



Full wwPDB X-ray Structure Validation Report ⓘ

May 28, 2020 – 07:33 pm BST

PDB ID : 1GOH
Title : NOVEL THIOETHER BOND REVEALED BY A 1.7 ANGSTROMS CRYSTAL STRUCTURE OF GALACTOSE OXIDASE
Authors : Ito, N.; Phillips, S.E.V.; Knowles, P.F.
Deposited on : 1993-09-30
Resolution : 2.20 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

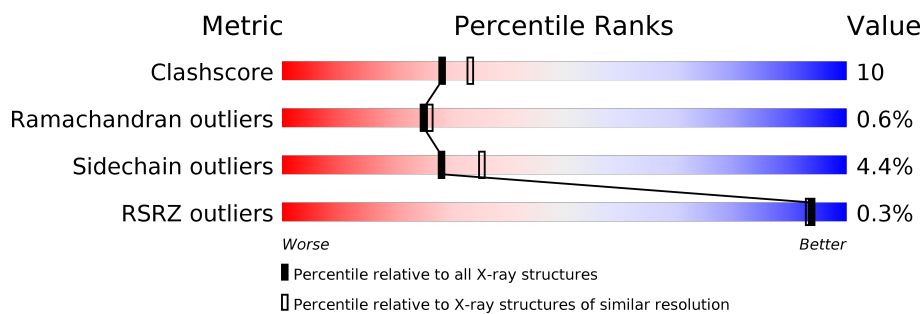
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 2.20 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))
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 on 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	639	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5141 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 GALACTOSE OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	639	Total	C	N	O	S	0	0	0
			4830	3017	840	954	19			

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		

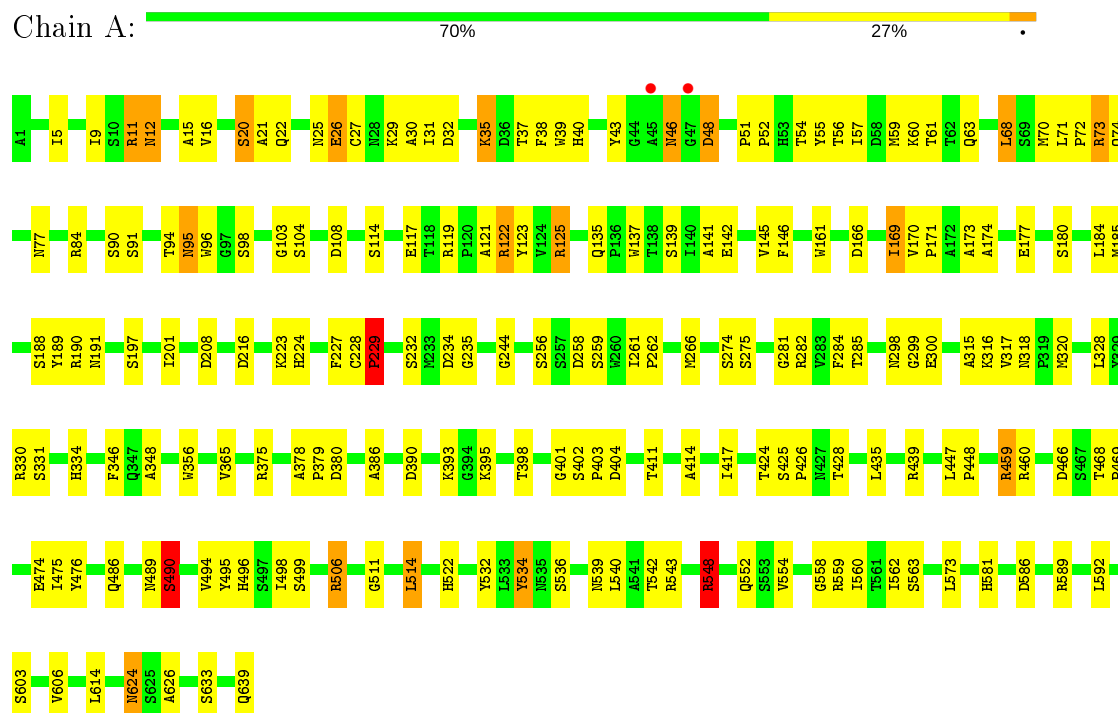
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	310	Total	O	0	0
			310	310		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: GALACTOSE OXIDASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	98.00Å 89.40Å 86.70Å 90.00° 117.80° 90.00°	Depositor
Resolution (Å)	10.00 – 2.20 9.99 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.20) 77.4 (9.99-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.26 (at 2.19Å)	Xtriage
Refinement program	PROLSQ	Depositor
R, R_{free}	0.156 , (Not available) 0.132 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	22.1	Xtriage
Anisotropy	0.175	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.19 , 43.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5141	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.96% 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: NA

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	0.99	1/4959 (0.0%)	1.73	66/6765 (1.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	490	SER	CA-CB	5.08	1.60	1.52

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	548	ARG	CD-NE-CZ	17.34	147.88	123.60
1	A	122	ARG	CD-NE-CZ	12.27	140.78	123.60
1	A	190	ARG	NE-CZ-NH1	11.00	125.80	120.30
1	A	122	ARG	NE-CZ-NH1	10.94	125.77	120.30
1	A	84	ARG	NE-CZ-NH2	-10.82	114.89	120.30
1	A	543	ARG	NE-CZ-NH1	9.85	125.22	120.30
1	A	460	ARG	NE-CZ-NH2	-9.25	115.67	120.30
1	A	476	TYR	CB-CG-CD2	-8.62	115.83	121.00
1	A	68	LEU	CA-CB-CG	8.51	134.88	115.30
1	A	548	ARG	NE-CZ-NH2	8.48	124.54	120.30
1	A	282	ARG	NE-CZ-NH1	8.32	124.46	120.30
1	A	543	ARG	NE-CZ-NH2	-8.00	116.30	120.30
1	A	119	ARG	NE-CZ-NH1	7.74	124.17	120.30
1	A	20	SER	N-CA-CB	7.69	122.03	110.50
1	A	119	ARG	NE-CZ-NH2	-7.64	116.48	120.30
1	A	499	SER	N-CA-CB	7.63	121.94	110.50
1	A	125	ARG	NE-CZ-NH2	-7.17	116.71	120.30
1	A	330	ARG	NE-CZ-NH2	-7.14	116.73	120.30
1	A	26	GLU	CA-CB-CG	7.13	129.09	113.40
1	A	459	ARG	NE-CZ-NH2	-7.09	116.75	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	234	ASP	CB-CG-OD1	6.81	124.43	118.30
1	A	459	ARG	NE-CZ-NH1	6.73	123.67	120.30
1	A	114	SER	N-CA-CB	6.51	120.27	110.50
1	A	48	ASP	CB-CG-OD2	-6.44	112.50	118.30
1	A	70	MET	CG-SD-CE	6.42	110.47	100.20
1	A	404	ASP	O-C-N	6.42	132.96	122.70
1	A	300	GLU	OE1-CD-OE2	6.40	130.97	123.30
1	A	534	TYR	CB-CG-CD1	-6.33	117.20	121.00
1	A	258	ASP	CB-CG-OD1	6.27	123.94	118.30
1	A	506	ARG	NE-CZ-NH1	6.12	123.36	120.30
1	A	559	ARG	NE-CZ-NH2	-6.03	117.28	120.30
1	A	190	ARG	NE-CZ-NH2	-6.02	117.29	120.30
1	A	633	SER	N-CA-CB	6.00	119.50	110.50
1	A	300	GLU	N-CA-CB	5.93	121.28	110.60
1	A	54	THR	N-CA-CB	-5.90	99.10	110.30
1	A	330	ARG	NE-CZ-NH1	5.89	123.25	120.30
1	A	586	ASP	CB-CG-OD1	-5.82	113.06	118.30
1	A	606	VAL	CA-CB-CG1	-5.82	102.17	110.90
1	A	375	ARG	NE-CZ-NH2	-5.76	117.42	120.30
1	A	460	ARG	CD-NE-CZ	-5.76	115.54	123.60
1	A	35	LYS	CA-CB-CG	5.74	126.03	113.40
1	A	188	SER	O-C-N	5.63	131.71	122.70
1	A	386	ALA	CB-CA-C	5.56	118.44	110.10
1	A	119	ARG	CD-NE-CZ	5.55	131.37	123.60
1	A	108	ASP	CB-CA-C	5.54	121.48	110.40
1	A	224	HIS	N-CA-CB	-5.54	100.63	110.60
1	A	404	ASP	CB-CG-OD2	-5.51	113.34	118.30
1	A	315	ALA	O-C-N	5.50	131.50	122.70
1	A	390	ASP	CB-CG-OD2	-5.42	113.42	118.30
1	A	20	SER	O-C-N	5.38	131.30	122.70
1	A	104	SER	O-C-N	5.33	131.23	122.70
1	A	121	ALA	CB-CA-C	5.29	118.03	110.10
1	A	624	ASN	CB-CA-C	5.29	120.97	110.40
1	A	266	MET	CA-CB-CG	-5.27	104.33	113.30
1	A	73	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	A	169	ILE	O-C-N	5.23	131.07	122.70
1	A	380	ASP	CB-CG-OD1	5.20	122.98	118.30
1	A	318	ASN	CB-CA-C	5.18	120.75	110.40
1	A	614	LEU	CA-CB-CG	5.17	127.20	115.30
1	A	586	ASP	CB-CG-OD2	5.15	122.94	118.30
1	A	589	ARG	NE-CZ-NH2	-5.14	117.73	120.30
1	A	208	ASP	CB-CG-OD2	5.13	122.92	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	435	LEU	O-C-N	5.07	130.81	122.70
1	A	11	ARG	CG-CD-NE	-5.03	101.24	111.80
1	A	84	ARG	NE-CZ-NH1	5.02	122.81	120.30
1	A	232	SER	N-CA-CB	5.01	118.01	110.50

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	4830	0	4604	92	1
2	A	1	0	0	0	0
3	A	310	0	0	6	1
All	All	5141	0	4604	92	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 10.

All (92) 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:22:GLN:HE21	1:A:46:ASN:HB2	1.32	0.94
1:A:35:LYS:HE3	1:A:71:LEU:HD21	1.58	0.86
1:A:22:GLN:NE2	1:A:46:ASN:HB2	1.92	0.84
1:A:122:ARG:HD3	1:A:123:TYR:CE1	2.22	0.75
1:A:417:ILE:HG12	1:A:428:THR:HG22	1.69	0.73
1:A:298:ASN:HD22	1:A:317:VAL:H	1.36	0.72
1:A:379:PRO:HD2	3:A:769:HOH:O	1.93	0.67
1:A:32:ASP:OD2	1:A:37:THR:OG1	2.11	0.65
1:A:60:LYS:HA	1:A:122:ARG:HH21	1.61	0.65
1:A:174:ALA:HB3	1:A:498:ILE:HD13	1.78	0.64
1:A:73:ARG:HD3	1:A:77:ASN:O	1.97	0.64
1:A:298:ASN:ND2	1:A:317:VAL:H	1.96	0.62
1:A:35:LYS:HE2	1:A:74:GLN:HG3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:ILE:HD12	1:A:145:VAL:HG12	1.85	0.58
1:A:189:TYR:CD1	1:A:197:SER:HB2	2.38	0.58
1:A:552:GLN:HG3	3:A:961:HOH:O	2.04	0.57
1:A:475:ILE:HD12	1:A:486:GLN:CD	2.24	0.57
1:A:495:TYR:O	1:A:496:HIS:HB2	2.04	0.57
1:A:30:ALA:HB3	1:A:55:TYR:OH	2.05	0.56
1:A:146:PHE:CD1	1:A:146:PHE:N	2.73	0.56
1:A:71:LEU:HD12	1:A:72:PRO:HD2	1.87	0.56
1:A:30:ALA:O	1:A:142:GLU:HA	2.05	0.56
1:A:171:PRO:HD2	1:A:511:GLY:HA2	1.89	0.55
1:A:174:ALA:HA	1:A:184:LEU:O	2.06	0.55
1:A:21:ALA:HA	1:A:40:HIS:O	2.06	0.55
1:A:56:THR:HG21	1:A:125:ARG:HH21	1.70	0.55
1:A:103:GLY:HA3	1:A:166:ASP:O	2.08	0.53
1:A:11:ARG:NH1	1:A:31:ILE:HA	2.23	0.53
1:A:22:GLN:NE2	1:A:43:TYR:O	2.42	0.52
1:A:573:LEU:HG	1:A:592:LEU:HD11	1.92	0.52
1:A:170:VAL:HB	1:A:514:LEU:HD13	1.91	0.52
1:A:95:ASN:N	1:A:95:ASN:HD22	2.08	0.52
1:A:38:PHE:HB3	1:A:141:ALA:HA	1.92	0.52
1:A:439:ARG:NH2	1:A:474:GLU:HG3	2.25	0.51
1:A:560:ILE:O	1:A:603:SER:HA	2.10	0.51
1:A:328:LEU:HA	1:A:331:SER:OG	2.11	0.51
1:A:522:HIS:HD2	3:A:820:HOH:O	1.94	0.50
1:A:5:ILE:HD12	1:A:490:SER:HB3	1.93	0.50
1:A:227:PHE:O	1:A:244:GLY:HA3	2.11	0.50
1:A:548:ARG:O	1:A:562:ILE:HA	2.11	0.50
1:A:16:VAL:HG12	1:A:57:ILE:HG12	1.92	0.50
1:A:90:SER:HB2	1:A:96:TRP:CE3	2.47	0.49
1:A:228:CYS:N	1:A:229:PRO:HD3	2.26	0.48
1:A:317:VAL:O	1:A:320:MET:HG2	2.13	0.48
1:A:522:HIS:CD2	3:A:820:HOH:O	2.66	0.48
1:A:94:THR:HB	1:A:95:ASN:HD22	1.77	0.48
1:A:285:THR:O	1:A:299:GLY:HA2	2.14	0.47
1:A:39:TRP:O	1:A:139:SER:HA	2.14	0.47
1:A:539:ASN:HB3	3:A:893:HOH:O	2.13	0.47
1:A:174:ALA:CB	1:A:498:ILE:HD13	2.42	0.47
1:A:177:GLU:HB2	1:A:180:SER:OG	2.15	0.47
1:A:235:GLY:HA3	1:A:281:GLY:HA3	1.97	0.47
1:A:401:GLY:O	1:A:411:THR:HG22	2.15	0.47
1:A:189:TYR:HB2	1:A:201:ILE:HG23	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:624:ASN:OD1	1:A:626:ALA:HB3	2.15	0.47
1:A:425:SER:HA	1:A:426:PRO:HD2	1.81	0.46
1:A:61:THR:O	1:A:122:ARG:HA	2.15	0.46
1:A:275:SER:HA	1:A:284:PHE:O	2.16	0.46
1:A:554:VAL:HG22	1:A:558:GLY:HA3	1.98	0.46
1:A:59:MET:C	1:A:61:THR:H	2.18	0.46
1:A:169:ILE:HG22	1:A:191:ASN:HB2	1.96	0.45
1:A:123:TYR:CD1	1:A:123:TYR:N	2.84	0.45
1:A:15:ALA:HB2	1:A:60:LYS:HZ2	1.82	0.45
1:A:298:ASN:ND2	1:A:316:LYS:HA	2.32	0.44
1:A:378:ALA:HA	1:A:379:PRO:HD3	1.85	0.44
1:A:46:ASN:HD22	1:A:46:ASN:HA	1.50	0.44
1:A:356:TRP:O	1:A:365:VAL:HA	2.17	0.43
1:A:402:SER:HB2	1:A:403:PRO:HD2	2.00	0.43
1:A:173:ALA:O	1:A:185:MET:HA	2.18	0.43
1:A:506:ARG:HD3	3:A:819:HOH:O	2.18	0.43
1:A:334:HIS:CE1	1:A:581:HIS:HB3	2.53	0.43
1:A:94:THR:HB	1:A:95:ASN:ND2	2.34	0.43
1:A:59:MET:O	1:A:61:THR:HG22	2.19	0.43
1:A:393:LYS:HB2	1:A:395:LYS:HG2	2.00	0.42
1:A:554:VAL:CG2	1:A:558:GLY:HA3	2.49	0.42
1:A:117:GLU:OE2	1:A:490:SER:HB2	2.19	0.42
1:A:534:TYR:CE1	1:A:540:LEU:HD23	2.55	0.42
1:A:26:GLU:OE1	1:A:29:LYS:HE2	2.20	0.41
1:A:468:THR:N	1:A:469:PRO:CD	2.83	0.41
1:A:261:ILE:HA	1:A:262:PRO:HD3	1.85	0.41
1:A:466:ASP:OD2	1:A:522:HIS:HE1	2.02	0.41
1:A:161:TRP:CE2	1:A:489:ASN:HB3	2.55	0.41
1:A:59:MET:O	1:A:60:LYS:HB2	2.21	0.41
1:A:43:TYR:HA	1:A:48:ASP:OD1	2.21	0.41
1:A:21:ALA:HB1	1:A:26:GLU:HA	2.03	0.41
1:A:398:THR:O	1:A:414:ALA:HA	2.21	0.41
1:A:51:PRO:HA	1:A:52:PRO:C	2.41	0.40
1:A:346:PHE:CE2	1:A:348:ALA:HB2	2.56	0.40
1:A:447:LEU:HA	1:A:448:PRO:HD3	1.89	0.40
1:A:61:THR:HG23	1:A:63:GLN:NE2	2.36	0.40
1:A:22:GLN:O	1:A:25:ASN:N	2.49	0.40
1:A:424:THR:O	1:A:426:PRO:HD3	2.22	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)
1:A:12:ASN:ND2	1:A:548:ARG:NH1[3_545]	2.14	0.06
3:A:836:HOH:O	3:A:958:HOH:O[2_656]	2.17	0.03

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	637/639 (100%)	604 (95%)	29 (5%)	4 (1%)	25 26

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	12	ASN
1	A	514	LEU
1	A	494	VAL
1	A	229	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	526/526 (100%)	503 (96%)	23 (4%)	28 35

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	SER

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Mol	Chain	Res	Type
1	A	27	CYS
1	A	46	ASN
1	A	68	LEU
1	A	91	SER
1	A	95	ASN
1	A	98	SER
1	A	135	GLN
1	A	137	TRP
1	A	216	ASP
1	A	223	LYS
1	A	229	PRO
1	A	256	SER
1	A	259	SER
1	A	274	SER
1	A	459	ARG
1	A	490	SER
1	A	532	TYR
1	A	536	SER
1	A	542	THR
1	A	548	ARG
1	A	563	SER
1	A	639	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	46	ASN
1	A	78	GLN
1	A	95	ASN
1	A	298	ASN
1	A	427	ASN
1	A	522	HIS
1	A	552	GLN
1	A	597	ASN
1	A	600	ASN

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

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 [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	636/639 (99%)	-1.25	2 (0%) 94 93	6, 21, 65, 100	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	47	GLY	3.1
1	A	45	ALA	2.8

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 carbohydrates 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(Å ²)	Q<0.9
2	NA	A	702	1/1	0.99	0.04	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.