



Full wwPDB X-ray Structure Validation Report ⓘ

May 12, 2020 – 11:12 pm BST

PDB ID : 5UN9
Title : The crystal structure of human O-GlcNAcase in complex with Thiamet-G
Authors : Li, B.; Jiang, J.
Deposited on : 2017-01-30
Resolution : 2.50 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
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

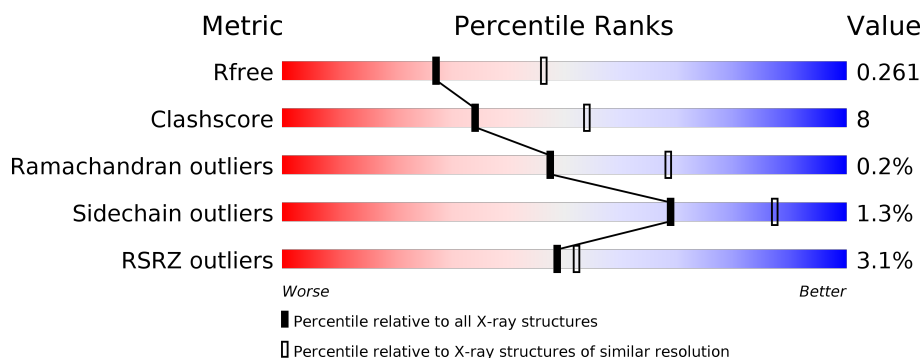
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	504	<div> <div>4%</div> <div> <div></div> <div>67%</div> <div>15%</div> <div>•</div> <div>17%</div> </div> </div>
1	B	504	<div> <div>%</div> <div> <div></div> <div>70%</div> <div>14%</div> <div>•</div> <div>15%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7189 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 Protein O-GlcNAcase.

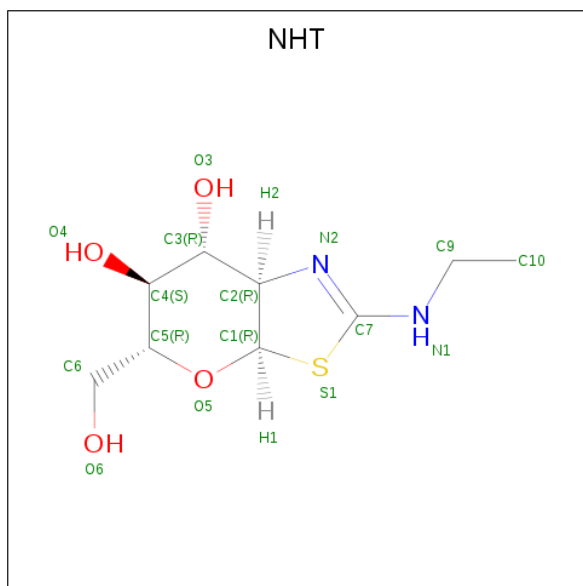
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	418	Total	C	N	O	S	0	0	0
			3426	2217	564	623	22			
1	B	426	Total	C	N	O	S	0	0	0
			3496	2264	577	632	23			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	59	HIS	-	expression tag	UNP O60502
A	543	GLY	-	linker	UNP O60502
A	544	GLY	-	linker	UNP O60502
A	545	GLY	-	linker	UNP O60502
A	546	GLY	-	linker	UNP O60502
A	547	SER	-	linker	UNP O60502
A	548	GLY	-	linker	UNP O60502
A	549	GLY	-	linker	UNP O60502
A	550	GLY	-	linker	UNP O60502
A	551	GLY	-	linker	UNP O60502
A	552	SER	-	linker	UNP O60502
B	59	HIS	-	expression tag	UNP O60502
B	543	GLY	-	linker	UNP O60502
B	544	GLY	-	linker	UNP O60502
B	545	GLY	-	linker	UNP O60502
B	546	GLY	-	linker	UNP O60502
B	547	SER	-	linker	UNP O60502
B	548	GLY	-	linker	UNP O60502
B	549	GLY	-	linker	UNP O60502
B	550	GLY	-	linker	UNP O60502
B	551	GLY	-	linker	UNP O60502
B	552	SER	-	linker	UNP O60502

- Molecule 2 is (3AR,5R,6S,7R,7AR)-2-(ETHYLAMINO)-5-(HYDROXYMETHYL)-5,6,7,7A-TETRAHYDRO-3AH-PYRANO[3,2-D][1,3]THIAZOLE-6,7-DIOL (three-letter code:

NHT) (formula: C₉H₁₆N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			16	9	2	4	1		
2	B	1	Total	C	N	O	S	0	0
			16	9	2	4	1		

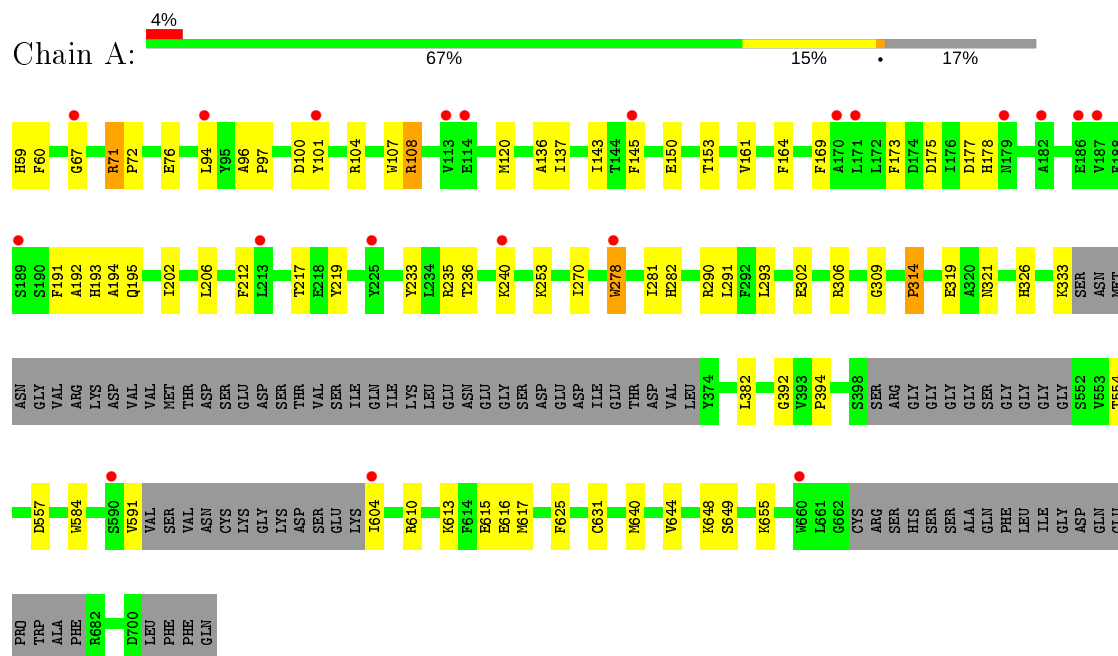
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	108	Total	O	0	0
			108	108		
3	B	127	Total	O	0	0
			127	127		

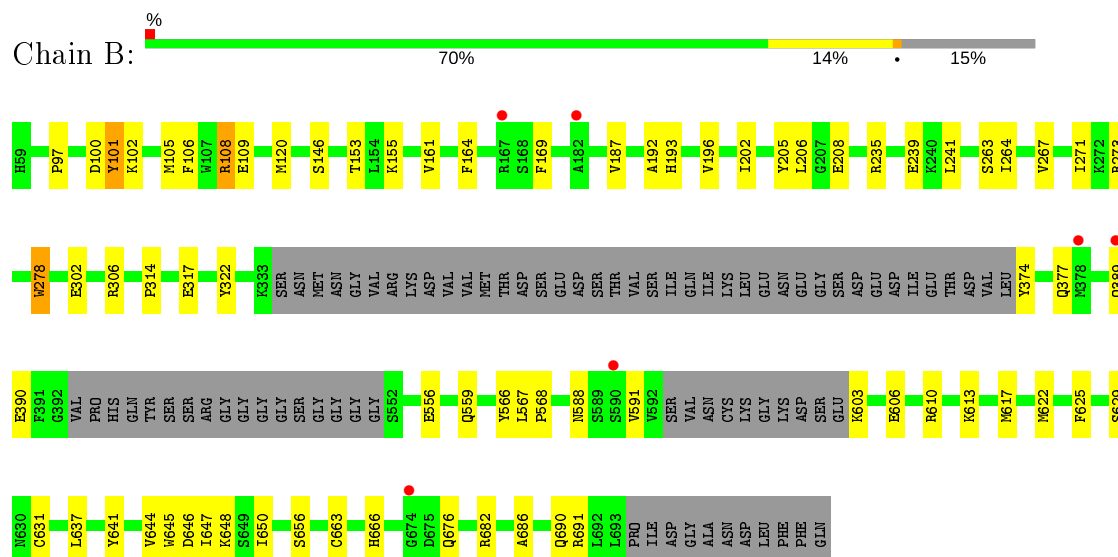
3 Residue-property plots

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: Protein O-GlcNAcase



• Molecule 1: Protein O-GlcNAcase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.57Å 96.11Å 89.53Å 90.00° 115.01° 90.00°	Depositor
Resolution (Å)	29.80 – 2.50 29.80 – 2.49	Depositor EDS
% Data completeness (in resolution range)	94.7 (29.80-2.50) 94.6 (29.80-2.49)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 2.48Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.210 , 0.260 0.210 , 0.261	Depositor DCC
R_{free} test set	2107 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	34.3	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.029 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7189	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% 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: NHT

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	0.61	0/3517	0.73	4/4760 (0.1%)
1	B	0.54	0/3590	0.70	4/4857 (0.1%)
All	All	0.57	0/7107	0.71	8/9617 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	108	ARG	NE-CZ-NH1	-11.39	114.61	120.30
1	B	108	ARG	NE-CZ-NH2	-7.90	116.35	120.30
1	B	108	ARG	NE-CZ-NH1	7.79	124.20	120.30
1	A	108	ARG	NE-CZ-NH2	7.65	124.13	120.30
1	A	71	ARG	NE-CZ-NH2	6.35	123.47	120.30
1	B	208	GLU	CA-CB-CG	6.24	127.13	113.40
1	B	389	GLN	CA-CB-CG	-6.02	100.15	113.40
1	A	71	ARG	NE-CZ-NH1	-5.21	117.69	120.30

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	3426	0	3348	66	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3496	0	3422	52	0
2	A	16	0	16	2	0
2	B	16	0	16	1	0
3	A	108	0	0	15	0
3	B	127	0	0	18	0
All	All	7189	0	6802	116	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 8.

All (116) 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:556:GLU:OE2	3:B:901:HOH:O	1.82	0.96
1:A:145:PHE:HD1	1:A:194:ALA:HB1	1.38	0.88
1:B:192:ALA:O	3:B:903:HOH:O	1.91	0.88
1:A:557:ASP:OD1	3:A:901:HOH:O	1.92	0.88
1:A:616:GLU:OE1	3:A:902:HOH:O	1.92	0.87
1:B:317:GLU:OE2	3:B:902:HOH:O	1.91	0.87
1:A:326:HIS:NE2	1:A:382:LEU:HD23	1.89	0.87
1:B:196:VAL:HB	3:B:903:HOH:O	1.77	0.84
1:A:193:HIS:NE2	1:A:240:LYS:HE2	1.91	0.84
1:B:374:TYR:N	3:B:908:HOH:O	2.10	0.83
1:A:290:ARG:NH1	3:A:905:HOH:O	2.11	0.82
1:A:59:HIS:N	3:A:906:HOH:O	2.12	0.81
1:B:390:GLU:OE1	3:B:904:HOH:O	1.97	0.80
1:A:143:ILE:HG23	1:A:145:PHE:HE2	1.46	0.79
1:B:105:MET:O	3:B:905:HOH:O	2.03	0.77
1:A:236:THR:HG23	1:A:240:LYS:HD2	1.66	0.75
1:A:319:GLU:OE2	3:A:903:HOH:O	2.03	0.75
1:A:108:ARG:NH1	1:A:150:GLU:OE1	2.20	0.74
1:B:559:GLN:OE1	3:B:906:HOH:O	2.05	0.74
1:A:178:HIS:HD2	1:A:191:PHE:HD2	1.37	0.73
1:A:591:VAL:HG12	1:A:604:ILE:HD13	1.72	0.71
1:B:101:TYR:CZ	1:B:102:LYS:HE3	2.26	0.70
1:B:278:TRP:CD2	2:B:801:NHT:H11	2.29	0.68
1:A:178:HIS:HD2	1:A:191:PHE:CD2	2.12	0.67
1:A:145:PHE:CD1	1:A:194:ALA:HB1	2.28	0.66
1:B:646:ASP:OD1	3:B:909:HOH:O	2.13	0.66
1:B:686:ALA:O	1:B:690:GLN:HG3	1.95	0.66
1:A:173:PHE:O	1:A:217:THR:HG22	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:ARG:HH12	1:A:150:GLU:CD	1.99	0.65
1:A:178:HIS:CD2	1:A:191:PHE:HD2	2.15	0.65
1:A:177:ASP:OD1	1:A:178:HIS:N	2.29	0.65
1:A:71:ARG:O	3:A:907:HOH:O	2.13	0.64
1:B:322:TYR:HB3	3:B:915:HOH:O	1.97	0.64
1:A:143:ILE:CG2	1:A:145:PHE:HE2	2.10	0.64
1:A:175:ASP:N	3:A:909:HOH:O	2.19	0.62
1:B:202:ILE:O	1:B:206:LEU:HD12	1.99	0.62
1:A:615:GLU:HB2	1:A:655:LYS:HG3	1.82	0.60
1:A:333:LYS:O	3:A:908:HOH:O	2.17	0.60
1:B:146:SER:HB3	1:B:187:VAL:HG21	1.82	0.60
1:A:97:PRO:HG2	1:A:100:ASP:HB2	1.85	0.59
1:A:281:ILE:HG23	1:A:282:HIS:CD2	2.37	0.59
1:B:629:SER:HA	1:B:637:LEU:HD11	1.85	0.58
1:B:241:LEU:O	1:B:273:ARG:NH2	2.36	0.58
1:A:625:PHE:HD1	1:A:644:VAL:HG12	1.67	0.58
1:A:143:ILE:HG23	1:A:145:PHE:CE2	2.33	0.58
1:A:326:HIS:HD2	3:A:945:HOH:O	1.87	0.57
1:A:278:TRP:CD2	2:A:801:NHT:H11	2.39	0.57
1:A:591:VAL:CG1	1:A:604:ILE:HD13	2.34	0.57
1:A:202:ILE:O	1:A:206:LEU:HD12	2.06	0.56
1:B:271:ILE:O	1:B:273:ARG:NH1	2.38	0.56
1:A:193:HIS:CD2	1:A:240:LYS:HE2	2.41	0.56
1:A:291:LEU:HD21	1:A:640:MET:HE3	1.87	0.55
1:A:145:PHE:HD1	1:A:194:ALA:CB	2.16	0.55
1:B:682:ARG:O	1:B:691:ARG:NH2	2.39	0.55
1:B:377:GLN:H	1:B:377:GLN:CD	2.10	0.53
1:B:663:CYS:HA	1:B:666:HIS:CE1	2.44	0.52
1:B:603:LYS:NZ	3:B:907:HOH:O	2.08	0.52
1:A:392:GLY:N	3:A:914:HOH:O	2.40	0.52
1:A:326:HIS:NE2	1:A:382:LEU:CD2	2.67	0.52
1:B:97:PRO:HG2	1:B:100:ASP:HB2	1.92	0.51
1:A:584:TRP:CD2	1:A:610:ARG:NH2	2.79	0.51
1:A:76:GLU:N	1:A:76:GLU:OE1	2.34	0.51
1:A:253:LYS:NZ	1:B:676:GLN:O	2.36	0.51
1:A:314:PRO:HB2	1:A:321:ASN:OD1	2.11	0.50
1:B:109:GLU:OE1	3:B:910:HOH:O	2.20	0.50
1:A:104:ARG:NH1	3:A:916:HOH:O	2.39	0.50
1:A:235:ARG:HG3	1:A:270:ILE:HD11	1.93	0.50
1:A:67:GLY:HA2	1:A:96:ALA:O	2.12	0.50
1:B:263:SER:O	3:B:911:HOH:O	2.20	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:TRP:HB2	1:A:137:ILE:HD11	1.95	0.48
1:B:235:ARG:O	1:B:239:GLU:HG3	2.13	0.48
1:B:161:VAL:HG13	1:B:169:PHE:HE2	1.79	0.48
1:B:271:ILE:O	1:B:273:ARG:HG3	2.15	0.47
1:B:606:GLU:HB3	3:B:928:HOH:O	2.15	0.47
1:B:588:ASN:O	1:B:591:VAL:HG22	2.14	0.47
1:B:637:LEU:CD2	1:B:641:TYR:HD2	2.28	0.46
1:A:169:PHE:O	1:A:212:PHE:HA	2.15	0.46
1:A:161:VAL:HA	1:A:164:PHE:CD2	2.51	0.46
1:B:622:MET:SD	1:B:648:LYS:HE3	2.55	0.46
1:A:192:ALA:HB2	1:A:233:TYR:CD2	2.51	0.45
1:B:613:LYS:O	1:B:617:MET:HG2	2.16	0.45
1:A:625:PHE:CD1	1:A:644:VAL:HG12	2.48	0.45
1:A:94:LEU:HD11	1:A:136:ALA:HB2	1.99	0.45
1:A:195:GLN:HE22	1:A:217:THR:CG2	2.30	0.45
1:B:606:GLU:O	1:B:610:ARG:HG3	2.16	0.44
1:B:645:TRP:CZ3	1:B:648:LYS:HD2	2.51	0.44
1:B:105:MET:HG2	3:B:963:HOH:O	2.17	0.44
1:B:647:ILE:HA	1:B:650:ILE:HG22	1.98	0.44
1:A:302:GLU:O	1:A:306:ARG:HD2	2.18	0.44
1:A:394:PRO:HG2	1:B:106:PHE:CE2	2.52	0.44
1:B:567:LEU:HB3	1:B:568:PRO:HD2	2.00	0.44
1:A:108:ARG:HB3	1:A:153:THR:HG21	2.00	0.43
1:A:616:GLU:CG	3:A:902:HOH:O	2.66	0.43
1:A:613:LYS:O	1:A:617:MET:HG2	2.17	0.43
1:B:566:TYR:CD2	1:B:567:LEU:HD12	2.53	0.43
1:A:219:TYR:CG	2:A:801:NHT:H9C2	2.54	0.43
1:B:101:TYR:OH	1:B:102:LYS:HE3	2.19	0.43
1:A:107:TRP:CH2	1:A:143:ILE:HD13	2.54	0.42
1:A:293:LEU:HD21	1:A:640:MET:HE3	2.01	0.42
1:B:109:GLU:HB3	3:B:910:HOH:O	2.19	0.42
1:B:625:PHE:HD1	1:B:644:VAL:HG12	1.85	0.42
1:B:146:SER:CB	1:B:187:VAL:HG21	2.48	0.42
1:B:161:VAL:HA	1:B:164:PHE:CD2	2.54	0.42
1:A:281:ILE:HD13	1:A:281:ILE:HG21	1.78	0.42
1:B:264:ILE:HA	1:B:267:VAL:HG12	2.02	0.42
1:A:649:SER:HA	3:A:950:HOH:O	2.20	0.42
1:A:616:GLU:HG3	3:A:902:HOH:O	2.20	0.41
1:B:193:HIS:HA	3:B:903:HOH:O	2.19	0.41
1:A:143:ILE:CD1	1:A:150:GLU:HB3	2.50	0.41
1:A:554:THR:OG1	3:A:904:HOH:O	2.09	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:PRO:HB3	1:A:97:PRO:HB3	2.01	0.41
1:B:302:GLU:O	1:B:306:ARG:HD2	2.20	0.41
1:B:108:ARG:HB3	1:B:153:THR:HG21	2.03	0.41
1:B:193:HIS:C	3:B:903:HOH:O	2.59	0.41
1:B:155:LYS:HB3	1:B:205:TYR:CE2	2.56	0.41
1:A:60:PHE:CZ	1:A:309:GLY:HA2	2.56	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	A	408/504 (81%)	397 (97%)	10 (2%)	1 (0%)	47	68
1	B	418/504 (83%)	406 (97%)	11 (3%)	1 (0%)	47	68
All	All	826/1008 (82%)	803 (97%)	21 (2%)	2 (0%)	47	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	314	PRO
1	B	314	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	370/442 (84%)	365 (99%)	5 (1%)	67	86
1	B	377/442 (85%)	372 (99%)	5 (1%)	69	87
All	All	747/884 (84%)	737 (99%)	10 (1%)	69	87

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	101	TYR
1	A	120	MET
1	A	278	TRP
1	A	631	CYS
1	A	648	LYS
1	B	101	TYR
1	B	120	MET
1	B	278	TRP
1	B	631	CYS
1	B	656	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	178	HIS
1	A	282	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

2 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	NHT	A	801	-	15,17,17	5.73	7 (46%)	14,24,24	1.62	3 (21%)
2	NHT	B	801	-	15,17,17	5.43	6 (40%)	14,24,24	1.67	3 (21%)

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
2	NHT	A	801	-	-	1/5/33/33	0/2/2/2
2	NHT	B	801	-	-	1/5/33/33	0/2/2/2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	NHT	C2-N2	17.64	1.67	1.47
2	B	801	NHT	C2-N2	17.19	1.66	1.47
2	A	801	NHT	C7-N2	8.59	1.49	1.29
2	B	801	NHT	C7-N2	8.30	1.48	1.29
2	A	801	NHT	C7-S1	-5.82	1.60	1.76
2	A	801	NHT	O5-C1	5.58	1.51	1.42
2	B	801	NHT	C7-S1	-4.94	1.63	1.76
2	A	801	NHT	C7-N1	4.67	1.44	1.34
2	B	801	NHT	O5-C1	4.48	1.49	1.42
2	B	801	NHT	C7-N1	3.94	1.43	1.34
2	A	801	NHT	C4-C3	-3.56	1.43	1.52
2	B	801	NHT	C4-C5	2.99	1.59	1.53
2	A	801	NHT	C4-C5	2.18	1.57	1.53

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	NHT	O5-C1-C2	-3.85	106.80	115.27
2	B	801	NHT	C9-N1-C7	-2.79	116.39	123.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	NHT	C9-N1-C7	-2.71	116.59	123.50
2	A	801	NHT	O5-C1-C2	-2.67	109.41	115.27
2	A	801	NHT	O3-C3-C2	2.43	114.70	109.14
2	B	801	NHT	N1-C7-N2	-2.42	121.10	124.26

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	NHT	C10-C9-N1-C7
2	B	801	NHT	C10-C9-N1-C7

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	NHT	2	0
2	B	801	NHT	1	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	418/504 (82%)	0.02	20 (4%) 30 32	16, 38, 67, 89	0
1	B	426/504 (84%)	-0.10	6 (1%) 75 77	21, 46, 65, 85	0
All	All	844/1008 (83%)	-0.04	26 (3%) 49 52	16, 43, 66, 89	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	187	VAL	4.4
1	A	590	SER	4.0
1	B	389	GLN	3.6
1	A	660	TRP	3.5
1	A	604	ILE	3.4
1	A	225	TYR	3.3
1	A	182	ALA	3.3
1	A	186	GLU	2.9
1	A	94	LEU	2.7
1	A	170	ALA	2.6
1	B	167	ARG	2.6
1	B	182	ALA	2.6
1	A	113	VAL	2.5
1	A	213	LEU	2.5
1	A	278	TRP	2.4
1	B	674	GLY	2.3
1	A	101	TYR	2.3
1	A	114	GLU	2.3
1	A	189	SER	2.2
1	A	171	LEU	2.2
1	A	179	ASN	2.1
1	A	240	LYS	2.1
1	B	590	SER	2.1
1	A	145	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	67	GLY	2.1
1	B	378	MET	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 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(\AA^2)	Q<0.9
2	NHT	A	801	16/16	0.97	0.20	21,28,35,35	0
2	NHT	B	801	16/16	0.97	0.12	19,29,35,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.