



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:18 PM UTC

PDB ID : 3ARY / pdb_00003ary
Title : Crystal Structure Analysis of Chitinase A from *Vibrio harveyi* with novel inhibitors - complex structure with 2-(imidazolin-2-yl)-5-isothiocyantobenzofuran
Authors : Pantoom, S.; Vetter, I.R.; Prinz, H.; Suginta, W.
Deposited on : 2010-12-09
Resolution : 1.35 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

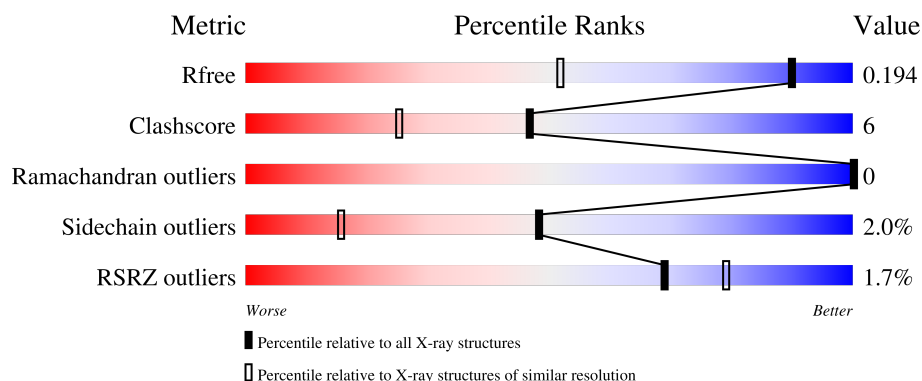
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

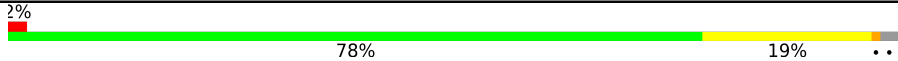
The reported resolution of this entry is 1.35 Å.

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	1216 (1.36-1.36)
Clashscore	190562	1232 (1.36-1.36)
Ramachandran outliers	187476	1220 (1.36-1.36)
Sidechain outliers	187428	1220 (1.36-1.36)
RSRZ outliers	180081	1214 (1.36-1.36)

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	584	

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	I5I	A	607	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5560 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.

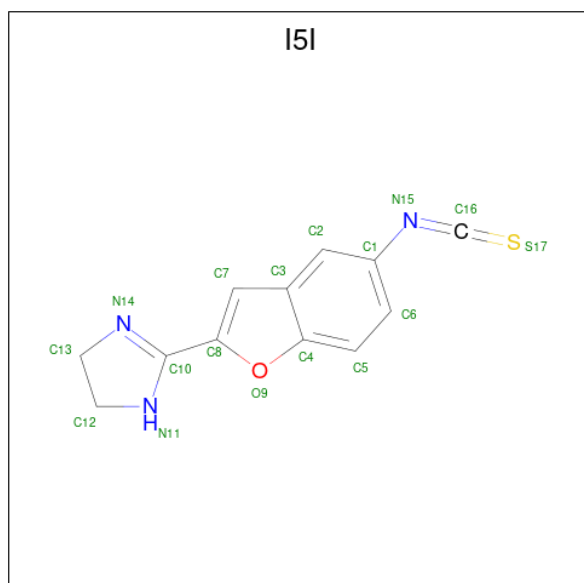
- Molecule 1 is a protein called Chitinase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	575	4551	2903	725	895	28	0	28	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	598	ARG	-	expression tag	UNP Q9AMP1
A	599	SER	-	expression tag	UNP Q9AMP1
A	600	HIS	-	expression tag	UNP Q9AMP1
A	601	HIS	-	expression tag	UNP Q9AMP1
A	602	HIS	-	expression tag	UNP Q9AMP1
A	603	HIS	-	expression tag	UNP Q9AMP1
A	604	HIS	-	expression tag	UNP Q9AMP1
A	605	HIS	-	expression tag	UNP Q9AMP1

- Molecule 2 is 2-(5-isothiocyanato-1-benzofuran-2-yl)-4,5-dihydro-1H-imidazole (CCD ID: I5I) (formula: C₁₂H₉N₃OS).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			17	12	3	1	1		
2	A	1	Total	C	N	O	S	0	0
			17	12	3	1	1		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	1
			12	6	6		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		

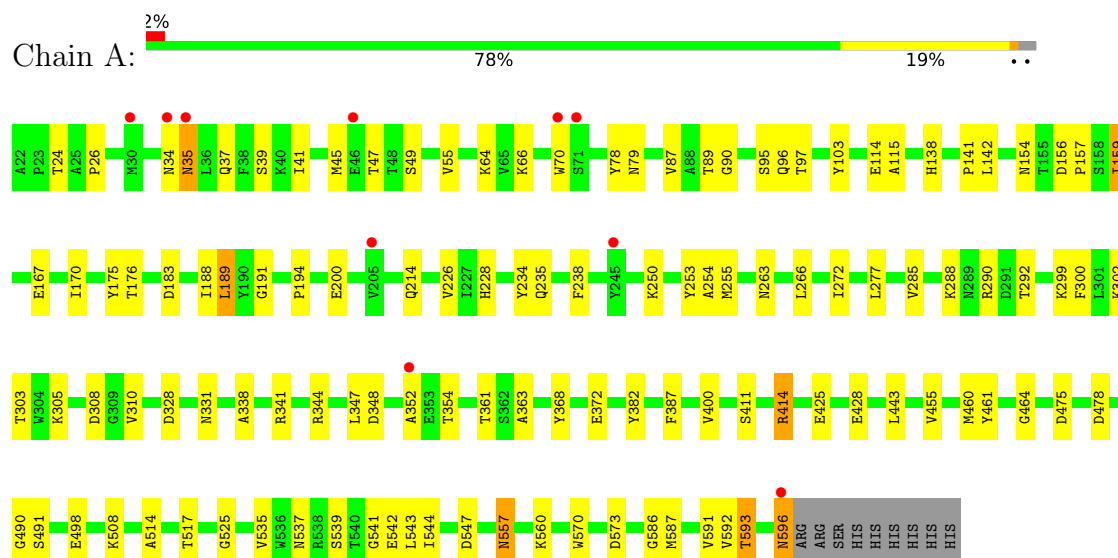
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	951	Total	O	0	0
			951	951		

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: Chitinase A



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	65.08Å 51.03Å 93.27Å 90.00° 99.55° 90.00°	Depositor
Resolution (Å)	19.97 – 1.35 19.97 – 1.35	Depositor EDS
% Data completeness (in resolution range)	97.4 (19.97-1.35) 97.4 (19.97-1.35)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 1.35Å)	Xtriage
Refinement program	REFMAC 5.6.0093	Depositor
R, R_{free}	0.164 , 0.194 0.163 , 0.194	Depositor DCC
R_{free} test set	6456 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	10.5	Xtriage
Anisotropy	0.248	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 47.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5560	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.18% 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: I5I, GOL

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	1.84	59/4742 (1.2%)	1.51	31/6450 (0.5%)

All (59) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	255	MET	CG-SD	-9.91	1.55	1.80
1	A	170	ILE	CA-CB	-9.10	1.44	1.54
1	A	191	GLY	CA-C	8.96	1.58	1.52
1	A	328	ASP	N-CA	7.59	1.54	1.46
1	A	253	TYR	N-CA	-7.54	1.37	1.46
1	A	302	LYS	N-CA	7.46	1.55	1.46
1	A	97	THR	CA-C	7.18	1.61	1.52
1	A	142	LEU	N-CA	-6.99	1.37	1.46
1	A	491	SER	N-CA	-6.67	1.38	1.45
1	A	525	GLY	N-CA	-6.63	1.38	1.45
1	A	95	SER	CA-C	6.56	1.61	1.52
1	A	250	LYS	CA-C	6.50	1.61	1.52
1	A	368	TYR	N-CA	6.15	1.54	1.46
1	A	460	MET	N-CA	6.03	1.52	1.46
1	A	41	ILE	CA-CB	5.98	1.63	1.54
1	A	352	ALA	N-CA	5.89	1.53	1.46
1	A	24	THR	CA-C	5.89	1.60	1.52
1	A	361	THR	N-CA	5.75	1.53	1.45
1	A	573	ASP	CA-C	-5.74	1.44	1.52
1	A	175	TYR	CA-C	-5.69	1.46	1.52
1	A	461	TYR	C-O	5.63	1.30	1.23
1	A	157	PRO	CA-CB	5.63	1.62	1.53
1	A	167	GLU	N-CA	5.61	1.53	1.46
1	A	400	VAL	N-CA	-5.61	1.42	1.46
1	A	547	ASP	C-O	5.61	1.30	1.24
1	A	348	ASP	C-O	5.58	1.30	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	537	ASN	N-CA	5.54	1.53	1.46
1	A	372	GLU	C-O	5.50	1.31	1.24
1	A	138	HIS	CA-C	-5.50	1.45	1.52
1	A	514	ALA	CA-CB	5.50	1.62	1.53
1	A	478	ASP	N-CA	5.48	1.52	1.46
1	A	159[A]	ILE	C-O	5.42	1.30	1.24
1	A	159[B]	ILE	C-O	5.42	1.30	1.24
1	A	354	THR	CA-CB	5.41	1.61	1.54
1	A	542	GLU	CA-CB	-5.39	1.45	1.53
1	A	498	GLU	C-O	5.38	1.30	1.23
1	A	254	ALA	C-O	5.26	1.30	1.24
1	A	455	VAL	CA-CB	-5.26	1.47	1.54
1	A	539	SER	N-CA	5.25	1.52	1.46
1	A	560	LYS	CE-NZ	5.25	1.65	1.49
1	A	517	THR	N-CA	-5.23	1.39	1.46
1	A	234	TYR	C-O	5.23	1.29	1.23
1	A	39	SER	C-O	5.21	1.30	1.23
1	A	475	ASP	CA-CB	5.20	1.59	1.53
1	A	443	LEU	C-O	5.20	1.30	1.24
1	A	535	VAL	CA-C	-5.19	1.46	1.52
1	A	142	LEU	CA-CB	5.18	1.62	1.53
1	A	490	GLY	N-CA	5.18	1.51	1.45
1	A	235	GLN	N-CA	5.13	1.52	1.46
1	A	200	GLU	C-O	5.12	1.30	1.24
1	A	176	THR	CA-CB	5.11	1.61	1.53
1	A	303	THR	C-N	5.11	1.40	1.33
1	A	414	ARG	N-CA	5.10	1.53	1.45
1	A	338	ALA	CA-C	-5.09	1.46	1.52
1	A	183	ASP	CG-OD1	-5.07	1.15	1.25
1	A	541	GLY	N-CA	5.04	1.52	1.45
1	A	45	MET	C-O	5.04	1.30	1.24
1	A	411	SER	N-CA	5.03	1.52	1.46
1	A	303	THR	CA-CB	5.00	1.61	1.53

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	49[A]	SER	CA-C-N	9.73	129.34	120.10
1	A	49[A]	SER	C-N-CA	9.73	129.34	120.10
1	A	49[B]	SER	CA-C-N	9.73	129.34	120.10
1	A	49[B]	SER	C-N-CA	9.73	129.34	120.10
1	A	238	PHE	O-C-N	7.83	130.05	121.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	114	GLU	N-CA-CB	-7.09	99.51	110.85
1	A	347	LEU	O-C-N	6.71	130.53	122.27
1	A	214	GLN	O-C-N	6.64	129.16	122.12
1	A	250	LYS	CA-C-O	-6.61	111.05	120.51
1	A	45	MET	N-CA-C	6.14	122.22	114.31
1	A	570	TRP	CA-C-N	6.00	132.51	121.70
1	A	570	TRP	C-N-CA	6.00	132.51	121.70
1	A	115	ALA	O-C-N	5.99	131.01	123.23
1	A	254	ALA	CA-C-O	-5.94	114.12	120.42
1	A	331	ASN	N-CA-C	5.91	120.76	113.50
1	A	191	GLY	CA-C-O	-5.88	115.79	120.91
1	A	464	GLY	O-C-N	5.82	130.30	123.66
1	A	344	ARG	O-C-N	5.66	128.12	122.12
1	A	508	LYS	CG-CD-CE	-5.59	98.44	111.30
1	A	490	GLY	O-C-N	-5.57	117.16	123.61
1	A	292	THR	CA-CB-OG1	-5.49	101.36	109.60
1	A	593	THR	CA-C-O	-5.45	115.26	119.46
1	A	290	ARG	NE-CZ-NH2	-5.44	114.31	119.20
1	A	250	LYS	O-C-N	5.29	129.63	122.59
1	A	234	TYR	CA-C-O	-5.26	112.74	118.52
1	A	300	PHE	O-C-N	5.24	127.67	122.12
1	A	263	ASN	O-C-N	-5.16	116.72	121.37
1	A	308	ASP	N-CA-C	5.10	119.03	112.41
1	A	475	ASP	CA-C-N	5.04	124.76	119.82
1	A	475	ASP	C-N-CA	5.04	124.76	119.82
1	A	428	GLU	CA-C-O	-5.03	115.97	120.19

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	4551	0	4362	44	0
2	A	34	0	18	7	0
3	A	24	0	32	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	951	0	0	19	2
All	All	5560	0	4412	51	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 6.

All (51) 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:607:I5I:H13	4:A:791:HOH:O	1.67	0.93
1:A:557:ASN:HB3	4:A:1532:HOH:O	1.72	0.89
1:A:87[A]:VAL:HG21	1:A:103:TYR:CE1	2.12	0.83
2:A:606:I5I:S17	2:A:607:I5I:H13A	2.24	0.77
1:A:87[A]:VAL:CG2	1:A:103:TYR:CE1	2.68	0.76
2:A:607:I5I:H2	4:A:765:HOH:O	1.86	0.74
2:A:607:I5I:S17	2:A:607:I5I:H6	2.29	0.73
1:A:159[B]:ILE:HD13	1:A:592[B]:VAL:HG23	1.77	0.66
1:A:159[B]:ILE:CD1	1:A:592[B]:VAL:HG23	2.26	0.66
1:A:154:ASN:ND2	4:A:1446:HOH:O	2.29	0.65
2:A:606:I5I:S17	2:A:607:I5I:C13	2.87	0.61
1:A:79:ASN:OD1	1:A:89[B]:THR:HG22	2.01	0.61
1:A:87[A]:VAL:CG2	1:A:103:TYR:HE1	2.14	0.58
1:A:288:LYS:HD2	4:A:943:HOH:O	2.03	0.58
1:A:299:LYS:HE2	4:A:969:HOH:O	2.06	0.56
1:A:592[B]:VAL:HG12	1:A:593:THR:N	2.21	0.56
1:A:188:ILE:HD12	1:A:266:LEU:HD21	1.88	0.55
1:A:414:ARG:HD2	4:A:1367:HOH:O	2.07	0.54
1:A:47:THR:CG2	4:A:646:HOH:O	2.55	0.54
1:A:70:TRP:CH2	1:A:96:GLN:NE2	2.76	0.54
1:A:425:GLU:HB2	4:A:715:HOH:O	2.09	0.53
1:A:586:GLY:HA3	4:A:1204:HOH:O	2.09	0.52
1:A:47:THR:HG23	4:A:646:HOH:O	2.08	0.51
1:A:414:ARG:NH1	4:A:1367:HOH:O	2.16	0.51
1:A:89[B]:THR:HG21	4:A:803:HOH:O	2.09	0.51
1:A:194:PRO:HB3	1:A:277:LEU:HD22	1.93	0.51
1:A:78:TYR:O	1:A:89[A]:THR:HG23	2.12	0.50
1:A:89[B]:THR:HG21	4:A:821:HOH:O	2.12	0.49
1:A:35:ASN:OD1	1:A:37:GLN:HG2	2.13	0.48
1:A:55:VAL:O	4:A:1149:HOH:O	2.20	0.48
1:A:66:LYS:HE2	4:A:705:HOH:O	2.13	0.48
1:A:194:PRO:HG3	1:A:228[B]:HIS:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:596:ASN:H	1:A:596:ASN:HD22	1.63	0.46
1:A:35:ASN:O	1:A:37:GLN:OE1	2.35	0.45
1:A:70:TRP:HH2	1:A:96:GLN:NE2	2.15	0.45
1:A:194:PRO:HD2	1:A:226:VAL:O	2.17	0.44
2:A:606:I5I:H5	4:A:1084:HOH:O	2.17	0.43
1:A:34:ASN:O	1:A:34:ASN:ND2	2.52	0.42
1:A:156:ASP:CG	1:A:591:VAL:HG12	2.44	0.42
1:A:285:VAL:HG13	4:A:1509:HOH:O	2.18	0.42
1:A:363:ALA:HA	1:A:387:PHE:O	2.20	0.41
1:A:341[A]:ARG:HD3	1:A:382:TYR:OH	2.21	0.41
1:A:543:LEU:C	1:A:544:ILE:HD12	2.45	0.41
1:A:89[A]:THR:HG22	1:A:90:GLY:N	2.35	0.41
2:A:607:I5I:S17	2:A:607:I5I:C6	3.01	0.41
1:A:592[B]:VAL:CG1	1:A:593:THR:N	2.82	0.40
1:A:89[A]:THR:HG22	4:A:904:HOH:O	2.21	0.40
1:A:272:ILE:HD12	1:A:310:VAL:CG2	2.52	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)
4:A:649:HOH:O	4:A:1492:HOH:O[2_555]	2.18	0.02
4:A:713:HOH:O	4:A:1329:HOH:O[2_545]	2.18	0.02

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	601/584 (103%)	582 (97%)	19 (3%)	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	488/469 (104%)	477 (98%)	11 (2%)	44 13

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

Mol	Chain	Res	Type
1	A	26	PRO
1	A	35	ASN
1	A	64	LYS
1	A	141	PRO
1	A	189[A]	LEU
1	A	189[B]	LEU
1	A	305	LYS
1	A	557	ASN
1	A	587[A]	MET
1	A	587[B]	MET
1	A	596	ASN

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

Mol	Chain	Res	Type
1	A	37	GLN
1	A	52	ASN
1	A	199	ASN
1	A	596	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	GOL	A	3500[B]	-	5,5,5	0.54	0	5,5,5	0.41	0
2	I5I	A	606	-	18,19,19	2.82	6 (33%)	22,26,26	7.02	10 (45%)
3	GOL	A	3501	-	5,5,5	0.48	0	5,5,5	0.61	0
3	GOL	A	3500[A]	-	5,5,5	0.45	0	5,5,5	0.86	0
3	GOL	A	3502	-	5,5,5	0.94	0	5,5,5	1.44	1 (20%)
2	I5I	A	607	-	18,19,19	1.92	4 (22%)	22,26,26	6.41	6 (27%)

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	3500[B]	-	-	0/4/4/4	-
2	I5I	A	606	-	-	2/6/14/14	0/3/3/3
3	GOL	A	3501	-	-	0/4/4/4	-
3	GOL	A	3500[A]	-	-	0/4/4/4	-
3	GOL	A	3502	-	-	1/4/4/4	-
2	I5I	A	607	-	-	0/6/14/14	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	606	I5I	C10-N14	9.56	1.39	1.29
2	A	607	I5I	C10-N14	4.80	1.34	1.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	607	I5I	C16-N15	3.92	1.33	1.17
2	A	606	I5I	C3-C7	-3.76	1.32	1.44
2	A	606	I5I	C16-N15	3.47	1.32	1.17
2	A	607	I5I	C3-C7	-3.28	1.33	1.44
2	A	606	I5I	C12-N11	2.52	1.51	1.46
2	A	607	I5I	C16-S17	2.11	1.69	1.62
2	A	606	I5I	O9-C8	-2.04	1.34	1.38
2	A	606	I5I	C1-N15	-2.03	1.32	1.41

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	606	I5I	S17-C16-N15	-31.03	121.81	176.71
2	A	607	I5I	S17-C16-N15	-29.07	125.28	176.71
2	A	606	I5I	C13-C12-N11	5.56	107.16	101.60
2	A	606	I5I	C3-C2-C1	-4.27	117.41	121.43
2	A	606	I5I	C1-N15-C16	-4.18	118.03	161.05
2	A	606	I5I	C5-C4-C3	-3.39	118.73	123.81
2	A	607	I5I	O9-C4-C5	3.13	133.08	126.49
2	A	607	I5I	C1-N15-C16	-3.02	129.91	161.05
2	A	606	I5I	O9-C4-C5	3.02	132.84	126.49
2	A	607	I5I	C4-C3-C7	2.96	108.69	105.64
2	A	606	I5I	C2-C3-C4	2.71	122.78	118.55
2	A	607	I5I	O9-C4-C3	-2.58	106.89	110.12
2	A	607	I5I	C5-C4-C3	-2.52	120.03	123.81
2	A	606	I5I	O9-C8-C7	2.30	113.72	111.34
2	A	606	I5I	C12-C13-N14	-2.25	101.73	105.46
3	A	3502	GOL	O1-C1-C2	2.14	120.02	110.38
2	A	606	I5I	C4-O9-C8	-2.11	103.43	105.91

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	606	I5I	C2-C1-N15-C16
2	A	606	I5I	C6-C1-N15-C16
3	A	3502	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	606	I5I	3	0
2	A	607	I5I	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](#)

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	575/584 (98%)	-0.13	10 (1%) 69 79	3, 12, 26, 58	28 (4%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	596	ASN	3.5
1	A	205	VAL	2.7
1	A	70	TRP	2.7
1	A	35	ASN	2.7
1	A	71	SER	2.5
1	A	34	ASN	2.3
1	A	352	ALA	2.3
1	A	30	MET	2.1
1	A	46	GLU	2.0
1	A	245	TYR	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
2	I5I	A	607	17/17	0.72	0.17	51,62,86,93	0
2	I5I	A	606	17/17	0.83	0.14	22,33,68,68	0
3	GOL	A	3502	6/6	0.93	0.10	19,28,37,53	0
3	GOL	A	3501	6/6	0.95	0.06	18,21,22,22	0
3	GOL	A	3500[A]	6/6	0.98	0.03	4,4,5,5	6
3	GOL	A	3500[B]	6/6	0.98	0.03	8,10,12,19	6

6.5 Other polymers [i](#)

There are no such residues in this entry.