



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:34 PM UTC

PDB ID : 1D0H / pdb_00001d0h
Title : THE HC FRAGMENT OF TETANUS TOXIN COMPLEXED WITH N-AC
ETYL-GALACTOSAMINE
Authors : Emsley, P.; Fotinou, C.; Black, I.; Fairweather, N.F.; Charles, I.G.; Watts, C.;
Hewitt, E.; Isaacs, N.W.
Deposited on : 1999-09-10
Resolution : 2.10 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

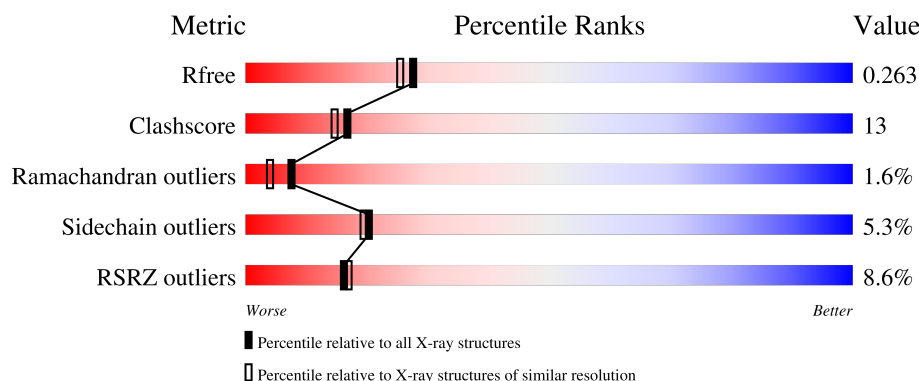
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	469	

2 Entry composition [i](#)

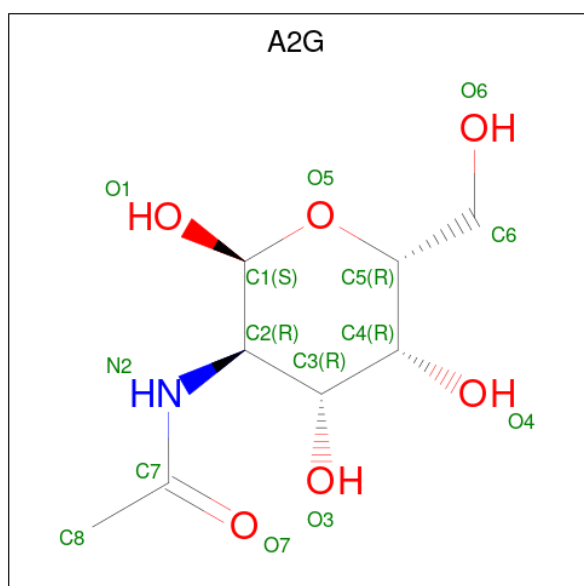
There are 4 unique types of molecules in this entry. The entry contains 4099 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 (TETANUS TOXIN HC).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	441	Total	C	N	O	S	0	0	0
			3563	2282	597	675	9			

- Molecule 2 is 2-acetamido-2-deoxy-alpha-D-galactopyranose (CCD ID: A2G) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			15	8	1	6		
2	A	1	Total	C	N	O	0	0
			15	8	1	6		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



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

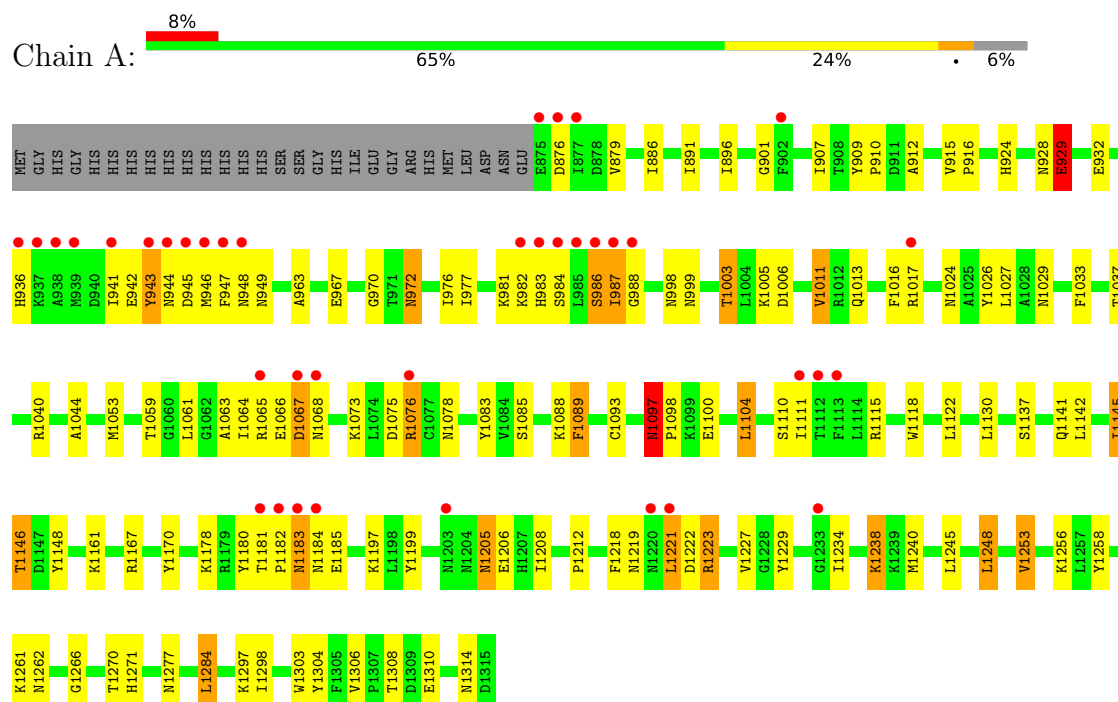
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	501	Total	O	0	0
			501	501		

3 Residue-property plots

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: PROTEIN (TETANUS TOXIN HC)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.08Å 70.88Å 122.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.10 40.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.00-2.10) 99.6 (40.00-2.10)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.68 (at 2.10Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.204 , 0.270 0.200 , 0.263	Depositor DCC
R_{free} test set	1764 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.639	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 64.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4099	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.42% 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: A2G, SO4

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	0.67	0/3644	1.59	33/4945 (0.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	1266	GLY	N-CA-C	8.75	120.97	112.04
1	A	1306	VAL	N-CA-C	8.58	116.78	107.60
1	A	1067	ASP	CA-CB-CG	8.31	120.91	112.60
1	A	1248	LEU	N-CA-C	8.12	120.13	111.28
1	A	928	ASN	CA-C-N	7.77	131.03	120.38
1	A	928	ASN	C-N-CA	7.77	131.03	120.38
1	A	1006	ASP	CA-CB-CG	7.40	120.00	112.60
1	A	901	GLY	N-CA-C	-7.16	105.19	114.85
1	A	1089	PHE	CA-CB-CG	-6.84	106.96	113.80
1	A	970	GLY	N-CA-C	6.48	121.73	113.24
1	A	1148	TYR	N-CA-C	6.41	119.88	110.59
1	A	1304	TYR	CA-C-O	-6.24	113.87	120.80
1	A	1184	ASN	N-CA-C	6.12	120.77	113.12
1	A	1024	ASN	CA-CB-CG	-6.03	106.57	112.60
1	A	1314	ASN	CA-CB-CG	5.87	118.47	112.60
1	A	1218	PHE	CA-CB-CG	-5.72	108.08	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1277	ASN	CA-CB-CG	-5.69	106.91	112.60
1	A	972	ASN	CA-CB-CG	5.57	118.17	112.60
1	A	929	GLU	CB-CG-CD	5.47	121.89	112.60
1	A	1253	VAL	N-CA-CB	-5.31	102.54	112.36
1	A	1097	ASN	N-CA-CB	5.26	115.43	109.49
1	A	1146	THR	N-CA-CB	5.22	119.31	110.49
1	A	1115	ARG	CD-NE-CZ	-5.20	117.13	124.40
1	A	1027	LEU	CA-C-N	-5.19	115.03	122.67
1	A	1027	LEU	C-N-CA	-5.19	115.03	122.67
1	A	1223	ARG	CD-NE-CZ	-5.17	117.16	124.40
1	A	1059	THR	N-CA-C	-5.15	103.24	110.50
1	A	1061	LEU	N-CA-C	-5.13	102.10	110.20
1	A	1059	THR	N-CA-CB	5.10	117.52	110.17
1	A	986	SER	CA-C-O	-5.08	113.24	120.51
1	A	1026	TYR	CA-CB-CG	-5.07	104.78	113.90
1	A	912	ALA	N-CA-C	-5.06	101.72	109.76
1	A	1266	GLY	CA-C-O	-5.02	118.84	122.45

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1065	ARG	Sidechain

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	3563	0	3516	90	0
2	A	30	0	23	2	0
3	A	5	0	0	0	0
4	A	501	0	0	13	0
All	All	4099	0	3539	91	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 13.

All (91) 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:1141:GLN:HE21	1:A:1142:LEU:H	1.05	0.98
1:A:929:GLU:H	1:A:929:GLU:CD	1.89	0.78
1:A:972:ASN:HD21	1:A:1078:ASN:H	1.31	0.78
1:A:1270:THR:HG22	1:A:1284:LEU:CD2	2.14	0.76
1:A:1093:CYS:HB3	4:A:207:HOH:O	1.89	0.71
1:A:1297:LYS:NZ	1:A:1298:ILE:HD11	2.04	0.71
1:A:1270:THR:HG22	1:A:1284:LEU:HD23	1.72	0.71
1:A:1141:GLN:NE2	1:A:1142:LEU:H	1.84	0.70
1:A:1223:ARG:HG2	1:A:1223:ARG:HH11	1.58	0.69
1:A:1182:PRO:O	1:A:1183:ASN:HB2	1.92	0.67
1:A:1053:MET:HE1	4:A:474:HOH:O	1.94	0.67
2:A:602:A2G:H8	4:A:252:HOH:O	1.94	0.67
1:A:1181:THR:HB	1:A:1182:PRO:HD2	1.76	0.67
1:A:1053:MET:SD	4:A:128:HOH:O	2.53	0.67
1:A:949:ASN:ND2	1:A:1040:ARG:H	1.98	0.62
1:A:1297:LYS:HZ3	1:A:1298:ILE:HD11	1.64	0.62
1:A:1017:ARG:HB2	4:A:301:HOH:O	2.00	0.61
1:A:944:ASN:O	1:A:948:ASN:HB3	2.02	0.60
1:A:945:ASP:HB3	4:A:77:HOH:O	2.02	0.60
1:A:1003:THR:HB	1:A:1013:GLN:HG2	1.84	0.60
1:A:1097:ASN:ND2	1:A:1100:GLU:H	2.01	0.59
1:A:987:ILE:HB	1:A:1005:LYS:HZ3	1.67	0.58
1:A:1083:TYR:OH	1:A:1310:GLU:HG3	2.05	0.56
1:A:1145:ILE:HD11	4:A:195:HOH:O	2.05	0.56
1:A:1180:TYR:HA	1:A:1197:LYS:NZ	2.20	0.56
1:A:907:ILE:HD12	1:A:936:HIS:CE1	2.41	0.56
1:A:1205:ASN:HB2	4:A:504:HOH:O	2.07	0.55
1:A:946:MET:HA	1:A:1064:ILE:O	2.07	0.55
1:A:876:ASP:HB2	1:A:879:VAL:HG23	1.88	0.54
1:A:1180:TYR:HA	1:A:1197:LYS:HZ2	1.72	0.54
1:A:891:ILE:HG12	1:A:896:ILE:HG12	1.89	0.54
1:A:1261:LYS:O	1:A:1262:ASN:HB2	2.08	0.53
1:A:924:HIS:HD2	1:A:1085:SER:OG	1.92	0.53
1:A:1270:THR:HG22	1:A:1284:LEU:HD22	1.90	0.52
1:A:1130:LEU:HD22	1:A:1303:TRP:HB3	1.91	0.52
1:A:987:ILE:HB	1:A:1005:LYS:NZ	2.25	0.51
1:A:1256:LYS:HE2	1:A:1258:TYR:CZ	2.46	0.51
1:A:948:ASN:ND2	1:A:949:ASN:H	2.07	0.51
1:A:932:GLU:OE1	1:A:1073:LYS:HD3	2.11	0.51
1:A:983:HIS:CD2	1:A:983:HIS:O	2.65	0.50
1:A:1067:ASP:OD1	1:A:1068:ASN:N	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1097:ASN:HB2	1:A:1098:PRO:HD2	1.94	0.49
1:A:1130:LEU:CD2	1:A:1303:TRP:HB3	2.42	0.49
1:A:1033:PHE:HZ	1:A:1104:LEU:HD13	1.76	0.49
1:A:1170:TYR:CE1	1:A:1308:THR:HA	2.48	0.49
1:A:1223:ARG:NH1	1:A:1223:ARG:CG	2.73	0.49
1:A:999:ASN:OD1	1:A:1017:ARG:HD2	2.13	0.47
1:A:1033:PHE:CZ	1:A:1104:LEU:HD13	2.48	0.47
1:A:1005:LYS:HG3	1:A:1011:VAL:HG23	1.97	0.47
1:A:1003:THR:HG23	4:A:2:HOH:O	2.15	0.47
1:A:1075:ASP:O	1:A:1076:ARG:HB2	2.14	0.46
1:A:1137:SER:HA	4:A:69:HOH:O	2.14	0.46
1:A:1097:ASN:HB2	1:A:1098:PRO:CD	2.46	0.46
1:A:1146:THR:N	1:A:1227:VAL:O	2.46	0.45
1:A:1212:PRO:HB3	1:A:1229:TYR:CE2	2.51	0.45
1:A:1208:ILE:HD13	1:A:1238:LYS:HG2	1.97	0.45
1:A:1221:LEU:O	1:A:1223:ARG:NH1	2.49	0.45
1:A:1145:ILE:HG12	1:A:1146:THR:N	2.32	0.45
1:A:1145:ILE:HD12	4:A:423:HOH:O	2.17	0.45
1:A:942:GLU:HG3	1:A:943:TYR:CD2	2.52	0.44
1:A:1017:ARG:HD2	1:A:1017:ARG:HH11	1.55	0.44
1:A:1029:ASN:HB3	1:A:1118:TRP:CZ3	2.53	0.44
1:A:1223:ARG:HG2	1:A:1223:ARG:NH1	2.28	0.44
1:A:947:PHE:CE1	1:A:1063:ALA:HB2	2.53	0.43
1:A:909:TYR:HB3	1:A:910:PRO:HD2	2.00	0.43
1:A:1181:THR:HB	1:A:1182:PRO:CD	2.46	0.43
1:A:1097:ASN:HD22	1:A:1097:ASN:C	2.26	0.43
1:A:963:ALA:O	1:A:967:GLU:HG3	2.19	0.43
1:A:981:LYS:HG3	1:A:1067:ASP:HB2	2.01	0.43
1:A:1199:TYR:HE2	2:A:602:A2G:C1	2.32	0.43
1:A:998:ASN:HD22	1:A:998:ASN:HA	1.66	0.42
1:A:976:ILE:HG13	1:A:977:ILE:HG13	2.01	0.42
1:A:1037:THR:O	1:A:1044:ALA:HA	2.20	0.42
1:A:987:ILE:HD12	1:A:1005:LYS:HE2	2.00	0.42
1:A:1016:PHE:HE1	1:A:1053:MET:HE2	1.84	0.42
1:A:1178:LYS:HD2	1:A:1199:TYR:CE1	2.54	0.42
1:A:1003:THR:HG21	4:A:303:HOH:O	2.20	0.42
1:A:915:VAL:HB	1:A:916:PRO:HD2	2.00	0.42
1:A:946:MET:HG3	1:A:1063:ALA:HB1	2.02	0.41
1:A:1110:SER:HB2	4:A:326:HOH:O	2.19	0.41
1:A:1088:LYS:O	1:A:1089:PHE:C	2.63	0.41
1:A:1205:ASN:HD22	1:A:1206:GLU:H	1.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1075:ASP:O	1:A:1076:ARG:CB	2.68	0.41
1:A:1222:ASP:OD2	1:A:1271:HIS:HD2	2.04	0.41
1:A:1182:PRO:O	1:A:1183:ASN:CB	2.63	0.41
1:A:1240:MET:HA	1:A:1256:LYS:O	2.21	0.41
1:A:876:ASP:HB2	1:A:879:VAL:CG2	2.50	0.41
1:A:972:ASN:ND2	1:A:1078:ASN:H	2.07	0.41
1:A:1066:GLU:H	1:A:1066:GLU:CD	2.29	0.41
1:A:1066:GLU:HG2	1:A:1067:ASP:O	2.21	0.41
1:A:886:ILE:HB	1:A:941:ILE:HG21	2.03	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	439/469 (94%)	411 (94%)	21 (5%)	7 (2%)	7 4

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	986	SER
1	A	988	GLY
1	A	1183	ASN
1	A	982	LYS
1	A	984	SER
1	A	1221	LEU
1	A	987	ILE

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	397/421 (94%)	376 (95%)	21 (5%)	20	19

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	929	GLU
1	A	943	TYR
1	A	1003	THR
1	A	1011	VAL
1	A	1076	ARG
1	A	1097	ASN
1	A	1104	LEU
1	A	1111	ILE
1	A	1122	LEU
1	A	1145	ILE
1	A	1161	LYS
1	A	1167	ARG
1	A	1185	GLU
1	A	1205	ASN
1	A	1219	ASN
1	A	1234	ILE
1	A	1238	LYS
1	A	1245	LEU
1	A	1248	LEU
1	A	1253	VAL
1	A	1284	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	892	ASN
1	A	919	ASN
1	A	924	HIS
1	A	936	HIS

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Mol	Chain	Res	Type
1	A	949	ASN
1	A	972	ASN
1	A	983	HIS
1	A	998	ASN
1	A	1013	GLN
1	A	1082	GLN
1	A	1097	ASN
1	A	1141	GLN
1	A	1171	ASN
1	A	1203	ASN
1	A	1205	ASN
1	A	1216	ASN
1	A	1271	HIS
1	A	1272	ASN
1	A	1280	ASN
1	A	1292	ASN

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](#)

3 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	A2G	A	601	-	15,15,15	0.35	0	21,21,21	0.99	2 (9%)
2	A2G	A	602	-	15,15,15	0.55	0	21,21,21	2.18	7 (33%)
3	SO4	A	701	-	4,4,4	0.75	0	6,6,6	0.19	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
2	A2G	A	601	-	-	1/6/26/26	0/1/1/1
2	A2G	A	602	-	-	2/6/26/26	0/1/1/1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	602	A2G	C2-N2-C7	4.88	134.54	123.11
2	A	602	A2G	C1-O5-C5	4.22	121.81	113.65
2	A	602	A2G	O7-C7-N2	3.40	128.00	121.98
2	A	602	A2G	C8-C7-N2	-3.05	111.05	116.12
2	A	601	A2G	O4-C4-C5	2.67	115.90	109.32
2	A	602	A2G	C1-C2-N2	2.54	113.67	110.73
2	A	602	A2G	O1-C1-C2	2.35	114.10	109.22
2	A	602	A2G	O5-C5-C4	2.20	113.66	109.70
2	A	601	A2G	O3-C3-C2	2.11	113.78	109.58

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	602	A2G	O5-C5-C6-O6
2	A	602	A2G	C4-C5-C6-O6
2	A	601	A2G	C3-C2-N2-C7

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	602	A2G	2	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	441/469 (94%)	0.39	38 (8%) 16 17	17, 30, 69, 100	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	985	LEU	8.8
1	A	987	ILE	6.8
1	A	1183	ASN	6.0
1	A	936	HIS	5.5
1	A	941	ILE	5.4
1	A	948	ASN	4.6
1	A	983	HIS	4.4
1	A	938	ALA	4.3
1	A	1182	PRO	4.1
1	A	944	ASN	4.0
1	A	877	ILE	3.9
1	A	937	LYS	3.7
1	A	943	TYR	3.6
1	A	946	MET	3.5
1	A	1113	PHE	3.4
1	A	1184	ASN	3.3
1	A	876	ASP	3.2
1	A	1181	THR	3.1
1	A	1067	ASP	3.0
1	A	1017	ARG	2.9
1	A	986	SER	2.9
1	A	1220	ASN	2.8
1	A	984	SER	2.8
1	A	1065	ARG	2.8
1	A	945	ASP	2.8
1	A	947	PHE	2.8
1	A	939	MET	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	1111	ILE	2.6
1	A	1068	ASN	2.5
1	A	988	GLY	2.4
1	A	902	PHE	2.4
1	A	1233	GLY	2.3
1	A	1112	THR	2.3
1	A	1076	ARG	2.3
1	A	875	GLU	2.2
1	A	1203	ASN	2.2
1	A	1221	LEU	2.1
1	A	982	LYS	2.1

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(Å ²)	Q<0.9
2	A2G	A	602	15/15	0.71	0.17	68,70,71,71	0
2	A2G	A	601	15/15	0.80	0.11	52,55,58,59	0
3	SO4	A	701	5/5	0.86	0.16	56,58,58,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.