



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2014 – 03:39 AM GMT

PDB ID : 1QX1  
Title : Golgi alpha-mannosidase II D341N mutant complex with 2-F-mannosyl-F  
Authors : Numao, S.; Kuntz, D.A.; Withers, S.G.; Rose, D.R.  
Deposited on : 2003-09-04  
Resolution : 1.30 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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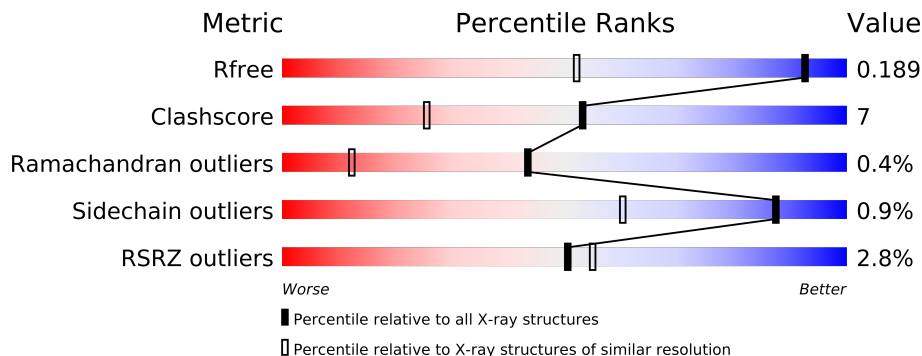
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance


The reported resolution of this entry is 1.30 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))
$R_{free}$	66092	1025 (1.34-1.26)
Clashscore	79885	1140 (1.34-1.26)
Ramachandran outliers	78287	1093 (1.34-1.26)
Sidechain outliers	78261	1092 (1.34-1.26)
RSRZ outliers	66119	1025 (1.34-1.26)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	1045	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	NAG	A	2001	-	X
4	FMF	A	2003[A]	-	X
4	FMF	A	2003[B]	-	X
5	MPD	A	2002	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9520 atoms, of which 0 are hydrogen and 0 are deuterium.

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
			Total	C	N	O	S			
1	A	1014	8433	5356	1482	1552	43	0	29	0

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	ARG	-	CLONING ARTIFACT	UNP Q24451
A	2	SER	-	CLONING ARTIFACT	UNP Q24451
A	3	SER	-	CLONING ARTIFACT	UNP Q24451
A	4	HIS	-	EXPRESSION TAG	UNP Q24451
A	5	HIS	-	EXPRESSION TAG	UNP Q24451
A	6	HIS	-	EXPRESSION TAG	UNP Q24451
A	7	HIS	-	EXPRESSION TAG	UNP Q24451
A	8	HIS	-	EXPRESSION TAG	UNP Q24451
A	9	HIS	-	EXPRESSION TAG	UNP Q24451
A	10	GLY	-	CLONING ARTIFACT	UNP Q24451
A	11	GLU	-	CLONING ARTIFACT	UNP Q24451
A	12	PHE	-	CLONING ARTIFACT	UNP Q24451
A	341	ASN	ASP	ENGINEERED	UNP Q24451
A	907	LYS	GLU	SEE REMARK 999	UNP Q24451

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).

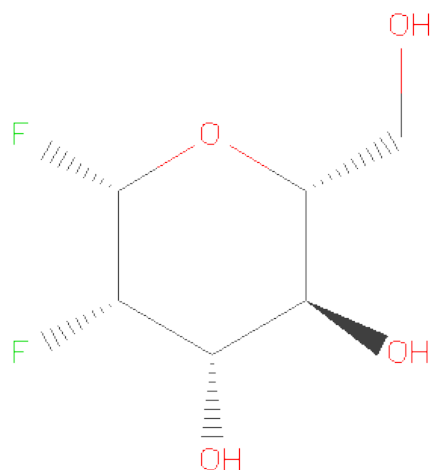


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

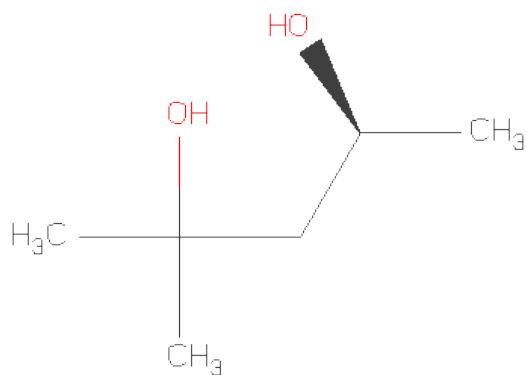
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is 2-DEOXY-2-FLUOROHXOPYRANOSYLFLUORIDE (three-letter code: FMF) (formula: C<sub>6</sub>H<sub>10</sub>F<sub>2</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	F	O	0	1
			22	12	2	8		

- Molecule 5 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>2</sub>).



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

- Molecule 6 is water.

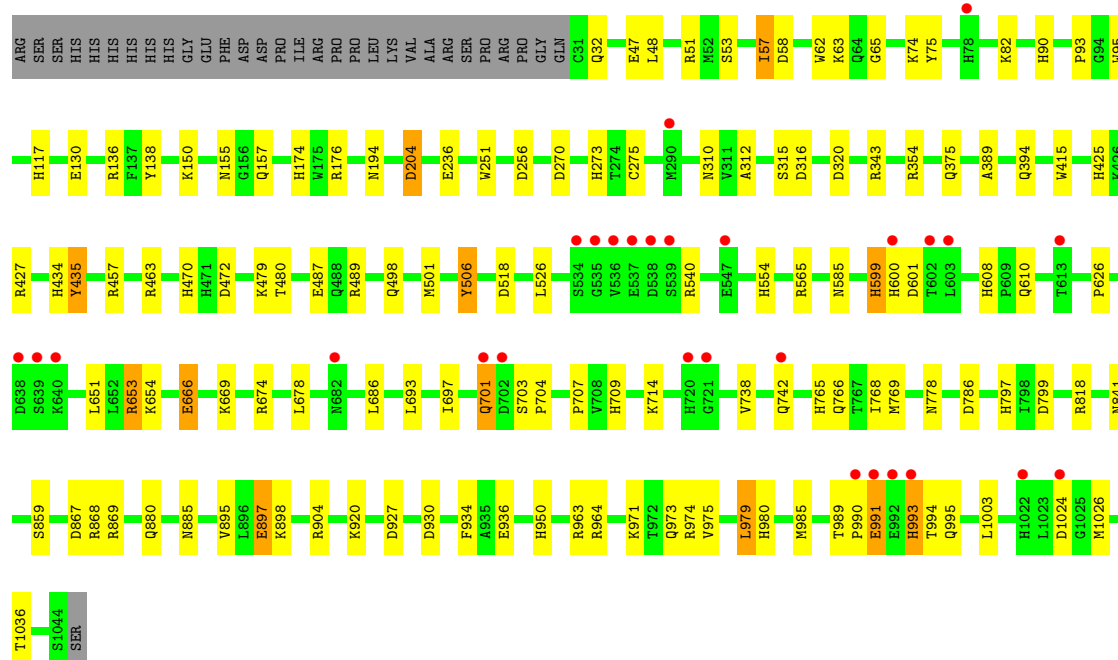
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1042	Total	O	0	0
			1042	1042		

### 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 errors 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 ( $\text{RSRZ} > 2$ ). 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

Chain A: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	69.05Å 109.83Å 138.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.23 – 1.30 29.86 – 1.30	Depositor EDS
% Data completeness (in resolution range)	92.7 (29.23-1.30) 92.9 (29.86-1.30)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.93 (at 1.30Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.170 , 0.189 0.170 , 0.189	Depositor DCC
$R_{free}$ test set	2180 reflections (0.92%)	DCC
Wilson B-factor (Å <sup>2</sup> )	11.5	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 40.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 257224 reflections	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	9520	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.03% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FMF, MPD, ZN, NAG

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	1.01	7/8666 (0.1%)	1.03	28/11760 (0.2%)

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

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	501	MET	SD-CE	-6.79	1.39	1.77
1	A	506[A]	TYR	CD2-CE2	6.00	1.48	1.39
1	A	506[B]	TYR	CD2-CE2	6.00	1.48	1.39
1	A	312	ALA	CA-CB	5.72	1.64	1.52
1	A	859	SER	CB-OG	-5.25	1.35	1.42
1	A	415	TRP	CG-CD1	5.09	1.43	1.36
1	A	897	GLU	CG-CD	5.07	1.59	1.51

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	427	ARG	NE-CZ-NH1	7.66	124.13	120.30
1	A	979[A]	LEU	CA-CB-CG	6.93	131.25	115.30
1	A	979[B]	LEU	CA-CB-CG	6.93	131.25	115.30
1	A	501	MET	CG-SD-CE	-6.93	89.12	100.20
1	A	963	ARG	NE-CZ-NH2	-6.54	117.03	120.30
1	A	427	ARG	NE-CZ-NH2	-6.29	117.15	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	565	ARG	NE-CZ-NH2	-6.25	117.17	120.30
1	A	457	ARG	NE-CZ-NH2	-6.23	117.19	120.30
1	A	343	ARG	NE-CZ-NH2	-6.01	117.30	120.30
1	A	974	ARG	NE-CZ-NH2	-5.99	117.30	120.30
1	A	674	ARG	NE-CZ-NH2	-5.97	117.31	120.30
1	A	818	ARG	NE-CZ-NH1	5.88	123.24	120.30
1	A	540	ARG	NE-CZ-NH1	5.82	123.21	120.30
1	A	320	ASP	CB-CG-OD2	-5.67	113.20	118.30
1	A	463	ARG	NE-CZ-NH1	-5.54	117.53	120.30
1	A	518	ASP	N-CA-C	-5.46	96.25	111.00
1	A	63	LYS	CD-CE-NZ	5.45	124.24	111.70
1	A	506[A]	TYR	CB-CG-CD1	-5.44	117.74	121.00
1	A	506[B]	TYR	CB-CG-CD1	-5.44	117.74	121.00
1	A	489	ARG	NE-CZ-NH2	-5.41	117.59	120.30
1	A	867	ASP	CB-CG-OD1	5.41	123.17	118.30
1	A	786	ASP	CB-CG-OD2	5.25	123.03	118.30
1	A	993	HIS	N-CA-C	5.20	125.04	111.00
1	A	818	ARG	NE-CZ-NH2	-5.11	117.74	120.30
1	A	868	ARG	NE-CZ-NH2	-5.11	117.75	120.30
1	A	868	ARG	NE-CZ-NH1	5.08	122.84	120.30
1	A	457	ARG	NE-CZ-NH1	5.06	122.83	120.30
1	A	799	ASP	CB-CG-OD2	5.05	122.84	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	435[B]	TYR	Mainchain
1	A	599[B]	HIS	Mainchain

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	8433	0	8183	112	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	14	0	13	7	0
3	A	1	0	0	0	0
4	A	22	0	15	2	0
5	A	8	0	14	3	0
6	A	1042	0	0	42	1
All	All	9520	0	8225	117	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

All (117) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:194:ASN:HD21	2:A:2001:NAG:C1	1.25	1.46
1:A:435[B]:TYR:CE1	6:A:2594:HOH:O	1.80	1.30
1:A:75[B]:TYR:CE1	6:A:2992:HOH:O	2.02	1.12
1:A:707:PRO:HG2	1:A:797[B]:HIS:CE1	1.87	1.09
1:A:82[A]:LYS:HD2	1:A:375:GLN:HE22	1.24	1.03
1:A:506[B]:TYR:OH	6:A:2896:HOH:O	1.55	1.02
1:A:75[B]:TYR:HD1	6:A:2781:HOH:O	1.44	0.99
1:A:435[B]:TYR:HE1	6:A:2594:HOH:O	1.30	0.95
1:A:709[A]:HIS:CD2	6:A:3045:HOH:O	2.16	0.95
1:A:707:PRO:HG2	1:A:797[B]:HIS:HE1	1.27	0.94
1:A:117[B]:HIS:HE1	1:A:354:ARG:HE	1.08	0.93
1:A:117[B]:HIS:CE1	1:A:354:ARG:HE	1.86	0.91
1:A:47:GLU:OE1	1:A:51:ARG:NH1	2.03	0.91
1:A:707:PRO:CG	1:A:797[B]:HIS:CE1	2.54	0.91
1:A:707:PRO:CG	1:A:797[B]:HIS:HE1	1.82	0.90
1:A:204:ASP:OD1	4:A:2003[A]:FMF:C1	2.17	0.90
1:A:75[B]:TYR:HE1	6:A:2992:HOH:O	1.43	0.88
1:A:707:PRO:CB	1:A:797[B]:HIS:CE1	2.58	0.87
1:A:82[A]:LYS:HD2	1:A:375:GLN:NE2	1.95	0.81
1:A:155:ASN:HD21	1:A:157:GLN:HE21	1.27	0.80
1:A:989:THR:HG22	1:A:991:GLU:HG3	1.66	0.78
1:A:136:ARG:HD3	6:A:2887:HOH:O	1.83	0.78
1:A:434:HIS:HE1	1:A:930:ASP:OD1	1.66	0.78
1:A:971[B]:LYS:HB2	1:A:971[B]:LYS:NZ	1.99	0.77
1:A:964:ARG:HH11	1:A:973:GLN:HE21	1.32	0.77
1:A:75[B]:TYR:CD1	6:A:2781:HOH:O	2.26	0.76
1:A:693:LEU:HD13	6:A:3029:HOH:O	1.85	0.76
1:A:707:PRO:HB2	1:A:797[B]:HIS:CE1	2.21	0.76
1:A:989:THR:CG2	1:A:991:GLU:HG3	2.16	0.76
1:A:117[B]:HIS:HE1	1:A:354:ARG:NE	1.82	0.74

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:669[B]:LYS:CD	1:A:669[B]:LYS:NZ	2.51	0.73
1:A:194:ASN:HD21	2:A:2001:NAG:C2	2.01	0.73
1:A:498:GLN:HE21	1:A:526:LEU:H	1.37	0.72
1:A:971[B]:LYS:HZ3	1:A:971[B]:LYS:HB2	1.54	0.71
1:A:256:ASP:HB2	6:A:3023:HOH:O	1.90	0.70
1:A:130:GLU:OE2	6:A:2891:HOH:O	2.10	0.69
1:A:82[B]:LYS:HE3	6:A:2745:HOH:O	1.93	0.69
1:A:82[B]:LYS:CE	6:A:2745:HOH:O	2.40	0.68
2:A:2001:NAG:N2	6:A:2914:HOH:O	2.27	0.68
5:A:2002:MPD:H13	6:A:2693:HOH:O	1.95	0.67
5:A:2002:MPD:H11	6:A:2844:HOH:O	1.96	0.65
1:A:434:HIS:HD2	1:A:927:ASP:OD1	1.79	0.65
1:A:194:ASN:ND2	2:A:2001:NAG:N2	2.45	0.65
1:A:75[B]:TYR:OH	6:A:3042:HOH:O	2.14	0.64
1:A:742:GLN:HG3	6:A:2682:HOH:O	1.97	0.64
1:A:990:PRO:O	1:A:991:GLU:O	2.17	0.63
1:A:765:HIS:HE1	6:A:2611:HOH:O	1.81	0.62
1:A:980:HIS:HE1	6:A:2083:HOH:O	1.83	0.61
1:A:707:PRO:CB	1:A:797[B]:HIS:HE1	2.05	0.61
1:A:703:SER:HB2	1:A:704:PRO:HD2	1.82	0.61
1:A:194:ASN:ND2	2:A:2001:NAG:C2	2.62	0.60
1:A:709[A]:HIS:HD2	6:A:3045:HOH:O	1.62	0.60
1:A:701:GLN:HE21	1:A:701:GLN:HA	1.66	0.60
1:A:904:ARG:HG2	1:A:985:MET:SD	2.43	0.59
1:A:651:LEU:CD1	1:A:653:ARG:HG2	2.31	0.59
1:A:980:HIS:HD2	1:A:1036:THR:OG1	1.86	0.58
1:A:950:HIS:HE1	6:A:2616:HOH:O	1.88	0.57
1:A:480:THR:H	1:A:880:GLN:HE22	1.53	0.57
1:A:74:LYS:HE3	6:A:3042:HOH:O	2.05	0.56
1:A:990:PRO:O	1:A:991:GLU:C	2.43	0.56
1:A:626:PRO:O	1:A:950:HIS:HD2	1.88	0.56
1:A:273:HIS:HE1	5:A:2002:MPD:O4	1.88	0.55
1:A:57:ILE:HD12	1:A:58:ASP:N	2.22	0.54
1:A:270:ASP:OD1	1:A:273:HIS:HD2	1.90	0.54
1:A:554:HIS:HD2	6:A:2799:HOH:O	1.91	0.54
1:A:425:HIS:HE1	1:A:487:GLU:OE1	1.91	0.54
1:A:765:HIS:HD2	1:A:778:ASN:OD1	1.91	0.53
1:A:994:THR:HA	6:A:2575:HOH:O	2.08	0.53
1:A:975:VAL:HG21	1:A:1003:LEU:CD1	2.39	0.53
1:A:599[A]:HIS:CD2	1:A:610:GLN:HG3	2.44	0.53
1:A:895:VAL:HG12	1:A:897:GLU:HG3	1.92	0.52
1:A:601:ASP:OD2	1:A:608:HIS:HE1	1.93	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:701:GLN:HE21	1:A:701:GLN:CA	2.22	0.52
1:A:654:LYS:HG3	1:A:742:GLN:NE2	2.25	0.52
1:A:174:HIS:CE1	1:A:176:ARG:HD3	2.44	0.51
1:A:766[B]:GLN:HG2	1:A:768:ILE:HG13	1.91	0.51
1:A:554:HIS:HE1	6:A:2803:HOH:O	1.93	0.51
1:A:62:TRP:CD2	1:A:65:GLY:HA3	2.46	0.50
1:A:599[A]:HIS:HD2	1:A:610:GLN:CG	2.24	0.50
1:A:678:LEU:HD12	1:A:769[A]:MET:HE1	1.93	0.50
1:A:434:HIS:CE1	1:A:930:ASP:OD1	2.57	0.50
1:A:600:HIS:N	1:A:600:HIS:ND1	2.59	0.50
1:A:57:ILE:HD12	1:A:58:ASP:O	2.11	0.49
1:A:251:TRP:CD1	1:A:251:TRP:C	2.85	0.49
1:A:599[A]:HIS:HD2	1:A:610:GLN:HG3	1.78	0.49
1:A:48:LEU:HD11	1:A:236:GLU:HG2	1.95	0.48
1:A:82[A]:LYS:NZ	6:A:2822:HOH:O	2.47	0.48
2:A:2001:NAG:H2	6:A:2914:HOH:O	2.13	0.48
1:A:155:ASN:HB3	6:A:2363:HOH:O	2.12	0.48
1:A:435[B]:TYR:CZ	6:A:2594:HOH:O	2.26	0.48
1:A:869:ARG:HH11	1:A:885:ASN:HD22	1.62	0.47
1:A:989:THR:HG21	1:A:991:GLU:HG3	1.96	0.47
1:A:498:GLN:NE2	1:A:526:LEU:H	2.10	0.46
1:A:714:LYS:HE3	1:A:738:VAL:HG22	1.97	0.46
1:A:389:ALA:HB1	1:A:394:GLN:CG	2.45	0.46
1:A:666[A]:GLU:CG	6:A:2198:HOH:O	2.64	0.46
1:A:608:HIS:HD2	6:A:2841:HOH:O	1.99	0.45
1:A:47:GLU:CD	1:A:51:ARG:HH11	2.21	0.44
1:A:138:TYR:OH	1:A:150:LYS:HE2	2.17	0.44
1:A:707:PRO:HG2	1:A:797[B]:HIS:ND1	2.28	0.43
1:A:934:PHE:CE2	1:A:936:GLU:HB2	2.54	0.43
2:A:2001:NAG:C2	6:A:2914:HOH:O	2.66	0.43
1:A:686:LEU:HD22	1:A:697:ILE:HG12	2.00	0.43
1:A:32:GLN:NE2	6:A:3023:HOH:O	2.51	0.42
1:A:701:GLN:NE2	1:A:701:GLN:HA	2.33	0.42
1:A:93:PRO:HD2	1:A:470:HIS:CD2	2.54	0.42
1:A:600:HIS:CE1	6:A:2660:HOH:O	2.73	0.42
1:A:920:LYS:HG3	6:A:2990:HOH:O	2.20	0.41
1:A:472:ASP:OD2	4:A:2003[B]:FMF:O3	2.38	0.41
1:A:599[B]:HIS:CE1	1:A:601:ASP:HA	2.54	0.41
1:A:310:ASN:HA	6:A:2824:HOH:O	2.21	0.41
1:A:315:SER:O	1:A:316[B]:ASP:C	2.57	0.41
1:A:90:HIS:HA	6:A:2891:HOH:O	2.21	0.40
1:A:1026[A]:MET:HE3	6:A:2720:HOH:O	2.20	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:479:LYS:HD3	6:A:2836:HOH:O	2.20	0.40
1:A:995:GLN:HG3	6:A:2575:HOH:O	2.22	0.40
1:A:841:ASN:O	1:A:898[A]:LYS:HE2	2.20	0.40

All (1) 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	Distance(Å)	Clash(Å)
1:A:53[B]:SER:OG	6:A:2448:HOH:O[1_455]	2.19	0.01

## 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	1041/1045 (100%)	1010 (97%)	27 (3%)	4 (0%)	43 12

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	991	GLU
1	A	95	TRP
1	A	993	HIS
1	A	204	ASP

### 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	930/929 (100%)	920 (99%)	10 (1%)	84 52

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

Mol	Chain	Res	Type
1	A	57	ILE
1	A	275	CYS
1	A	585	ASN
1	A	653	ARG
1	A	666[A]	GLU
1	A	666[B]	GLU
1	A	701	GLN
1	A	979[A]	LEU
1	A	979[B]	LEU
1	A	1024	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	HIS
1	A	91	ASN
1	A	121	ASN
1	A	148	GLN
1	A	157	GLN
1	A	194	ASN
1	A	240	GLN
1	A	249	GLN
1	A	273	HIS
1	A	347	ASN
1	A	388	GLN
1	A	394	GLN
1	A	425	HIS
1	A	434	HIS
1	A	460	GLN
1	A	469	GLN
1	A	470	HIS
1	A	488	GLN
1	A	498	GLN
1	A	554	HIS
1	A	608	HIS
1	A	655	ASN
1	A	701	GLN
1	A	742	GLN
1	A	765	HIS
1	A	809	ASN
1	A	812	GLN

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Mol	Chain	Res	Type
1	A	880	GLN
1	A	885	ASN
1	A	901	ASN
1	A	919	HIS
1	A	950	HIS
1	A	973	GLN
1	A	980	HIS
1	A	986	GLN
1	A	1018	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

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	$\# Z  > 2$
2	NAG	A	2001	1	12,14,15	0.52	0	15,19,21	0.87	0
5	MPD	A	2002	-	7,7,7	0.77	0	10,10,10	0.62	0
4	FMF	A	2003[A]	3	9,11,12	4.24	4 (44%)	11,15,17	3.46	5 (45%)
4	FMF	A	2003[B]	1,3	9,11,12	3.78	4 (44%)	11,15,17	2.23	3 (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
2	NAG	A	2001	1	1/1/5/7	0/6/23/26	1/1/1/1
5	MPD	A	2002	-	1/1/2/2	0/5/5/5	0/0/0/0
4	FMF	A	2003[A]	3	-	0/2/19/22	0/1/1/1
4	FMF	A	2003[B]	1,3	-	0/2/19/22	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	2003[A]	FMF	C2-C3	11.46	1.61	1.51
4	A	2003[B]	FMF	C2-C3	10.05	1.60	1.51
4	A	2003[B]	FMF	C4-C3	3.47	1.61	1.52
4	A	2003[A]	FMF	O4-C4	-3.27	1.35	1.43
4	A	2003[B]	FMF	O3-C3	3.01	1.50	1.43
4	A	2003[A]	FMF	O-C5	-2.19	1.41	1.45
4	A	2003[A]	FMF	O3-C3	2.12	1.48	1.43
4	A	2003[B]	FMF	O4-C4	-2.08	1.37	1.43

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	2003[A]	FMF	C2-C3-C4	7.99	119.34	109.36
4	A	2003[A]	FMF	C3-C4-C5	-5.67	100.07	110.20
4	A	2003[B]	FMF	C3-C4-C5	-4.52	102.13	110.20
4	A	2003[B]	FMF	C2-C3-C4	4.50	114.97	109.36
4	A	2003[A]	FMF	F2-C2-C3	4.14	111.85	108.55
4	A	2003[A]	FMF	C6-C5-C4	-2.85	106.11	113.00
4	A	2003[B]	FMF	F2-C2-C3	2.76	110.75	108.55
4	A	2003[A]	FMF	O4-C4-C5	2.37	115.54	109.28

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	A	2002	MPD	C4
2	A	2001	NAG	C1

There are no torsion outliers.

All (1) ring outliers are listed below:



Mol	Chain	Res	Type	Atoms
2	A	2001	NAG	C1-C2-C3-C4-C5-O5

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	1014/1045 (97%)	-0.15	28 (2%) 50 55	7, 13, 27, 49	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	702	ASP	6.0
1	A	993	HIS	5.9
1	A	701	GLN	4.9
1	A	992	GLU	4.4
1	A	638	ASP	4.2
1	A	78	HIS	4.2
1	A	720	HIS	3.9
1	A	991	GLU	3.7
1	A	603	LEU	3.6
1	A	536	VAL	3.4
1	A	602	THR	3.4
1	A	600	HIS	3.2
1	A	547	GLU	3.2
1	A	538[A]	ASP	3.0
1	A	1024	ASP	3.0
1	A	537	GLU	2.8
1	A	990	PRO	2.7
1	A	539	SER	2.7
1	A	290	MET	2.6
1	A	534	SER	2.6
1	A	721	GLY	2.6
1	A	682	ASN	2.4
1	A	742	GLN	2.3
1	A	613	THR	2.3
1	A	1022	HIS	2.2
1	A	535	GLY	2.1
1	A	640	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	639	SER	2.1

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	NAG	A	2001	14/15	0.36	8.81	35,47,50,55	0
4	FMF	A	2003[A]	11/12	0.11	6.11	6,12,16,16	11
4	FMF	A	2003[B]	11/12	0.11	6.11	11,14,18,21	11
5	MPD	A	2002	8/8	0.12	2.40	14,17,22,25	0
3	ZN	A	2004	1/1	0.04	-0.95	12,12,12,12	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.