



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 08:50 AM UTC

PDB ID : 2F1A / pdb_00002f1a
Title : GOLGI ALPHA-MANNOSIDASE II COMPLEX WITH (2R,3R,4S)-2-({[(1S)-2-hydroxy-1-phenylethyl]amino}methyl)pyrrolidine-3,4-diol
Authors : Kuntz, D.A.; Rose, D.R.
Deposited on : 2005-11-14
Resolution : 1.45 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

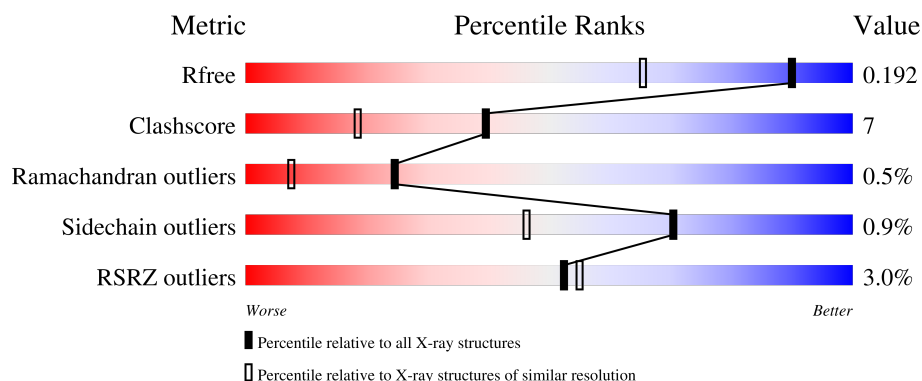
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

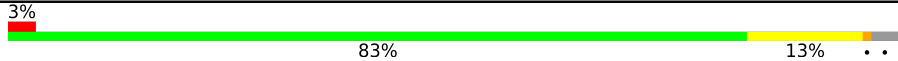
The reported resolution of this entry is 1.45 Å.

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	1756 (1.46-1.46)
Clashscore	190562	1795 (1.46-1.46)
Ramachandran outliers	187476	1776 (1.46-1.46)
Sidechain outliers	187428	1776 (1.46-1.46)
RSRZ outliers	180081	1756 (1.46-1.46)

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	1045	

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
6	MPD	A	1801	X	-	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 9452 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.

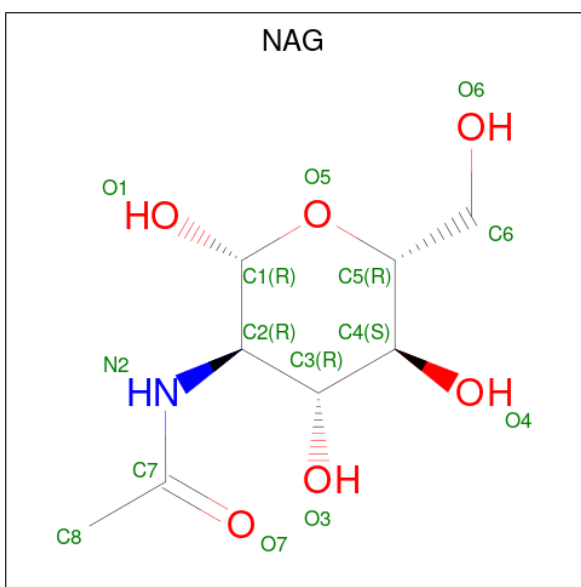
- Molecule 1 is a protein called alpha-mannosidase II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1015	Total	C	N	O	S	0	29	0
			8307	5277	1455	1532	43			

There are 12 discrepancies between the modelled and reference sequences:

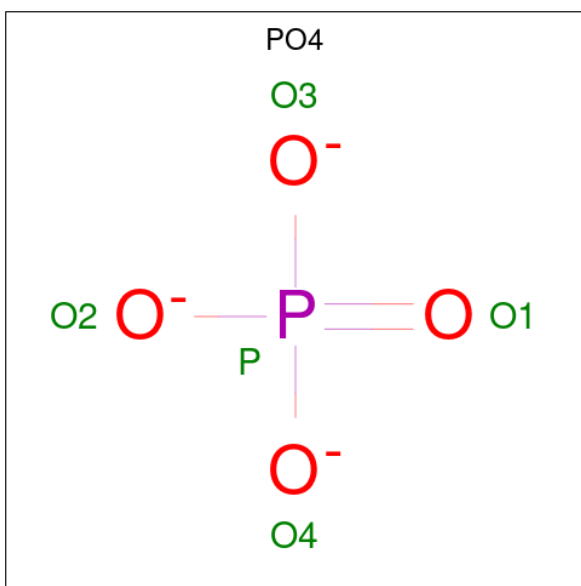
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	ARG	-	expression tag	GB 517481
A	2	SER	-	expression tag	GB 517481
A	3	SER	-	expression tag	GB 517481
A	4	HIS	-	expression tag	GB 517481
A	5	HIS	-	expression tag	GB 517481
A	6	HIS	-	expression tag	GB 517481
A	7	HIS	-	expression tag	GB 517481
A	8	HIS	-	expression tag	GB 517481
A	9	HIS	-	expression tag	GB 517481
A	10	GLY	-	expression tag	GB 517481
A	11	GLU	-	expression tag	GB 517481
A	12	PHE	-	expression tag	GB 517481

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).

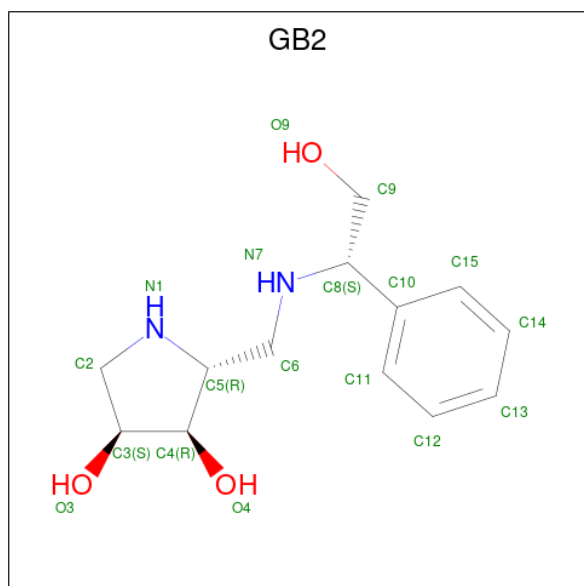


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

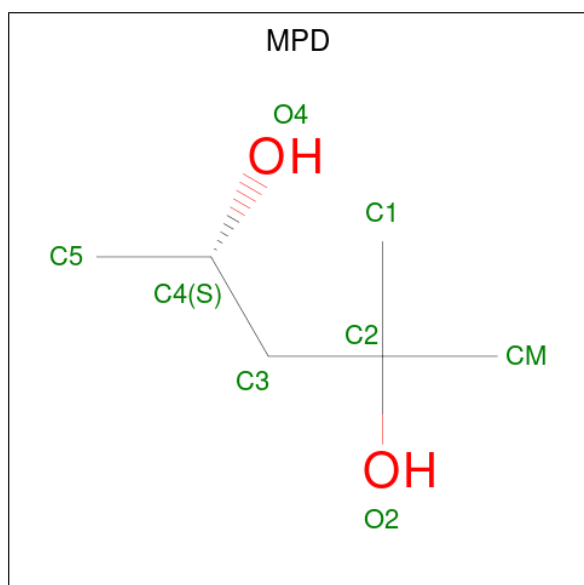
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	Zn		0	0
			1	1			

- Molecule 5 is (2R,3R,4S)-2-([(1S)-2-HYDROXY-1-PHENYLETHYL]AMINO}METHYL) PYRROLIDINE-3,4-DIOL (CCD ID: GB2) (formula: $C_{13}H_{20}N_2O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			18	13	2	3		

- Molecule 6 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: $C_6H_{14}O_2$).



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

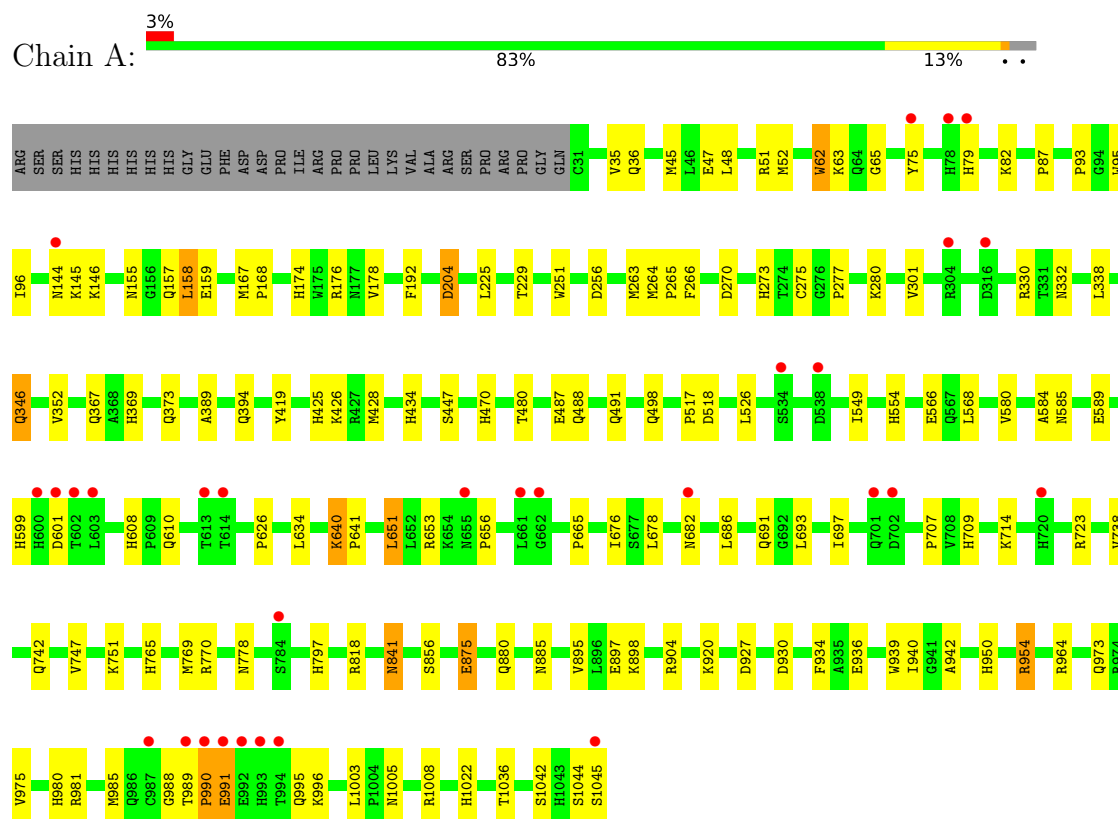
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1099	Total	O	0	0
			1099	1099		

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: alpha-mannosidase II



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.90Å 109.36Å 138.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.78 – 1.45 29.78 – 1.45	Depositor EDS
% Data completeness (in resolution range)	93.9 (29.78-1.45) 94.1 (29.78-1.45)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 1.45Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.168 , 0.189 0.173 , 0.192	Depositor DCC
R_{free} test set	4146 reflections (2.24%)	wwPDB-VP
Wilson B-factor (Å ²)	16.3	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9452	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.43% 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: NAG, ZN, MPD, GB2, PO4

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	0.81	0/8665	1.04	30/11759 (0.3%)

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	2

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	517	PRO	N-CA-C	7.76	123.38	111.19
1	A	665	PRO	N-CA-C	7.52	123.22	114.20
1	A	168	PRO	N-CA-C	6.82	122.22	111.38
1	A	518	ASP	N-CA-C	-6.60	96.31	107.99
1	A	35	VAL	N-CA-C	6.48	117.15	111.90
1	A	265	PRO	N-CA-C	6.32	125.49	112.47
1	A	45	MET	N-CA-C	6.27	118.92	111.71
1	A	584	ALA	N-CA-C	-6.16	105.35	112.92
1	A	35	VAL	CB-CA-C	-5.94	105.60	111.30
1	A	640	LYS	CA-C-N	5.82	125.73	119.85
1	A	640	LYS	C-N-CA	5.82	125.73	119.85
1	A	277	PRO	N-CA-C	5.82	121.18	114.03
1	A	192	PHE	N-CA-C	5.72	120.39	113.41
1	A	330	ARG	N-CA-C	5.67	120.26	113.23
1	A	566	GLU	N-CA-C	-5.56	101.11	109.95
1	A	682	ASN	N-CA-C	-5.49	106.16	112.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	266	PHE	N-CA-C	5.44	117.79	109.95
1	A	875	GLU	N-CA-C	5.34	119.08	112.24
1	A	447[A]	SER	N-CA-C	-5.33	101.01	109.76
1	A	447[B]	SER	N-CA-C	-5.33	101.01	109.76
1	A	36	GLN	N-CA-C	5.29	119.74	113.28
1	A	738	VAL	N-CA-C	-5.26	102.79	109.58
1	A	63	LYS	N-CA-C	5.25	117.75	111.71
1	A	62	TRP	N-CA-C	-5.25	98.10	107.98
1	A	841	ASN	N-CA-C	5.23	120.31	113.88
1	A	178	VAL	N-CA-C	-5.19	105.33	110.62
1	A	167	MET	N-CA-C	-5.14	98.37	109.01
1	A	1044	SER	CA-C-N	5.07	130.82	121.70
1	A	1044	SER	C-N-CA	5.07	130.82	121.70
1	A	204	ASP	N-CA-C	5.00	120.86	109.81

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	709[B]	HIS	Mainchain
1	A	75[B]	TYR	Mainchain

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	8307	0	8086	105	0
2	A	14	0	13	0	0
3	A	5	0	0	0	0
4	A	1	0	0	0	0
5	A	18	0	18	3	0
6	A	8	0	14	3	0
7	A	1099	0	0	31	0
All	All	9452	0	8131	110	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 7.

All (110) 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:954:ARG:NH2	1:A:981:ARG:HH21	1.50	1.10
1:A:954:ARG:HH21	1:A:981:ARG:NH2	1.56	1.01
1:A:954:ARG:HH21	1:A:981:ARG:HH21	1.03	0.97
5:A:1804:GB2:H11	7:A:2904:HOH:O	1.65	0.95
1:A:989:THR:HG22	1:A:991:GLU:HG3	1.53	0.91
1:A:678:LEU:HD12	1:A:769[B]:MET:HE1	1.57	0.87
1:A:434:HIS:HE1	1:A:930:ASP:OD1	1.62	0.83
1:A:491[B]:GLN:HG3	7:A:2871:HOH:O	1.80	0.82
1:A:964:ARG:HH11	1:A:973:GLN:HE21	1.28	0.81
1:A:96:ILE:HG22	7:A:2351:HOH:O	1.81	0.79
1:A:742:GLN:HG3	7:A:2023:HOH:O	1.85	0.77
5:A:1804:GB2:H92	7:A:2460:HOH:O	1.85	0.77
1:A:498:GLN:HE21	1:A:526:LEU:H	1.35	0.74
1:A:693:LEU:HD13	7:A:2018:HOH:O	1.91	0.70
6:A:1801:MPD:H11	7:A:2827:HOH:O	1.90	0.70
1:A:653:ARG:HD2	1:A:656:PRO:HA	1.74	0.70
1:A:155:ASN:HD21	1:A:157:GLN:HE21	1.39	0.69
1:A:678:LEU:CD1	1:A:769[B]:MET:HE1	2.21	0.69
1:A:346:GLN:HE21	1:A:346:GLN:H	1.39	0.69
1:A:256:ASP:HB2	7:A:1969:HOH:O	1.91	0.69
1:A:144:ASN:HB3	7:A:2451:HOH:O	1.93	0.69
1:A:82:LYS:HD3	1:A:373:GLN:HB3	1.73	0.68
1:A:651:LEU:HD13	1:A:653:ARG:HG2	1.76	0.68
6:A:1801:MPD:H13	7:A:2534:HOH:O	1.94	0.68
1:A:714:LYS:HD2	7:A:2756:HOH:O	1.95	0.67
1:A:707:PRO:HG2	1:A:797[A]:HIS:CE1	2.30	0.66
1:A:434:HIS:HD2	1:A:927:ASP:OD1	1.78	0.66
1:A:589:GLU:OE2	1:A:751:LYS:HD3	1.96	0.65
1:A:904:ARG:HG2	1:A:985:MET:SD	2.37	0.64
1:A:954:ARG:NH2	1:A:981:ARG:NH2	2.24	0.63
1:A:225:LEU:HD21	1:A:264[B]:MET:HE2	1.80	0.62
1:A:950:HIS:HE1	7:A:2209:HOH:O	1.83	0.62
1:A:47:GLU:OE2	1:A:51:ARG:CD	2.47	0.62
1:A:273:HIS:HE1	6:A:1801:MPD:O4	1.84	0.60
5:A:1804:GB2:C11	7:A:2904:HOH:O	2.37	0.60
1:A:765:HIS:HE1	7:A:2901:HOH:O	1.85	0.60
1:A:818[B]:ARG:HD3	1:A:856:SER:O	2.03	0.59
1:A:480:THR:H	1:A:880:GLN:HE22	1.50	0.58
1:A:995:GLN:HG3	7:A:2636:HOH:O	2.03	0.58
1:A:980:HIS:HD2	1:A:1036:THR:OG1	1.87	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:980:HIS:HE1	7:A:1939:HOH:O	1.86	0.57
1:A:96:ILE:CG2	7:A:2351:HOH:O	2.47	0.57
1:A:626:PRO:O	1:A:950:HIS:HD2	1.87	0.57
1:A:651:LEU:CD1	1:A:653:ARG:HG2	2.35	0.57
1:A:82:LYS:HD2	7:A:1864:HOH:O	2.05	0.57
1:A:47:GLU:OE2	1:A:51:ARG:HD3	2.05	0.56
1:A:145:LYS:HE3	7:A:2420:HOH:O	2.06	0.54
1:A:93:PRO:HD2	1:A:470:HIS:CD2	2.42	0.54
1:A:723:ARG:HG2	1:A:723:ARG:HH11	1.73	0.54
1:A:765:HIS:HD2	1:A:778:ASN:OD1	1.90	0.54
1:A:934:PHE:CE2	1:A:936:GLU:HB2	2.43	0.54
1:A:954:ARG:HG2	1:A:954:ARG:NH1	2.21	0.53
1:A:47:GLU:OE2	1:A:51:ARG:HD2	2.09	0.53
1:A:895:VAL:HG12	1:A:897:GLU:HG3	1.89	0.53
1:A:174:HIS:CE1	1:A:176:ARG:HD3	2.43	0.53
1:A:62:TRP:CD2	1:A:65:GLY:HA3	2.45	0.52
1:A:146:LYS:NZ	7:A:2430:HOH:O	2.31	0.52
1:A:975:VAL:HG21	1:A:1003:LEU:CD1	2.39	0.52
1:A:936:GLU:HG2	7:A:2676:HOH:O	2.10	0.52
1:A:940:ILE:HD12	7:A:2684:HOH:O	2.09	0.52
1:A:841:ASN:O	1:A:898:LYS:HE2	2.10	0.51
1:A:954:ARG:CG	1:A:954:ARG:HH11	2.22	0.51
1:A:601:ASP:HB2	1:A:608:HIS:CE1	2.45	0.51
1:A:434:HIS:CE1	1:A:930:ASP:OD1	2.53	0.51
1:A:601:ASP:HB2	1:A:608:HIS:HE1	1.76	0.50
1:A:1005:ASN:OD1	1:A:1045:SER:HA	2.12	0.50
1:A:367:GLN:HB3	1:A:369[A]:HIS:NE2	2.27	0.50
1:A:428:MET:HE1	1:A:491[B]:GLN:HG2	1.94	0.49
1:A:79:HIS:HE1	7:A:2506:HOH:O	1.94	0.49
1:A:51:ARG:HG3	7:A:2850:HOH:O	2.12	0.49
1:A:875:GLU:CD	7:A:2336:HOH:O	2.56	0.48
1:A:270:ASP:OD1	1:A:273:HIS:HD2	1.96	0.48
1:A:251:TRP:C	1:A:251:TRP:CD1	2.91	0.48
1:A:954:ARG:HG2	1:A:954:ARG:HH11	1.78	0.48
1:A:425:HIS:HE1	1:A:487:GLU:OE1	1.96	0.48
1:A:920:LYS:HG3	7:A:2631:HOH:O	2.13	0.48
1:A:47:GLU:CD	1:A:51:ARG:HD3	2.39	0.48
1:A:1008:ARG:NH1	1:A:1022:HIS:NE2	2.62	0.47
1:A:488:GLN:HG3	7:A:2872:HOH:O	2.14	0.47
1:A:158[A]:LEU:HD23	1:A:159:GLU:N	2.30	0.47
1:A:498:GLN:NE2	1:A:526:LEU:H	2.08	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:367:GLN:HB3	1:A:369[A]:HIS:CD2	2.50	0.47
1:A:554:HIS:HD2	7:A:2682:HOH:O	1.99	0.46
1:A:580:VAL:HG22	1:A:634:LEU:HD22	1.98	0.46
1:A:989:THR:O	1:A:990:PRO:C	2.58	0.45
1:A:939:TRP:CD2	1:A:942:ALA:HB2	2.53	0.44
1:A:996:LYS:HE2	7:A:2217:HOH:O	2.16	0.44
1:A:964:ARG:HH11	1:A:973:GLN:NE2	2.07	0.43
1:A:988:GLY:O	1:A:990:PRO:HD3	2.19	0.42
1:A:895:VAL:HG12	1:A:897:GLU:CG	2.49	0.42
1:A:640:LYS:HA	1:A:641:PRO:HD3	1.86	0.42
1:A:332:ASN:H	1:A:394:GLN:NE2	2.18	0.42
1:A:280:LYS:HZ3	1:A:301:VAL:HG21	1.85	0.42
1:A:554:HIS:HE1	7:A:2086:HOH:O	2.02	0.42
1:A:229:THR:HG21	1:A:263:MET:HE2	2.02	0.41
1:A:389:ALA:O	1:A:394:GLN:HG2	2.20	0.41
1:A:48:LEU:HG	1:A:52[A]:MET:SD	2.61	0.41
1:A:599:HIS:HD2	1:A:610:GLN:CG	2.34	0.41
1:A:256:ASP:CB	7:A:1969:HOH:O	2.61	0.41
1:A:419:TYR:C	1:A:426:LYS:HE2	2.45	0.41
1:A:885:ASN:ND2	1:A:885:ASN:H	2.19	0.41
1:A:87:PRO:HA	1:A:338:LEU:O	2.21	0.40
1:A:568:LEU:HD12	1:A:770[A]:ARG:HD3	2.04	0.40
1:A:332:ASN:H	1:A:394:GLN:HE22	1.68	0.40
1:A:686:LEU:HD22	1:A:697:ILE:HG12	2.02	0.40
1:A:723:ARG:HG2	1:A:723:ARG:NH1	2.36	0.40
1:A:1008:ARG:HB3	1:A:1042:SER:HB2	2.03	0.40
1:A:676:ILE:HD13	1:A:747:VAL:HG21	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	1042/1045 (100%)	1007 (97%)	30 (3%)	5 (0%)	24 7

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	95	TRP
1	A	990	PRO
1	A	991	GLU
1	A	204	ASP
1	A	549	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	931/929 (100%)	921 (99%)	10 (1%)	65 38

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

Mol	Chain	Res	Type
1	A	158[A]	LEU
1	A	158[B]	LEU
1	A	275	CYS
1	A	346	GLN
1	A	352	VAL
1	A	585[A]	ASN
1	A	585[B]	ASN
1	A	651	LEU
1	A	691	GLN
1	A	954	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	HIS

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Mol	Chain	Res	Type
1	A	119	HIS
1	A	148	GLN
1	A	157	GLN
1	A	191	GLN
1	A	240	GLN
1	A	273	HIS
1	A	346	GLN
1	A	347	ASN
1	A	371	ASN
1	A	388	GLN
1	A	394	GLN
1	A	425	HIS
1	A	434	HIS
1	A	460	GLN
1	A	470	HIS
1	A	488	GLN
1	A	498	GLN
1	A	554	HIS
1	A	599	HIS
1	A	608	HIS
1	A	742	GLN
1	A	765	HIS
1	A	778	ASN
1	A	812	GLN
1	A	866	GLN
1	A	880	GLN
1	A	885	ASN
1	A	923	GLN
1	A	950	HIS
1	A	973	GLN
1	A	980	HIS
1	A	986	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

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$
5	GB2	A	1804	4	19,19,19	1.55	5 (26%)	20,25,25	2.31	7 (35%)
3	PO4	A	1803	-	4,4,4	1.74	1 (25%)	6,6,6	1.05	0
6	MPD	A	1801	-	7,7,7	0.85	0	9,10,10	0.43	0
2	NAG	A	1802	1	14,14,15	0.56	0	17,19,21	0.89	1 (5%)

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
6	MPD	A	1801	-	1/1/2/2	2/5/5/5	-
5	GB2	A	1804	4	-	9/11/24/24	0/2/2/2
2	NAG	A	1802	1	-	3/6/23/26	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1804	GB2	C15-C10	3.45	1.44	1.39
3	A	1803	PO4	P-O1	3.25	1.58	1.50
5	A	1804	GB2	C11-C10	2.98	1.43	1.39
5	A	1804	GB2	C14-C15	2.68	1.43	1.38
5	A	1804	GB2	C12-C11	2.25	1.42	1.38
5	A	1804	GB2	C10-C8	2.20	1.55	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1804	GB2	C6-N7-C8	6.58	126.25	114.00
5	A	1804	GB2	C13-C12-C11	3.80	124.93	120.24
5	A	1804	GB2	C5-C6-N7	-2.97	106.53	112.68
5	A	1804	GB2	C15-C10-C8	2.89	126.66	120.79
2	A	1802	NAG	C2-N2-C7	-2.88	119.04	122.90
5	A	1804	GB2	O4-C4-C5	-2.52	106.18	112.92
5	A	1804	GB2	C11-C10-C8	-2.40	115.93	120.79
5	A	1804	GB2	C12-C13-C14	-2.22	116.83	119.87

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
6	A	1801	MPD	C4

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1802	NAG	C8-C7-N2-C2
2	A	1802	NAG	O7-C7-N2-C2
5	A	1804	GB2	C10-C8-N7-C6
5	A	1804	GB2	C9-C8-N7-C6
5	A	1804	GB2	N7-C8-C9-O9
5	A	1804	GB2	N1-C5-C6-N7
6	A	1801	MPD	C2-C3-C4-O4
6	A	1801	MPD	C2-C3-C4-C5
5	A	1804	GB2	C15-C10-C8-C9
5	A	1804	GB2	C11-C10-C8-C9
5	A	1804	GB2	C15-C10-C8-N7
5	A	1804	GB2	C11-C10-C8-N7
2	A	1802	NAG	O5-C5-C6-O6
5	A	1804	GB2	C4-C5-C6-N7

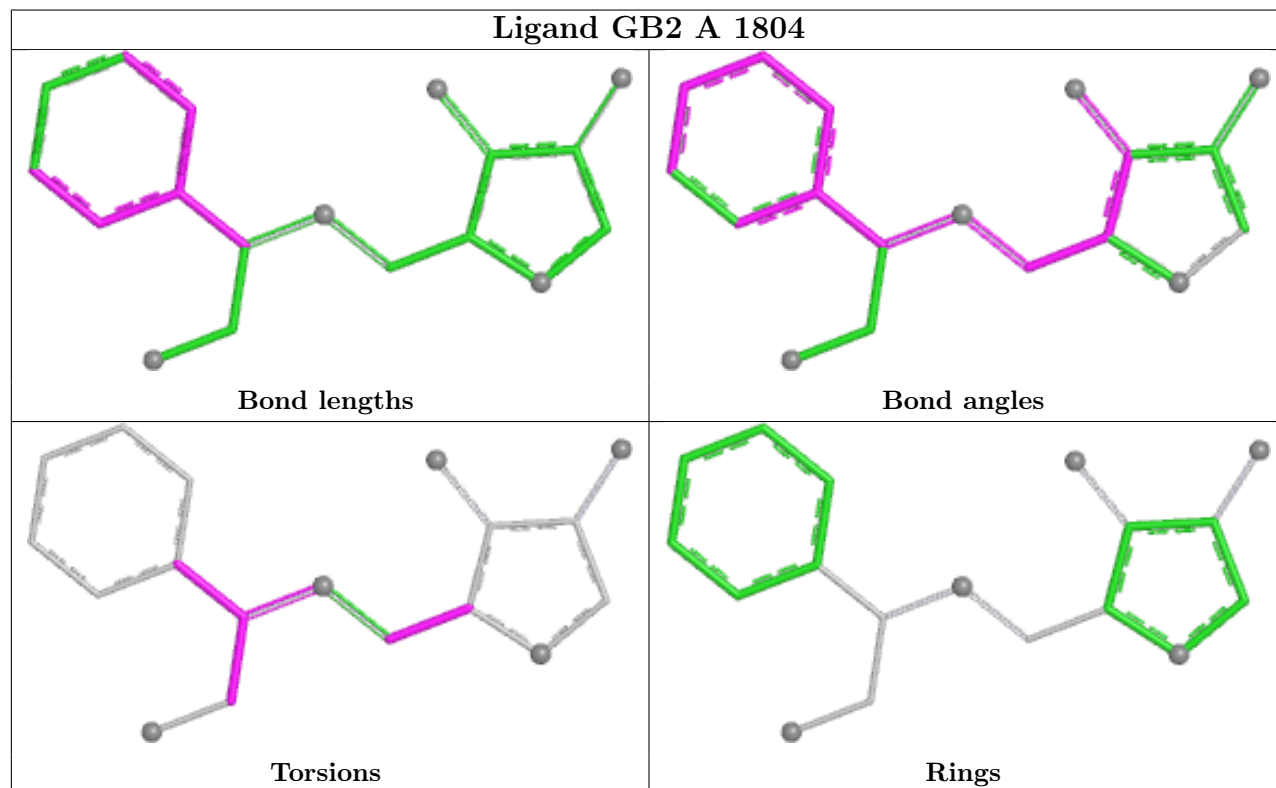
There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1804	GB2	3	0
6	A	1801	MPD	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	1015/1045 (97%)	-0.06	30 (2%) 52 55	8, 17, 32, 68	29 (2%)

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	990	PRO	5.7
1	A	993	HIS	5.3
1	A	1045	SER	4.8
1	A	534	SER	4.2
1	A	613	THR	3.5
1	A	992	GLU	3.4
1	A	144	ASN	3.0
1	A	602	THR	3.0
1	A	989	THR	3.0
1	A	991	GLU	2.8
1	A	600	HIS	2.6
1	A	614	THR	2.5
1	A	720	HIS	2.5
1	A	603	LEU	2.4
1	A	655	ASN	2.3
1	A	682	ASN	2.2
1	A	987	CYS	2.2
1	A	662	GLY	2.2
1	A	304	ARG	2.2
1	A	538	ASP	2.2
1	A	661	LEU	2.1
1	A	702	ASP	2.1
1	A	701	GLN	2.1
1	A	994	THR	2.0
1	A	75[A]	TYR	2.0
1	A	784	SER	2.0
1	A	601	ASP	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	78	HIS	2.0
1	A	79	HIS	2.0
1	A	316	ASP	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](#)

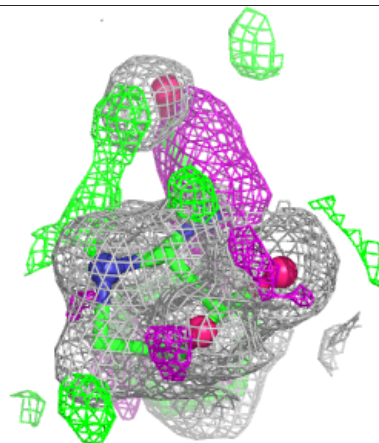
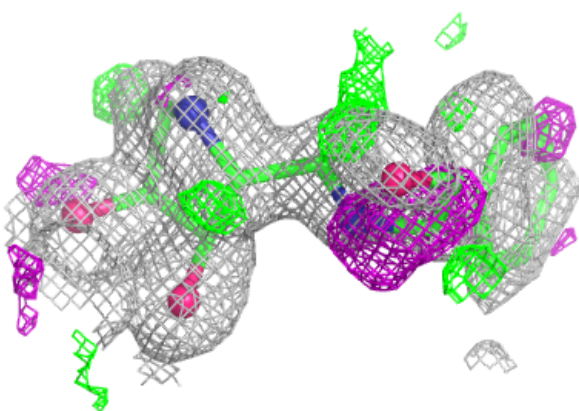
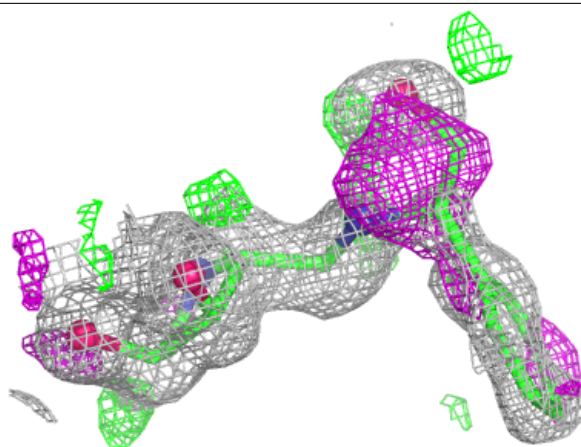
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	NAG	A	1802	14/15	0.65	0.14	40,47,52,54	0
5	GB2	A	1804	18/18	0.89	0.12	14,27,33,34	0
3	PO4	A	1803	5/5	0.90	0.11	32,36,38,38	0
6	MPD	A	1801	8/8	0.95	0.09	19,24,26,27	0
4	ZN	A	1805	1/1	1.00	0.01	13,13,13,13	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around GB2 A 1804:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.