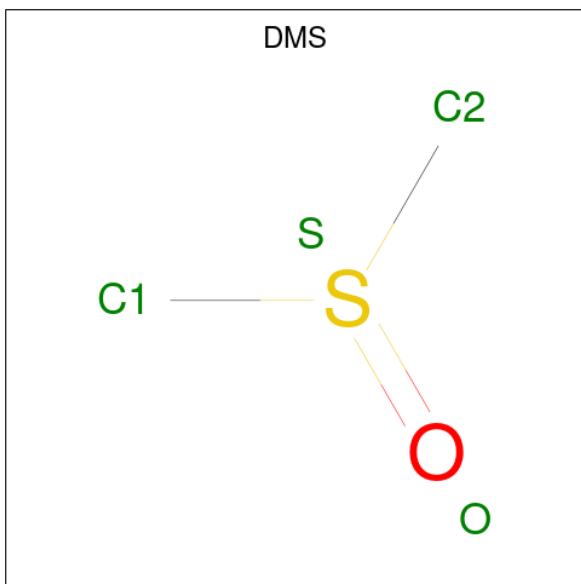
- Molecule 1 is a protein called BIFUNCTIONAL EPOXIDE HYDROLASE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	547	Total	C	N	O	S	0	2	1
			4346	2787	732	790	37			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP P34913

- Molecule 2 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C₂H₆OS).



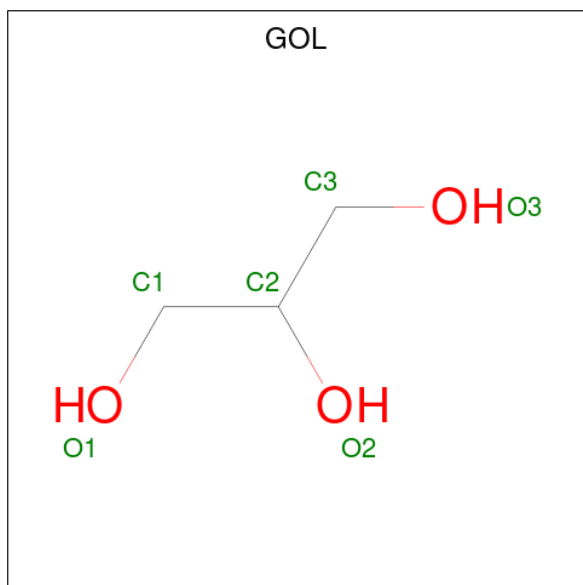
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	S	0	0
			4	2	1	1		
2	A	1	Total	C	O	S	0	0
			4	2	1	1		
2	A	1	Total	C	O	S	0	0
			4	2	1	1		

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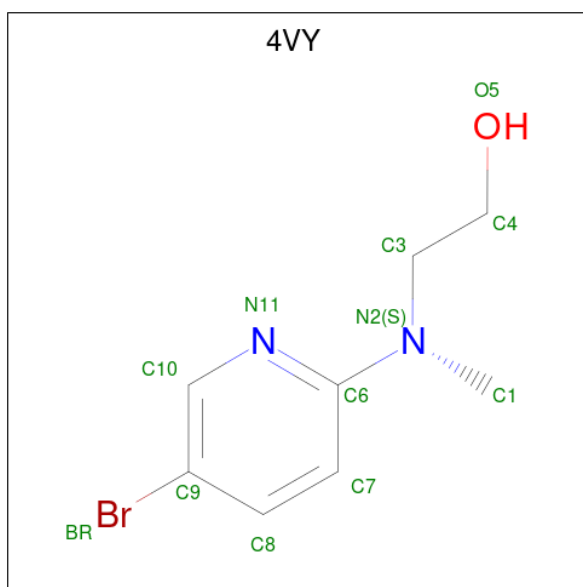
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	S	0	0
			4	2	1	1		
2	A	1	Total	C	O	S	0	0
			4	2	1	1		
2	A	1	Total	C	O	S	0	0
			4	2	1	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 4 is 2-[(5-BROMO-2-PYRIDYL)-METHYL-AMINO]ETHANOL (three-letter code: 4VY) (formula: $C_8H_{11}BrN_2O$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	Br	C	N	O	0	0
			12	1	8	2	1		
4	A	1	Total	Br	C	N	O	0	0
			12	1	8	2	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	487	Total	O	0	0
			487	487		

● Molecule 1: BIFUNCTIONAL EPOXIDE HYDROLASE 2

C477	P246	GLY
L480	R247	MET
L484	V248	T2
	R249	L3
		R4
K495	P259	
L499	G266	V19
	P267	F20
M503	P268	G21
		V22
		L23
	Q277	G24
W510		R25
F511	D292	
		L36
R516	F331	L37
W525	D335	K43
F526		
Q527	V341	P65
Q536	R353	R72
	A354	K73
	V355	
		N85
A546	P364	
R547		F92
N548	M369	
	S370	A95
	P371	
	L372	R99
	E373	
		N102
	K376	R103
	A377	P104
	N378	
	P379	H146
	V380	
		Q164
	Y383	
	Q384	D184
	L385	
	Y386	L190
		K191
	A417	R194
	S418	D195
	M419	L196
	H420	
	K421	V199
	V422	
		D207
	L428	
	F429	E214
	V430	R215
	M431	V216
	S432	
	N472	I219
	W473	
		V215

4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	92.48Å 92.48Å 244.93Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	80.09 – 1.93 76.13 – 1.93	Depositor EDS
% Data completeness (in resolution range)	(Not available) (80.09-1.93) 99.4 (76.13-1.93)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 1.92Å)	Xtriage
Refinement program	BUSTER 2.11.1	Depositor
R, R_{free}	0.181 , 0.208 0.191 , 0.219	Depositor DCC
R_{free} test set	2387 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	26.7	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 57.0	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	0.90	EDS
Total number of atoms	4887	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 5.29% 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: GOL, DMS, 4VY

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	0.47	0/4453	0.67	1/6035 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	292	ASP	N-CA-C	-5.44	96.31	111.00

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	4346	0	4346	65	0
2	A	24	0	36	6	0
3	A	6	0	8	1	0
4	A	24	0	22	10	0
5	A	487	0	0	10	2
All	All	4887	0	4412	70	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 8.

The worst 5 of 70 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:370:SER:HB2	1:A:373:GLU:HG3	1.39	1.01
1:A:547:ARG:HG2	1:A:547:ARG:HH11	1.26	1.00
1:A:472:ASN:HB3	3:A:1554:GOL:H31	1.45	0.99
1:A:331:PHE:HB3	1:A:341:VAL:HG22	1.46	0.94
1:A:383:TYR:HA	1:A:422:VAL:HG21	1.54	0.86

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:2052:HOH:O	5:A:2314:HOH:O[10_665]	1.60	0.60
5:A:2064:HOH:O	5:A:2183:HOH:O[10_665]	2.13	0.07

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	547/549 (100%)	534 (98%)	11 (2%)	2 (0%)	34 24

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	547	ARG
1	A	268	PRO

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	A	474/474 (100%)	452 (95%)	22 (5%)	27 12

5 of 22 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	419	MET
1	A	432	SER
1	A	430	VAL
1	A	480	LEU
1	A	102	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	102	ASN
1	A	146	HIS
1	A	204	GLN
1	A	384	GLN
1	A	452	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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

9 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	DMS	A	1549	-	3,3,3	0.90	0	3,3,3	0.46	0
3	GOL	A	1554	-	5,5,5	0.62	0	5,5,5	1.02	0
4	4VY	A	1556	-	12,12,12	1.81	3 (25%)	15,15,15	5.05	9 (60%)
4	4VY	A	1555	-	12,12,12	2.26	5 (41%)	15,15,15	4.33	6 (40%)
2	DMS	A	1552	-	3,3,3	0.65	0	3,3,3	1.35	1 (33%)
2	DMS	A	1550	-	3,3,3	0.89	0	3,3,3	0.50	0
2	DMS	A	1553	-	3,3,3	0.82	0	3,3,3	0.73	0
2	DMS	A	1548	-	3,3,3	1.02	0	3,3,3	0.12	0
2	DMS	A	1551	-	3,3,3	0.83	0	3,3,3	0.62	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
3	GOL	A	1554	-	-	4/4/4/4	-
4	4VY	A	1556	-	-	6/7/7/7	0/1/1/1
4	4VY	A	1555	-	-	3/7/7/7	0/1/1/1

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1555	4VY	C6-N11	3.80	1.42	1.34
4	A	1555	4VY	C7-C6	3.77	1.48	1.39
4	A	1556	4VY	C6-N11	3.53	1.41	1.34
4	A	1555	4VY	C3-C4	-3.02	1.39	1.51
4	A	1556	4VY	BR-C9	-2.96	1.84	1.90

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1555	4VY	C3-N2-C6	-14.40	105.79	120.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1556	4VY	C3-N2-C6	-13.61	106.62	120.94
4	A	1556	4VY	N11-C6-N2	8.87	127.63	116.39
4	A	1556	4VY	C7-C6-N2	-6.42	111.45	121.97
4	A	1555	4VY	C1-N2-C3	-5.46	100.57	115.42

There are no chirality outliers.

5 of 13 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1555	4VY	C4-C3-N2-C1
4	A	1555	4VY	C7-C6-N2-C3
4	A	1555	4VY	N11-C6-N2-C3
4	A	1556	4VY	C4-C3-N2-C1
4	A	1556	4VY	C4-C3-N2-C6

There are no ring outliers.

4 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1554	GOL	1	0
4	A	1556	4VY	9	0
4	A	1555	4VY	3	0
2	A	1552	DMS	6	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.