



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 03:04 am GMT

PDB ID : 1IVG
Title : STRUCTURES OF AROMATIC INHIBITORS OF INFLUENZA VIRUS
NEURAMINIDASE
Authors : Jedrzejewski, M.J.; Luo, M.
Deposited on : 1994-12-12
Resolution : 1.90 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

<http://wwpdb.org/validation/2016/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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

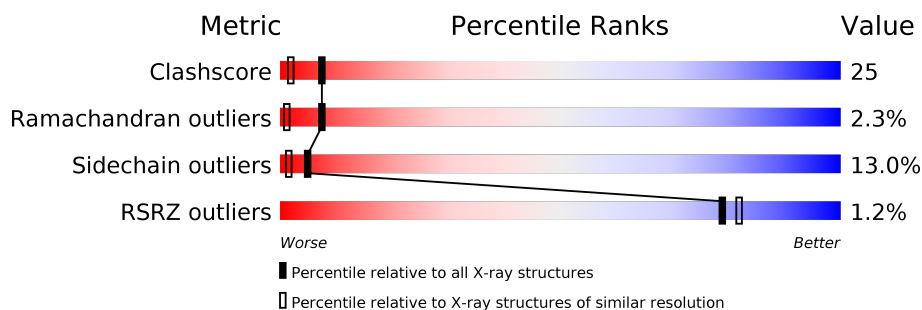
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	112137	5731 (1.90-1.90)
Ramachandran outliers	110173	5669 (1.90-1.90)
Sidechain outliers	110143	5670 (1.90-1.90)
RSRZ outliers	101464	5100 (1.90-1.90)

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. 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	388	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> % <div style="width: 100%; height: 10px; background-color: green;"></div> 61% 31% 7% . </div> </div>
1	B	388	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> % <div style="width: 100%; height: 10px; background-color: green;"></div> 63% 29% 7% . </div> </div>

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
4	NAG	B	478	-	-	X	X
4	MAN	B	481	-	-	X	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	MAN	B	482	-	-	-	X
4	MAN	B	483	-	-	X	X
4	MAN	B	487	-	-	X	X
4	MAN	B	488	-	-	X	X
4	MAN	B	489	-	-	-	X
5	FUC	B	477	X	-	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 8557 atoms, of which 2032 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 INFLUENZA A SUBTYPE N2 NEURAMINIDASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	388	Total	C	H	N	O	S	0	0	0
			3745	1866	723	545	588	23			
1	B	388	Total	C	H	N	O	S	0	0	0
			3745	1866	723	545	588	23			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	339	ASP	ASN	CONFLICT	UNP P06820
B	339	ASP	ASN	CONFLICT	UNP P06820

- Molecule 2 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	2	Total	C	H	N	O	0	0
			55	16	27	2	10		
2	A	2	Total	C	H	N	O	0	0
			55	16	27	2	10		
2	B	2	Total	C	H	N	O	0	0
			55	16	27	2	10		
2	B	2	Total	C	H	N	O	0	0
			55	16	27	2	10		

- Molecule 3 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	4	Total	C	H	N	O	0	0
			96	28	47	2	19		

- Molecule 4 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	6	Total	C	H	N	O	0	0
			139	40	67	2	30		
4	B	6	Total	C	H	N	O	0	0
			139	40	67	2	30		

- Molecule 5 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	4	Total	C	H	N	O	0	0
			96	28	47	2	19		

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Ca	0	0
			1	1		
6	A	1	Total	Ca	0	0
			1	1		

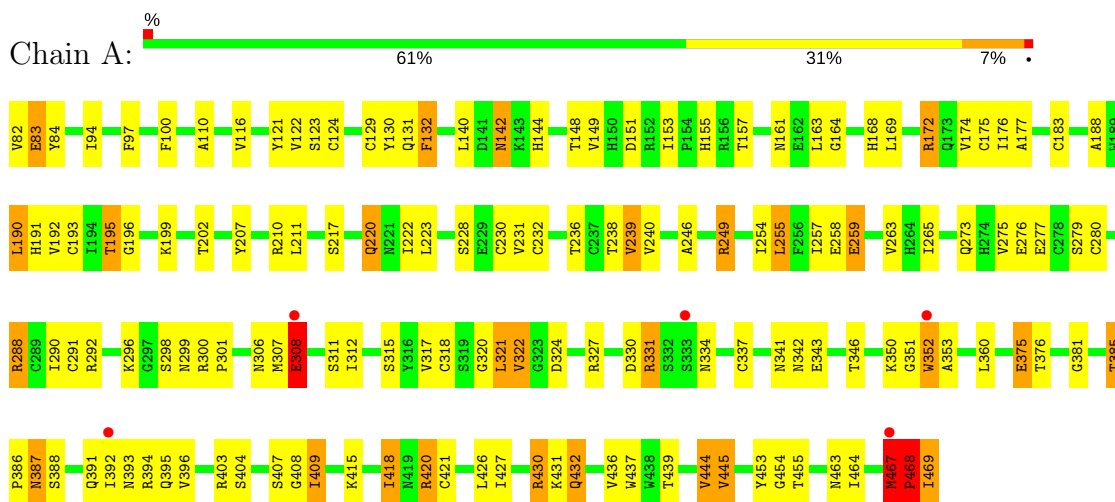
- Molecule 7 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	120	Total	H	O	0	0
			360	240	120		
7	B	5	Total	H	O	0	0
			15	10	5		

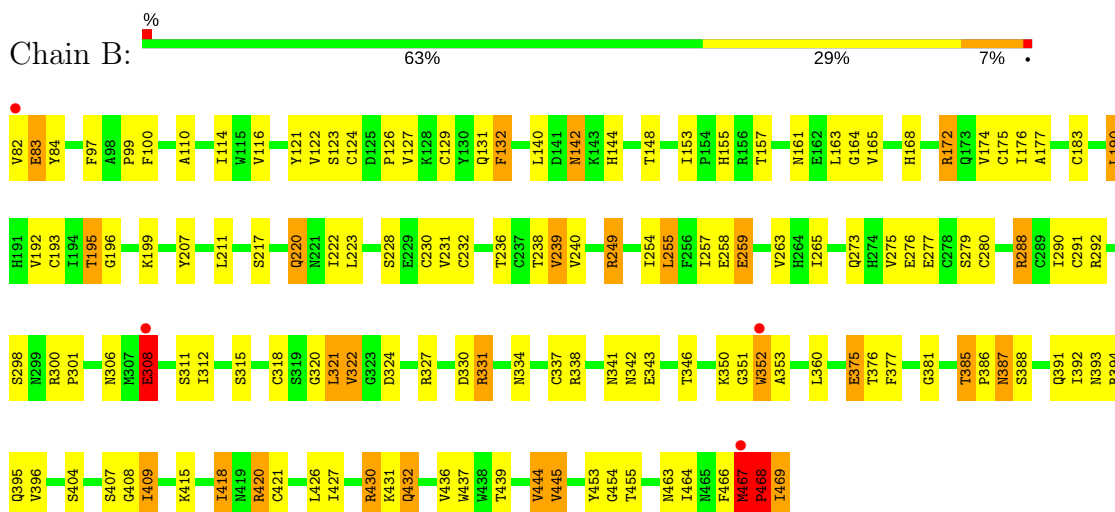
3 Residue-property plots

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: INFLUENZA A SUBTYPE N2 NEURAMINIDASE



• Molecule 1: INFLUENZA A SUBTYPE N2 NEURAMINIDASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	120.97Å 141.14Å 142.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	5.00 – 1.90 9.65 – 1.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (5.00-1.90) 34.1 (9.65-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 1.80Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.205 , (Not available) 0.239 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	16.7	Xtriage
Anisotropy	0.461	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.47 , 78.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	8557	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.06% 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: BMA, NAG, CA, FUC, FUL, MAN

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	0.55	0/3092	0.84	3/4194 (0.1%)
1	B	0.55	0/3092	0.84	3/4194 (0.1%)
All	All	0.55	0/6184	0.84	6/8388 (0.1%)

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
5	B	1	0

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	467	MET	N-CA-C	6.16	127.63	111.00
1	B	467	MET	N-CA-C	6.16	127.63	111.00
1	A	468	PRO	CA-N-CD	-5.38	103.96	111.50
1	B	468	PRO	CA-N-CD	-5.38	103.96	111.50
1	A	190	LEU	CA-CB-CG	5.15	127.15	115.30
1	B	190	LEU	CA-CB-CG	5.15	127.15	115.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	B	477	FUC	C1

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3022	723	2851	136	75
1	B	3022	723	2850	167	4
2	A	56	54	50	3	0
2	B	56	54	50	3	0
3	A	49	47	43	1	0
4	B	144	134	108	51	74
5	B	49	47	43	0	0
6	A	1	0	0	1	0
6	B	1	0	0	0	0
7	A	120	240	0	18	3
7	B	5	10	0	2	0
All	All	6525	2032	5995	306	78

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 25.

All (306) 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:453:TYR:HD1	4:B:478:NAG:C8	1.33	1.41
1:B:453:TYR:CD1	4:B:478:NAG:C8	2.04	1.38
1:B:453:TYR:CD1	4:B:478:NAG:H81	1.64	1.27
1:B:377:PHE:HB3	4:B:481:MAN:H3	1.25	1.14
1:B:394:ARG:HD3	4:B:480:BMA:H3	1.14	1.14
1:B:453:TYR:CD1	4:B:478:NAG:H82	1.81	1.11
1:A:467:MET:HB3	1:A:468:PRO:HD2	1.15	1.10
1:B:467:MET:HB3	1:B:468:PRO:HD2	1.15	1.10
1:A:172:ARG:HG2	1:B:165:VAL:HG12	1.10	1.09
1:B:352:TRP:CD1	1:B:409:ILE:HB	1.92	1.04
1:A:352:TRP:CD1	1:A:409:ILE:HB	1.92	1.03
1:B:453:TYR:HD1	4:B:478:NAG:H81	1.00	1.02
1:B:455:THR:OG1	4:B:483:MAN:H2	1.62	1.00
1:B:454:GLY:O	4:B:478:NAG:C8	2.09	1.00
1:B:177:ALA:HB2	1:B:193:CYS:HB3	1.44	0.99
1:A:177:ALA:HB2	1:A:193:CYS:HB3	1.44	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:172:ARG:CG	1:B:165:VAL:HG12	1.94	0.96
1:B:455:THR:HB	4:B:483:MAN:H3	1.48	0.95
1:B:454:GLY:O	4:B:478:NAG:C7	2.14	0.95
1:A:437:TRP:H	1:A:469:ILE:HG21	1.31	0.94
1:B:455:THR:CB	4:B:483:MAN:H3	1.96	0.94
1:B:437:TRP:H	1:B:469:ILE:HG21	1.31	0.93
1:B:394:ARG:HD3	4:B:480:BMA:C3	1.98	0.93
1:A:467:MET:HB3	1:A:468:PRO:CD	1.99	0.92
1:B:454:GLY:CA	4:B:478:NAG:H83	1.99	0.92
1:B:454:GLY:N	4:B:478:NAG:H83	1.84	0.91
1:B:467:MET:HB3	1:B:468:PRO:CD	1.99	0.91
1:B:279:SER:HB3	1:B:352:TRP:CE2	2.06	0.91
1:A:279:SER:HB3	1:A:352:TRP:CE2	2.06	0.90
1:B:352:TRP:NE1	1:B:409:ILE:HB	1.86	0.90
1:A:352:TRP:NE1	1:A:409:ILE:HB	1.86	0.89
1:B:454:GLY:O	4:B:478:NAG:H83	1.72	0.89
1:B:453:TYR:CE1	4:B:478:NAG:H81	2.08	0.89
1:B:288:ARG:HH11	1:B:288:ARG:HG3	1.38	0.87
1:A:288:ARG:HH11	1:A:288:ARG:HG3	1.38	0.87
1:B:455:THR:CB	4:B:483:MAN:C3	2.47	0.86
1:B:391:GLN:C	4:B:481:MAN:O6	2.13	0.86
1:B:352:TRP:CE3	1:B:408:GLY:HA2	2.11	0.85
1:A:352:TRP:CE3	1:A:408:GLY:HA2	2.11	0.85
1:B:436:VAL:HA	1:B:469:ILE:HG22	1.60	0.84
1:A:436:VAL:HA	1:A:469:ILE:HG22	1.60	0.84
1:B:455:THR:OG1	4:B:483:MAN:C2	2.25	0.84
1:A:169:LEU:HD13	1:B:114:ILE:HD11	1.59	0.83
1:B:455:THR:HB	4:B:483:MAN:C3	2.08	0.83
1:A:418:ILE:HD11	1:A:420:ARG:NH2	1.94	0.83
1:B:394:ARG:CD	4:B:480:BMA:H3	2.05	0.82
1:B:418:ILE:HD11	1:B:420:ARG:NH2	1.94	0.82
1:A:172:ARG:HG2	1:B:165:VAL:CG1	2.04	0.82
1:B:394:ARG:HD2	4:B:481:MAN:O4	1.79	0.82
1:B:392:ILE:N	4:B:481:MAN:O6	2.12	0.81
1:B:377:PHE:HB3	4:B:481:MAN:C3	2.11	0.78
1:B:352:TRP:CZ3	1:B:408:GLY:HA2	2.19	0.78
1:A:352:TRP:CZ3	1:A:408:GLY:HA2	2.19	0.78
1:B:394:ARG:O	4:B:482:MAN:H3	1.82	0.77
1:B:394:ARG:O	4:B:482:MAN:C3	2.31	0.76
1:B:454:GLY:C	4:B:478:NAG:H83	2.06	0.76
1:B:453:TYR:CE1	4:B:478:NAG:C8	2.66	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:279:SER:HA	1:B:352:TRP:CZ2	2.22	0.75
1:A:279:SER:HA	1:A:352:TRP:CZ2	2.22	0.75
1:A:296:LYS:HD2	7:A:589:HOH:O	1.86	0.74
1:B:376:THR:O	4:B:482:MAN:O6	2.03	0.74
1:B:377:PHE:CB	4:B:481:MAN:H3	2.12	0.73
1:B:255:LEU:HD13	1:B:265:ILE:HG12	1.72	0.71
1:A:312:ILE:HG13	7:A:527:HOH:O	1.89	0.71
1:B:376:THR:OG1	4:B:482:MAN:O6	2.04	0.71
1:A:255:LEU:HD13	1:A:265:ILE:HG12	1.72	0.71
1:A:467:MET:CB	1:A:468:PRO:HD2	2.07	0.71
1:B:467:MET:CB	1:B:468:PRO:HD2	2.07	0.70
1:A:375:GLU:HG3	1:A:394:ARG:NH1	2.07	0.69
1:B:375:GLU:HG3	1:B:394:ARG:NH1	2.07	0.69
1:A:148:THR:HG22	1:A:439:THR:HB	1.75	0.69
1:B:279:SER:HB3	1:B:352:TRP:CD2	2.27	0.69
1:A:279:SER:HB3	1:A:352:TRP:CD2	2.27	0.69
1:A:144:HIS:HE2	1:B:463:ASN:H	1.38	0.68
1:B:148:THR:HG22	1:B:439:THR:HB	1.75	0.68
1:A:409:ILE:HG12	1:A:421:CYS:O	1.96	0.66
1:B:409:ILE:HG12	1:B:421:CYS:O	1.96	0.66
1:A:321:LEU:O	1:A:322:VAL:HB	1.97	0.65
1:B:321:LEU:O	1:B:322:VAL:HB	1.97	0.64
1:B:418:ILE:HD11	1:B:420:ARG:CZ	2.29	0.63
1:B:258:GLU:HG3	1:B:263:VAL:HG11	1.80	0.63
1:A:249:ARG:HH22	1:A:273:GLN:NE2	1.97	0.62
1:A:418:ILE:HD11	1:A:420:ARG:CZ	2.29	0.62
1:A:258:GLU:HG3	1:A:263:VAL:HG11	1.80	0.62
1:A:207:TYR:CE2	1:A:259:GLU:HG3	2.35	0.62
1:B:127:VAL:O	7:B:513:HOH:O	2.16	0.62
1:B:454:GLY:HA2	4:B:478:NAG:HN2	1.64	0.61
1:B:249:ARG:HH22	1:B:273:GLN:NE2	1.97	0.61
1:B:207:TYR:CE2	1:B:259:GLU:HG3	2.35	0.61
1:B:455:THR:OG1	4:B:483:MAN:H3	1.99	0.61
1:A:279:SER:HB2	1:A:290:ILE:O	2.02	0.60
1:B:455:THR:CG2	4:B:483:MAN:O3	2.50	0.60
1:B:394:ARG:CD	4:B:481:MAN:H2	1.78	0.60
1:A:97:PHE:H	1:A:395:GLN:HE22	1.49	0.59
1:B:97:PHE:H	1:B:395:GLN:HE22	1.49	0.59
1:B:279:SER:HB2	1:B:290:ILE:O	2.02	0.59
1:B:392:ILE:N	4:B:481:MAN:HO6	2.01	0.58
1:B:426:LEU:HD11	1:B:444:VAL:HG22	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:352:TRP:NE1	1:B:409:ILE:CB	2.64	0.58
1:A:436:VAL:HA	1:A:469:ILE:CG2	2.32	0.58
1:B:288:ARG:NH1	1:B:288:ARG:HG3	2.15	0.58
6:A:470:CA:CA	7:A:592:HOH:O	1.79	0.57
1:B:436:VAL:HA	1:B:469:ILE:CG2	2.32	0.57
1:A:426:LEU:HD11	1:A:444:VAL:HG22	1.85	0.57
1:A:83:GLU:CD	1:A:84:TYR:H	2.08	0.57
1:B:83:GLU:CD	1:B:84:TYR:H	2.08	0.57
1:B:300:ARG:CZ	1:B:351:GLY:HA3	2.35	0.56
1:B:350:LYS:HE3	1:B:352:TRP:CZ3	2.40	0.56
1:A:130:TYR:HB2	7:A:506:HOH:O	2.04	0.56
1:A:317:VAL:HG23	7:A:543:HOH:O	2.04	0.56
1:B:157:THR:HG22	1:B:176:ILE:HA	1.87	0.56
1:A:249:ARG:HH22	1:A:273:GLN:CD	2.09	0.56
1:A:351:GLY:O	1:A:352:TRP:CE3	2.59	0.56
1:B:436:VAL:CA	1:B:469:ILE:HG22	2.36	0.56
1:A:300:ARG:CZ	1:A:351:GLY:HA3	2.35	0.56
1:B:454:GLY:HA2	4:B:478:NAG:N2	2.21	0.56
1:A:157:THR:HG22	1:A:176:ILE:HA	1.87	0.55
1:A:239:VAL:HG12	1:A:255:LEU:HB2	1.88	0.55
1:A:436:VAL:CA	1:A:469:ILE:HG22	2.36	0.55
1:B:351:GLY:O	1:B:352:TRP:CE3	2.59	0.55
1:A:306:ASN:HD21	1:A:308:GLU:CD	2.09	0.55
1:A:350:LYS:HE3	1:A:352:TRP:CZ3	2.40	0.55
1:B:249:ARG:HH22	1:B:273:GLN:CD	2.09	0.55
1:B:324:ASP:O	1:B:327:ARG:HD3	2.06	0.55
1:B:306:ASN:HD21	1:B:308:GLU:CD	2.09	0.55
1:A:324:ASP:O	1:A:327:ARG:HD3	2.06	0.55
1:A:352:TRP:NE1	1:A:409:ILE:CB	2.64	0.55
1:B:239:VAL:HG12	1:B:255:LEU:HB2	1.88	0.54
1:B:279:SER:CB	1:B:352:TRP:CE2	2.87	0.54
1:B:148:THR:HG21	1:B:439:THR:O	2.07	0.54
1:A:334:ASN:HB3	1:A:387:ASN:HD21	1.73	0.54
1:A:148:THR:HG21	1:A:439:THR:O	2.07	0.54
1:B:157:THR:HB	1:B:175:CYS:O	2.08	0.54
1:A:437:TRP:H	1:A:469:ILE:CG2	2.14	0.53
1:A:157:THR:HB	1:A:175:CYS:O	2.08	0.53
1:A:437:TRP:N	1:A:469:ILE:HG21	2.13	0.53
1:A:191:HIS:CE1	7:A:507:HOH:O	2.62	0.53
1:B:276:GLU:O	1:B:291:CYS:HB3	2.09	0.53
1:B:279:SER:CA	1:B:352:TRP:CZ2	2.92	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:404:SER:O	1:B:427:ILE:HG22	2.09	0.53
1:B:430:ARG:HD3	1:B:436:VAL:O	2.08	0.53
1:A:404:SER:O	1:A:427:ILE:HG22	2.09	0.53
1:B:334:ASN:HB3	1:B:387:ASN:HD21	1.73	0.53
1:A:430:ARG:HD3	1:A:436:VAL:O	2.08	0.52
1:B:453:TYR:C	4:B:478:NAG:C8	2.77	0.52
1:A:276:GLU:O	1:A:291:CYS:HB3	2.09	0.52
1:A:279:SER:CA	1:A:352:TRP:CZ2	2.92	0.52
1:A:246:ALA:HB3	7:A:599:HOH:O	2.09	0.52
1:A:110:ALA:HA	1:A:468:PRO:HG2	1.92	0.52
1:B:453:TYR:O	7:B:512:HOH:O	2.18	0.52
1:A:352:TRP:HE1	1:A:409:ILE:CG2	2.22	0.52
1:B:352:TRP:CD2	1:B:408:GLY:HA2	2.43	0.52
1:B:352:TRP:HE1	1:B:409:ILE:CG2	2.22	0.51
1:A:279:SER:CB	1:A:352:TRP:CE2	2.87	0.51
4:B:485:NAG:H5	4:B:486:BMA:O5	2.10	0.51
1:B:110:ALA:HA	1:B:468:PRO:HG2	1.92	0.51
1:A:352:TRP:CD2	1:A:408:GLY:HA2	2.43	0.51
4:B:478:NAG:O3	4:B:479:NAG:N2	2.41	0.51
1:B:454:GLY:CA	4:B:478:NAG:C8	2.83	0.50
1:A:210:ARG:NH2	1:B:126:PRO:O	2.45	0.50
4:B:479:NAG:H5	4:B:480:BMA:O5	2.10	0.50
4:B:484:NAG:O3	4:B:485:NAG:N2	2.41	0.50
1:B:437:TRP:H	1:B:469:ILE:CG2	2.14	0.50
1:A:321:LEU:O	1:A:322:VAL:CB	2.59	0.50
1:A:352:TRP:HZ2	7:A:525:HOH:O	1.94	0.50
1:B:121:TYR:CG	1:B:228:SER:HA	2.47	0.50
1:B:321:LEU:O	1:B:322:VAL:CB	2.59	0.50
1:B:350:LYS:NZ	1:B:352:TRP:CH2	2.78	0.49
2:B:472:NAG:H61	2:B:473:NAG:O7	2.12	0.49
1:A:288:ARG:NH1	1:A:288:ARG:HG3	2.15	0.49
1:B:377:PHE:CB	4:B:481:MAN:C3	2.82	0.49
2:A:472:NAG:H61	2:A:473:NAG:O7	2.12	0.49
1:A:121:TYR:CG	1:A:228:SER:HA	2.47	0.49
1:A:320:GLY:N	1:A:388:SER:O	2.46	0.49
1:A:259:GLU:HA	7:A:520:HOH:O	2.12	0.49
1:A:258:GLU:HG3	1:A:263:VAL:CG1	2.43	0.48
1:A:353:ALA:HB1	1:A:360:LEU:HD11	1.95	0.48
1:B:350:LYS:HE3	1:B:352:TRP:HZ3	1.78	0.48
1:B:453:TYR:C	4:B:478:NAG:H83	2.31	0.48
1:B:430:ARG:NH2	1:B:437:TRP:CZ3	2.82	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:394:ARG:HD3	4:B:481:MAN:H2	1.85	0.48
1:A:273:GLN:HE21	1:A:273:GLN:HA	1.79	0.48
1:A:199:LYS:HA	1:A:220:GLN:O	2.14	0.48
1:A:430:ARG:NH2	1:A:437:TRP:CZ3	2.82	0.48
1:B:142:ASN:HD22	1:B:144:HIS:H	1.62	0.48
1:B:353:ALA:HB1	1:B:360:LEU:HD11	1.95	0.48
1:A:403:ARG:HG2	7:A:604:HOH:O	2.14	0.48
1:A:426:LEU:HD11	1:A:444:VAL:CG2	2.44	0.48
1:B:273:GLN:HA	1:B:273:GLN:HE21	1.79	0.48
1:B:426:LEU:HD11	1:B:444:VAL:CG2	2.44	0.48
1:B:258:GLU:HG3	1:B:263:VAL:CG1	2.43	0.48
1:B:320:GLY:N	1:B:388:SER:O	2.46	0.48
1:B:437:TRP:N	1:B:469:ILE:HG21	2.13	0.48
1:B:392:ILE:C	4:B:481:MAN:O3	2.52	0.48
1:A:142:ASN:HD22	1:A:144:HIS:H	1.62	0.47
2:A:472:NAG:C6	2:A:473:NAG:O7	2.62	0.47
1:B:199:LYS:HA	1:B:220:GLN:O	2.14	0.47
1:A:176:ILE:HD12	1:B:99:PRO:HB3	1.94	0.47
1:B:273:GLN:HA	1:B:273:GLN:NE2	2.29	0.47
1:A:273:GLN:HA	1:A:273:GLN:NE2	2.29	0.47
1:A:82:VAL:HG12	1:A:83:GLU:N	2.29	0.47
1:B:254:ILE:N	1:B:254:ILE:HD12	2.29	0.47
3:A:477:FUL:HO3	1:B:466:PHE:HZ	1.62	0.47
1:A:142:ASN:HD21	1:B:110:ALA:HB3	1.78	0.47
1:A:254:ILE:N	1:A:254:ILE:HD12	2.29	0.47
1:A:350:LYS:NZ	1:A:352:TRP:CH2	2.78	0.47
1:B:279:SER:HA	1:B:352:TRP:CH2	2.49	0.47
2:B:472:NAG:C6	2:B:473:NAG:O7	2.62	0.47
1:A:279:SER:HA	1:A:352:TRP:CH2	2.49	0.47
1:B:82:VAL:HG12	1:B:83:GLU:N	2.29	0.47
1:B:217:SER:HA	1:B:223:LEU:HD22	1.96	0.47
1:A:217:SER:HA	1:A:223:LEU:HD22	1.96	0.47
1:A:307:MET:N	7:A:529:HOH:O	2.48	0.47
1:B:392:ILE:C	4:B:481:MAN:HO3	2.18	0.47
1:A:467:MET:CB	1:A:468:PRO:CD	2.81	0.46
1:A:321:LEU:HD11	1:A:391:GLN:HB3	1.98	0.46
1:B:131:GLN:NE2	1:B:164:GLY:H	2.13	0.46
1:B:334:ASN:CB	1:B:387:ASN:HD21	2.28	0.46
1:B:392:ILE:HG12	1:B:393:ASN:H	1.80	0.46
1:A:195:THR:HG22	1:A:196:GLY:H	1.80	0.46
1:A:311:SER:O	1:A:312:ILE:HD13	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:PRO:HA	1:A:315:SER:O	2.16	0.46
1:B:311:SER:O	1:B:312:ILE:HD13	2.16	0.46
1:B:195:THR:HG22	1:B:196:GLY:H	1.80	0.46
1:A:131:GLN:NE2	1:A:164:GLY:H	2.13	0.46
1:A:300:ARG:NH2	1:A:351:GLY:HA3	2.30	0.46
1:B:124:CYS:HA	1:B:129:CYS:HA	1.97	0.46
1:A:94:ILE:HG23	7:A:553:HOH:O	2.16	0.46
1:B:275:VAL:O	1:B:276:GLU:HG2	2.16	0.46
1:B:334:ASN:HA	1:B:387:ASN:HD21	1.81	0.46
1:A:334:ASN:CB	1:A:387:ASN:HD21	2.28	0.46
1:B:321:LEU:HD11	1:B:391:GLN:HB3	1.98	0.46
1:A:392:ILE:HG12	1:A:393:ASN:H	1.80	0.45
1:B:301:PRO:HA	1:B:315:SER:O	2.16	0.45
1:A:124:CYS:HA	1:A:129:CYS:HA	1.97	0.45
1:B:300:ARG:NH2	1:B:351:GLY:HA3	2.30	0.45
1:A:334:ASN:HA	1:A:387:ASN:HD21	1.81	0.45
1:B:455:THR:OG1	4:B:483:MAN:C3	2.59	0.45
1:A:275:VAL:O	1:A:276:GLU:HG2	2.16	0.45
1:B:131:GLN:HG2	1:B:163:LEU:HA	1.98	0.45
1:A:350:LYS:HE3	1:A:352:TRP:HZ3	1.78	0.44
1:B:337:CYS:SG	1:B:386:PRO:HB3	2.57	0.44
1:A:131:GLN:HG2	1:A:163:LEU:HA	1.98	0.44
1:A:337:CYS:SG	1:A:386:PRO:HB3	2.57	0.44
1:B:140:LEU:CD2	1:B:468:PRO:HB3	2.48	0.44
1:A:409:ILE:HD12	7:A:523:HOH:O	2.18	0.44
1:A:352:TRP:HE1	1:A:409:ILE:HG22	1.82	0.44
1:A:183:CYS:HB3	1:A:230:CYS:O	2.18	0.43
1:A:392:ILE:HG12	1:A:393:ASN:N	2.34	0.43
1:B:352:TRP:HE1	1:B:409:ILE:HG22	1.82	0.43
1:A:140:LEU:CD2	1:A:468:PRO:HB3	2.48	0.43
1:A:232:CYS:HA	1:A:236:THR:O	2.18	0.43
1:A:299:ASN:N	1:A:299:ASN:OD1	2.51	0.43
1:B:132:PHE:CD1	1:B:132:PHE:N	2.86	0.43
1:B:331:ARG:HD2	1:B:331:ARG:N	2.34	0.43
1:A:432:GLN:HE21	1:A:432:GLN:C	2.22	0.43
1:B:392:ILE:HG12	1:B:393:ASN:N	2.34	0.43
1:B:183:CYS:HB3	1:B:230:CYS:O	2.18	0.43
1:A:169:LEU:HA	1:A:169:LEU:HD23	1.84	0.43
1:A:331:ARG:HD2	1:A:331:ARG:N	2.34	0.43
1:B:116:VAL:HG13	1:B:439:THR:O	2.19	0.43
2:A:485:NAG:O7	2:A:485:NAG:H3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:232:CYS:HA	1:B:236:THR:O	2.18	0.42
1:A:116:VAL:HG13	1:A:439:THR:O	2.19	0.42
1:A:132:PHE:N	1:A:132:PHE:CD1	2.86	0.42
1:B:153:ILE:HG13	1:B:155:HIS:H	1.84	0.42
1:B:280:CYS:HA	1:B:288:ARG:O	2.20	0.42
1:A:151:ASP:N	7:A:586:HOH:O	2.53	0.42
1:A:376:THR:O	1:A:394:ARG:HA	2.20	0.42
1:A:352:TRP:CZ3	1:A:407:SER:O	2.73	0.42
1:B:381:GLY:O	1:B:385:THR:HG23	2.19	0.42
1:B:376:THR:O	1:B:394:ARG:HA	2.20	0.42
1:A:188:ALA:HB1	7:A:510:HOH:O	2.20	0.42
1:B:82:VAL:HG12	1:B:83:GLU:O	2.20	0.42
2:B:491:NAG:H3	2:B:491:NAG:O7	2.19	0.42
1:A:341:ASN:O	1:A:343:GLU:HG3	2.20	0.42
1:B:279:SER:OG	1:B:351:GLY:O	2.38	0.42
1:B:432:GLN:C	1:B:432:GLN:HE21	2.22	0.42
1:A:350:LYS:CE	1:A:352:TRP:HZ3	2.33	0.42
1:B:318:CYS:HB3	1:B:385:THR:O	2.20	0.42
1:B:437:TRP:HB2	1:B:469:ILE:HD13	2.02	0.42
1:A:153:ILE:HG13	1:A:155:HIS:H	1.84	0.41
1:A:318:CYS:HB3	1:A:385:THR:O	2.20	0.41
1:B:350:LYS:CE	1:B:352:TRP:HZ3	2.33	0.41
1:A:381:GLY:O	1:A:385:THR:HG23	2.19	0.41
1:A:437:TRP:HB2	1:A:469:ILE:HD13	2.02	0.41
1:B:350:LYS:CE	1:B:352:TRP:CZ3	3.03	0.41
1:B:453:TYR:C	4:B:478:NAG:H82	2.40	0.41
1:A:279:SER:OG	1:A:351:GLY:O	2.38	0.41
1:A:280:CYS:HA	1:A:288:ARG:O	2.20	0.41
1:B:350:LYS:NZ	1:B:352:TRP:HH2	2.17	0.41
1:B:352:TRP:CZ3	1:B:407:SER:O	2.73	0.41
1:B:308:GLU:HG3	1:B:308:GLU:H	1.67	0.41
1:A:149:VAL:HB	7:A:572:HOH:O	2.21	0.41
1:A:409:ILE:HG21	7:A:523:HOH:O	2.20	0.41
1:A:202:THR:O	1:A:202:THR:HG23	2.21	0.41
1:A:352:TRP:NE1	1:A:409:ILE:CG2	2.84	0.41
1:A:350:LYS:CE	1:A:352:TRP:CZ3	3.03	0.40
1:B:467:MET:CB	1:B:468:PRO:CD	2.81	0.40
1:A:100:PHE:HB3	1:A:445:VAL:HG22	2.02	0.40
1:A:168:HIS:H	1:A:168:HIS:CD2	2.40	0.40
1:A:82:VAL:HG12	1:A:83:GLU:O	2.20	0.40
1:B:168:HIS:CD2	1:B:168:HIS:H	2.40	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:PHE:HE2	7:A:507:HOH:O	2.01	0.40
1:B:341:ASN:O	1:B:343:GLU:HG3	2.20	0.40
1:B:100:PHE:HB3	1:B:445:VAL:HG22	2.02	0.40

All (78) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:394:ARG:CB	4:B:487:MAN:C2[4_555]	0.65	1.55
1:A:394:ARG:N	4:B:488:MAN:C1[4_555]	0.82	1.38
1:A:394:ARG:CA	4:B:487:MAN:O2[4_555]	0.91	1.29
1:A:394:ARG:CA	4:B:488:MAN:C1[4_555]	1.10	1.10
1:A:391:GLN:OE1	4:B:487:MAN:C4[4_555]	1.11	1.09
1:A:393:ASN:C	4:B:488:MAN:C2[4_555]	1.20	1.00
1:B:338:ARG:HH11	7:A:489:HOH:H2[7_544]	0.70	0.90
1:A:393:ASN:C	4:B:488:MAN:O2[4_555]	1.34	0.86
1:A:395:GLN:N	4:B:488:MAN:C5[4_555]	1.37	0.83
1:A:391:GLN:OE1	4:B:487:MAN:C5[4_555]	1.37	0.83
1:A:394:ARG:CB	4:B:487:MAN:O2[4_555]	1.38	0.82
1:A:393:ASN:C	4:B:488:MAN:C1[4_555]	1.41	0.79
1:A:393:ASN:O	4:B:488:MAN:O2[4_555]	1.41	0.79
1:A:395:GLN:CB	4:B:488:MAN:C6[4_555]	1.41	0.79
1:A:395:GLN:CG	4:B:488:MAN:O4[4_555]	1.41	0.79
1:A:394:ARG:C	4:B:488:MAN:H5[4_555]	0.83	0.77
1:A:394:ARG:N	4:B:487:MAN:O2[4_555]	1.43	0.77
1:A:394:ARG:C	4:B:488:MAN:C5[4_555]	1.47	0.73
1:A:455:THR:CG2	4:B:489:MAN:O2[4_555]	1.48	0.72
1:A:394:ARG:CB	4:B:487:MAN:H2[4_555]	0.90	0.70
1:A:391:GLN:CG	4:B:487:MAN:H61[4_555]	0.91	0.69
1:A:394:ARG:N	4:B:488:MAN:C2[4_555]	1.54	0.66
1:A:393:ASN:CA	4:B:488:MAN:O2[4_555]	1.56	0.64
1:A:391:GLN:CG	4:B:487:MAN:C6[4_555]	1.57	0.63
1:A:395:GLN:CA	4:B:488:MAN:H62[4_555]	0.98	0.62
1:A:395:GLN:N	4:B:488:MAN:H5[4_555]	1.02	0.58
1:A:394:ARG:CG	4:B:487:MAN:H2[4_555]	1.03	0.57
1:A:394:ARG:CA	4:B:488:MAN:O5[4_555]	1.65	0.55
1:A:391:GLN:OE1	4:B:487:MAN:H4[4_555]	1.06	0.54
1:A:376:THR:HG1	4:B:488:MAN:HO6[4_555]	1.08	0.52
1:A:394:ARG:C	4:B:487:MAN:O2[4_555]	1.69	0.51
1:A:395:GLN:CB	4:B:488:MAN:H62[4_555]	1.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:391:GLN:CD	4:B:487:MAN:H61[4_555]	1.10	0.50
1:A:391:GLN:OE1	4:B:487:MAN:C6[4_555]	1.70	0.50
1:A:391:GLN:CD	4:B:487:MAN:C6[4_555]	1.71	0.49
1:A:395:GLN:N	4:B:488:MAN:C6[4_555]	1.72	0.48
1:A:395:GLN:CG	4:B:488:MAN:C4[4_555]	1.74	0.46
1:A:394:ARG:CB	4:B:487:MAN:C3[4_555]	1.75	0.45
1:A:395:GLN:CA	4:B:488:MAN:C6[4_555]	1.75	0.45
1:A:394:ARG:C	4:B:488:MAN:O5[4_555]	1.75	0.45
1:A:393:ASN:CB	4:B:488:MAN:O2[4_555]	1.78	0.42
1:A:394:ARG:N	4:B:488:MAN:H1[4_555]	1.22	0.38
1:B:172:ARG:HH12	7:A:505:HOH:H2[4_555]	1.24	0.36
1:A:393:ASN:O	4:B:488:MAN:C1[4_555]	1.87	0.33
1:A:394:ARG:CA	4:B:487:MAN:C2[4_555]	1.88	0.32
1:A:376:THR:OG1	4:B:488:MAN:HO6[4_555]	1.29	0.31
1:A:391:GLN:CB	4:B:487:MAN:O6[4_555]	1.90	0.30
1:A:395:GLN:CB	4:B:488:MAN:H61[4_555]	1.31	0.29
1:A:463:ASN:H	1:B:144:HIS:HE2[4_555]	1.31	0.29
1:A:391:GLN:CG	4:B:487:MAN:O6[4_555]	1.91	0.29
1:A:394:ARG:O	4:B:488:MAN:H5[4_555]	1.32	0.28
1:A:393:ASN:O	4:B:488:MAN:C2[4_555]	1.92	0.28
1:A:393:ASN:CA	4:B:488:MAN:C2[4_555]	1.92	0.28
1:A:455:THR:OG1	4:B:489:MAN:H2[4_555]	1.37	0.23
1:A:391:GLN:OE1	4:B:487:MAN:O4[4_555]	1.97	0.23
1:A:394:ARG:C	4:B:488:MAN:C1[4_555]	1.97	0.23
1:A:393:ASN:O	4:B:488:MAN:O5[4_555]	1.98	0.22
1:A:391:GLN:C	4:B:487:MAN:HO6[4_555]	1.39	0.21
1:A:393:ASN:C	4:B:488:MAN:H1[4_555]	1.40	0.20
1:A:391:GLN:CB	4:B:487:MAN:HO6[4_555]	1.40	0.20
1:A:394:ARG:CG	4:B:487:MAN:C2[4_555]	2.00	0.20
1:A:395:GLN:N	4:B:488:MAN:O5[4_555]	2.02	0.18
1:A:395:GLN:CG	4:B:488:MAN:C6[4_555]	2.03	0.17
1:A:376:THR:O	4:B:488:MAN:O6[4_555]	2.03	0.17
1:A:453:TYR:CD1	4:B:484:NAG:C8[4_555]	2.04	0.16
1:A:376:THR:OG1	4:B:488:MAN:O6[4_555]	2.04	0.16
1:A:394:ARG:O	4:B:488:MAN:C5[4_555]	2.06	0.14
1:A:395:GLN:CG	4:B:488:MAN:H61[4_555]	1.47	0.13
1:A:454:GLY:O	4:B:484:NAG:C8[4_555]	2.09	0.11
1:A:395:GLN:CA	4:B:488:MAN:C5[4_555]	2.10	0.10
1:A:392:ILE:N	4:B:487:MAN:O6[4_555]	2.12	0.08
1:A:391:GLN:CA	4:B:487:MAN:HO6[4_555]	1.53	0.07
1:A:394:ARG:O	4:B:487:MAN:O2[4_555]	2.13	0.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:391:GLN:C	4:B:487:MAN:O6[4_555]	2.13	0.07
1:A:454:GLY:O	4:B:484:NAG:C7[4_555]	2.14	0.06
1:B:338:ARG:NH1	7:A:489:HOH:H2[7_544]	1.58	0.02
1:A:391:GLN:CD	4:B:487:MAN:C5[4_555]	2.19	0.01
1:A:394:ARG:CA	4:B:488:MAN:H1[4_555]	1.60	0.00

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	386/388 (100%)	344 (89%)	33 (8%)	9 (2%)	7	1
1	B	386/388 (100%)	344 (89%)	33 (8%)	9 (2%)	7	1
All	All	772/776 (100%)	688 (89%)	66 (8%)	18 (2%)	7	1

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	322	VAL
1	A	467	MET
1	A	468	PRO
1	B	322	VAL
1	B	467	MET
1	B	468	PRO
1	A	321	LEU
1	A	330	ASP
1	B	321	LEU
1	B	330	ASP
1	A	430	ARG
1	B	430	ARG
1	A	222	ILE
1	A	277	GLU
1	A	308	GLU

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Mol	Chain	Res	Type
1	B	222	ILE
1	B	277	GLU
1	B	308	GLU

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	338/338 (100%)	294 (87%)	44 (13%)	5	1
1	B	338/338 (100%)	294 (87%)	44 (13%)	5	1
All	All	676/676 (100%)	588 (87%)	88 (13%)	5	1

All (88) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	83	GLU
1	A	122	VAL
1	A	123	SER
1	A	132	PHE
1	A	142	ASN
1	A	161	ASN
1	A	172	ARG
1	A	174	VAL
1	A	190	LEU
1	A	192	VAL
1	A	195	THR
1	A	211	LEU
1	A	220	GLN
1	A	231	VAL
1	A	238	THR
1	A	239	VAL
1	A	240	VAL
1	A	249	ARG
1	A	255	LEU
1	A	257	ILE

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Mol	Chain	Res	Type
1	A	259	GLU
1	A	288	ARG
1	A	292	ARG
1	A	298	SER
1	A	308	GLU
1	A	331	ARG
1	A	342	ASN
1	A	346	THR
1	A	352	TRP
1	A	375	GLU
1	A	385	THR
1	A	387	ASN
1	A	396	VAL
1	A	409	ILE
1	A	415	LYS
1	A	418	ILE
1	A	420	ARG
1	A	431	LYS
1	A	432	GLN
1	A	444	VAL
1	A	445	VAL
1	A	464	ILE
1	A	467	MET
1	A	469	ILE
1	B	83	GLU
1	B	122	VAL
1	B	123	SER
1	B	132	PHE
1	B	142	ASN
1	B	161	ASN
1	B	172	ARG
1	B	174	VAL
1	B	190	LEU
1	B	192	VAL
1	B	195	THR
1	B	211	LEU
1	B	220	GLN
1	B	231	VAL
1	B	238	THR
1	B	239	VAL
1	B	240	VAL
1	B	249	ARG

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Mol	Chain	Res	Type
1	B	255	LEU
1	B	257	ILE
1	B	259	GLU
1	B	288	ARG
1	B	292	ARG
1	B	298	SER
1	B	308	GLU
1	B	331	ARG
1	B	342	ASN
1	B	346	THR
1	B	352	TRP
1	B	375	GLU
1	B	385	THR
1	B	387	ASN
1	B	396	VAL
1	B	409	ILE
1	B	415	LYS
1	B	418	ILE
1	B	420	ARG
1	B	431	LYS
1	B	432	GLN
1	B	444	VAL
1	B	445	VAL
1	B	464	ILE
1	B	467	MET
1	B	469	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	104	ASN
1	A	131	GLN
1	A	142	ASN
1	A	173	GLN
1	A	273	GLN
1	A	387	ASN
1	A	393	ASN
1	A	395	GLN
1	A	402	ASN
1	A	419	ASN
1	A	432	GLN
1	A	465	ASN

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Mol	Chain	Res	Type
1	B	104	ASN
1	B	131	GLN
1	B	142	ASN
1	B	273	GLN
1	B	387	ASN
1	B	393	ASN
1	B	395	GLN
1	B	402	ASN
1	B	419	ASN
1	B	432	GLN
1	B	465	ASN

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 ⓘ

28 carbohydrates 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	NAG	A	472	1,2	14,14,15	1.43	3 (21%)	15,19,21	2.59	7 (46%)
2	NAG	A	473	2	14,14,15	1.60	3 (21%)	15,19,21	2.90	6 (40%)
3	NAG	A	474	1,3	14,14,15	1.24	1 (7%)	15,19,21	3.47	3 (20%)
3	NAG	A	475	3	14,14,15	1.46	2 (14%)	15,19,21	3.09	6 (40%)
3	BMA	A	476	3	11,11,12	1.32	2 (18%)	13,15,17	1.70	4 (30%)
3	FUL	A	477	3	9,10,11	1.34	2 (22%)	13,14,16	1.25	2 (15%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	A	484	1,2	14,14,15	1.39	2 (14%)	15,19,21	2.11	4 (26%)
2	NAG	A	485	2	14,14,15	1.09	2 (14%)	15,19,21	1.98	3 (20%)
2	NAG	B	472	1,2	14,14,15	1.43	3 (21%)	15,19,21	2.59	7 (46%)
2	NAG	B	473	2	14,14,15	1.60	3 (21%)	15,19,21	2.90	6 (40%)
5	NAG	B	474	1,5	14,14,15	1.24	1 (7%)	15,19,21	3.47	3 (20%)
5	NAG	B	475	5	14,14,15	1.46	2 (14%)	15,19,21	3.09	6 (40%)
5	BMA	B	476	5	11,11,12	1.32	2 (18%)	13,15,17	1.70	4 (30%)
5	FUC	B	477	5	9,10,11	1.34	2 (22%)	13,14,16	1.25	2 (15%)
4	NAG	B	478	1,4	14,14,15	1.62	3 (21%)	15,19,21	3.67	5 (33%)
4	NAG	B	479	4	14,14,15	2.34	6 (42%)	15,19,21	3.86	8 (53%)
4	BMA	B	480	4	11,11,12	3.81	8 (72%)	13,15,17	2.10	4 (30%)
4	MAN	B	481	4	11,11,12	2.14	5 (45%)	13,15,17	2.82	5 (38%)
4	MAN	B	482	4	11,11,12	1.52	1 (9%)	13,15,17	1.37	2 (15%)
4	MAN	B	483	4	11,11,12	1.74	2 (18%)	13,15,17	3.55	8 (61%)
4	NAG	B	484	1,4	14,14,15	1.62	3 (21%)	15,19,21	3.67	5 (33%)
4	NAG	B	485	4	14,14,15	2.34	6 (42%)	15,19,21	3.86	8 (53%)
4	BMA	B	486	4	11,11,12	3.81	8 (72%)	13,15,17	2.10	4 (30%)
4	MAN	B	487	4	11,11,12	2.14	5 (45%)	13,15,17	2.82	5 (38%)
4	MAN	B	488	4	11,11,12	1.52	1 (9%)	13,15,17	1.37	2 (15%)
4	MAN	B	489	4	11,11,12	1.74	2 (18%)	13,15,17	3.55	8 (61%)
2	NAG	B	490	1,2	14,14,15	1.39	2 (14%)	15,19,21	2.11	4 (26%)
2	NAG	B	491	2	14,14,15	1.09	2 (14%)	15,19,21	1.98	3 (20%)

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	472	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	473	2	-	0/6/23/26	0/1/1/1
3	NAG	A	474	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	475	3	-	0/6/23/26	0/1/1/1
3	BMA	A	476	3	-	0/2/19/22	0/1/1/1
3	FUL	A	477	3	-	0/0/17/20	0/1/1/1
2	NAG	A	484	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	485	2	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	472	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	473	2	-	0/6/23/26	0/1/1/1
5	NAG	B	474	1,5	-	0/6/23/26	0/1/1/1
5	NAG	B	475	5	-	0/6/23/26	0/1/1/1
5	BMA	B	476	5	-	0/2/19/22	0/1/1/1
5	FUC	B	477	5	1/1/5/5	0/0/17/20	0/1/1/1
4	NAG	B	478	1,4	-	0/6/23/26	0/1/1/1
4	NAG	B	479	4	-	0/6/23/26	0/1/1/1
4	BMA	B	480	4	-	0/2/19/22	1/1/1/1
4	MAN	B	481	4	-	0/2/19/22	0/1/1/1
4	MAN	B	482	4	-	0/2/19/22	0/1/1/1
4	MAN	B	483	4	-	0/2/19/22	0/1/1/1
4	NAG	B	484	1,4	-	0/6/23/26	0/1/1/1
4	NAG	B	485	4	-	0/6/23/26	0/1/1/1
4	BMA	B	486	4	-	0/2/19/22	1/1/1/1
4	MAN	B	487	4	-	0/2/19/22	0/1/1/1
4	MAN	B	488	4	-	0/2/19/22	0/1/1/1
4	MAN	B	489	4	-	0/2/19/22	0/1/1/1
2	NAG	B	490	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	491	2	-	0/6/23/26	0/1/1/1

All (84) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	472	NAG	C1-C2	-3.80	1.47	1.52
2	B	472	NAG	C1-C2	-3.80	1.47	1.52
4	B	484	NAG	O3-C3	-3.78	1.34	1.43
4	B	478	NAG	O3-C3	-3.78	1.34	1.43
4	B	484	NAG	O5-C1	-2.62	1.39	1.43
4	B	478	NAG	O5-C1	-2.62	1.39	1.43
4	B	484	NAG	C3-C2	-2.40	1.47	1.52
4	B	478	NAG	C3-C2	-2.40	1.47	1.52
2	A	485	NAG	C4-C5	-2.18	1.48	1.53
2	B	491	NAG	C4-C5	-2.18	1.48	1.53
3	A	477	FUL	C4-C3	2.01	1.57	1.52
5	B	477	FUC	C4-C3	2.01	1.57	1.52
2	B	490	NAG	C1-C2	2.01	1.55	1.52
2	A	484	NAG	C1-C2	2.01	1.55	1.52
5	B	476	BMA	C4-C3	2.01	1.57	1.52
3	A	476	BMA	C4-C3	2.01	1.57	1.52
2	A	485	NAG	C3-C2	2.03	1.56	1.52
2	B	491	NAG	C3-C2	2.03	1.56	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	486	BMA	O5-C5	2.03	1.47	1.43
4	B	480	BMA	O5-C5	2.03	1.47	1.43
2	A	472	NAG	C4-C5	2.05	1.57	1.53
2	B	472	NAG	C4-C5	2.05	1.57	1.53
4	B	479	NAG	C3-C2	2.15	1.57	1.52
4	B	485	NAG	C3-C2	2.15	1.57	1.52
4	B	481	MAN	C2-C3	2.18	1.55	1.52
4	B	487	MAN	C2-C3	2.18	1.55	1.52
2	B	473	NAG	C1-C2	2.19	1.55	1.52
2	A	473	NAG	C1-C2	2.19	1.55	1.52
4	B	479	NAG	O4-C4	2.24	1.48	1.43
4	B	485	NAG	O4-C4	2.24	1.48	1.43
3	A	477	FUL	O5-C1	2.32	1.47	1.43
5	B	477	FUC	O5-C1	2.32	1.47	1.43
2	A	472	NAG	O4-C4	2.36	1.48	1.43
2	B	472	NAG	O4-C4	2.36	1.48	1.43
4	B	486	BMA	C6-C5	2.45	1.60	1.51
4	B	480	BMA	C6-C5	2.45	1.60	1.51
5	B	475	NAG	O5-C5	2.47	1.48	1.43
3	A	475	NAG	O5-C5	2.47	1.48	1.43
4	B	481	MAN	O5-C1	2.62	1.48	1.43
4	B	487	MAN	O5-C1	2.62	1.48	1.43
5	B	475	NAG	C4-C5	2.64	1.58	1.53
3	A	475	NAG	C4-C5	2.64	1.58	1.53
4	B	486	BMA	O5-C1	2.68	1.48	1.43
4	B	480	BMA	O5-C1	2.68	1.48	1.43
2	B	473	NAG	C3-C2	2.72	1.58	1.52
2	A	473	NAG	C3-C2	2.72	1.58	1.52
2	B	490	NAG	O4-C4	2.72	1.49	1.43
2	A	484	NAG	O4-C4	2.72	1.49	1.43
4	B	483	MAN	C4-C5	2.75	1.58	1.53
4	B	489	MAN	C4-C5	2.75	1.58	1.53
4	B	481	MAN	C1-C2	2.77	1.58	1.52
4	B	487	MAN	C1-C2	2.77	1.58	1.52
4	B	479	NAG	C4-C5	2.89	1.59	1.53
4	B	485	NAG	C4-C5	2.89	1.59	1.53
5	B	476	BMA	C2-C3	3.17	1.56	1.52
3	A	476	BMA	C2-C3	3.17	1.56	1.52
4	B	481	MAN	O2-C2	3.36	1.50	1.43
4	B	487	MAN	O2-C2	3.36	1.50	1.43
5	B	474	NAG	O5-C5	3.37	1.50	1.43
3	A	474	NAG	O5-C5	3.37	1.50	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	486	BMA	C4-C3	3.38	1.61	1.52
4	B	480	BMA	C4-C3	3.38	1.61	1.52
2	B	473	NAG	C4-C3	3.46	1.61	1.52
2	A	473	NAG	C4-C3	3.46	1.61	1.52
4	B	479	NAG	C6-C5	3.48	1.63	1.51
4	B	485	NAG	C6-C5	3.48	1.63	1.51
4	B	479	NAG	C1-C2	3.60	1.57	1.52
4	B	485	NAG	C1-C2	3.60	1.57	1.52
4	B	481	MAN	O5-C5	3.89	1.51	1.43
4	B	487	MAN	O5-C5	3.89	1.51	1.43
4	B	482	MAN	C2-C3	4.29	1.58	1.52
4	B	488	MAN	C2-C3	4.29	1.58	1.52
4	B	486	BMA	C1-C2	4.34	1.62	1.52
4	B	480	BMA	C1-C2	4.34	1.62	1.52
4	B	486	BMA	O3-C3	4.41	1.53	1.43
4	B	480	BMA	O3-C3	4.41	1.53	1.43
4	B	483	MAN	C1-C2	4.49	1.62	1.52
4	B	489	MAN	C1-C2	4.49	1.62	1.52
4	B	479	NAG	C4-C3	4.88	1.64	1.52
4	B	485	NAG	C4-C3	4.88	1.64	1.52
4	B	486	BMA	C4-C5	6.25	1.66	1.53
4	B	480	BMA	C4-C5	6.25	1.66	1.53
4	B	486	BMA	C2-C3	7.12	1.62	1.52
4	B	480	BMA	C2-C3	7.12	1.62	1.52

All (134) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	485	NAG	C1-C2-N2	-5.91	100.39	110.49
2	B	491	NAG	C1-C2-N2	-5.91	100.39	110.49
5	B	475	NAG	C4-C3-C2	-5.91	102.36	111.02
3	A	475	NAG	C4-C3-C2	-5.91	102.36	111.02
4	B	483	MAN	O5-C1-C2	-5.66	101.91	110.79
4	B	489	MAN	O5-C1-C2	-5.66	101.91	110.79
5	B	475	NAG	O5-C1-C2	-5.38	103.98	111.47
3	A	475	NAG	O5-C1-C2	-5.38	103.98	111.47
4	B	481	MAN	C1-C2-C3	-5.24	103.01	109.65
4	B	487	MAN	C1-C2-C3	-5.24	103.01	109.65
4	B	486	BMA	C2-C3-C4	-5.10	101.98	110.88
4	B	480	BMA	C2-C3-C4	-5.10	101.98	110.88
2	B	490	NAG	C1-C2-N2	-5.01	101.94	110.49
2	A	484	NAG	C1-C2-N2	-5.01	101.94	110.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	473	NAG	O7-C7-C8	-5.00	112.96	122.06
2	A	473	NAG	O7-C7-C8	-5.00	112.96	122.06
4	B	483	MAN	O4-C4-C3	-4.97	99.54	110.36
4	B	489	MAN	O4-C4-C3	-4.97	99.54	110.36
2	A	472	NAG	C3-C4-C5	-4.25	102.73	110.22
2	B	472	NAG	C3-C4-C5	-4.25	102.73	110.22
4	B	479	NAG	O5-C1-C2	-4.21	105.62	111.47
4	B	485	NAG	O5-C1-C2	-4.21	105.62	111.47
4	B	484	NAG	C3-C4-C5	-4.11	102.98	110.22
4	B	478	NAG	C3-C4-C5	-4.11	102.98	110.22
4	B	483	MAN	O2-C2-C3	-3.90	102.51	110.17
4	B	489	MAN	O2-C2-C3	-3.90	102.51	110.17
4	B	479	NAG	O7-C7-C8	-3.63	115.45	122.06
4	B	485	NAG	O7-C7-C8	-3.63	115.45	122.06
2	A	472	NAG	O5-C1-C2	-3.44	106.69	111.47
2	B	472	NAG	O5-C1-C2	-3.44	106.69	111.47
2	B	490	NAG	O5-C1-C2	-3.33	106.83	111.47
2	A	484	NAG	O5-C1-C2	-3.33	106.83	111.47
4	B	483	MAN	O3-C3-C4	-2.93	103.98	110.36
4	B	489	MAN	O3-C3-C4	-2.93	103.98	110.36
5	B	474	NAG	C6-C5-C4	-2.80	106.44	113.00
3	A	474	NAG	C6-C5-C4	-2.80	106.44	113.00
4	B	481	MAN	C6-C5-C4	-2.74	106.59	113.00
4	B	487	MAN	C6-C5-C4	-2.74	106.59	113.00
5	B	476	BMA	C3-C4-C5	-2.62	105.59	110.22
3	A	476	BMA	C3-C4-C5	-2.62	105.59	110.22
2	A	472	NAG	O7-C7-C8	-2.39	117.71	122.06
2	B	472	NAG	O7-C7-C8	-2.39	117.71	122.06
3	A	477	FUL	O3-C3-C2	-2.09	106.23	110.02
5	B	477	FUC	O3-C3-C2	-2.09	106.23	110.02
2	A	472	NAG	C2-N2-C7	-2.07	119.92	122.94
2	B	472	NAG	C2-N2-C7	-2.07	119.92	122.94
2	A	485	NAG	C3-C4-C5	2.01	113.75	110.22
2	B	491	NAG	C3-C4-C5	2.01	113.75	110.22
2	A	472	NAG	C1-O5-C5	2.11	115.08	112.17
2	B	472	NAG	C1-O5-C5	2.11	115.08	112.17
4	B	486	BMA	C3-C4-C5	2.15	114.00	110.22
4	B	480	BMA	C3-C4-C5	2.15	114.00	110.22
4	B	479	NAG	O6-C6-C5	2.27	118.99	111.34
4	B	485	NAG	O6-C6-C5	2.27	118.99	111.34
3	A	477	FUL	C3-C4-C5	2.27	113.25	109.68
5	B	477	FUC	C3-C4-C5	2.27	113.25	109.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	490	NAG	C1-O5-C5	2.31	115.35	112.17
2	A	484	NAG	C1-O5-C5	2.31	115.35	112.17
4	B	486	BMA	O3-C3-C4	2.33	115.42	110.36
4	B	480	BMA	O3-C3-C4	2.33	115.42	110.36
4	B	483	MAN	C6-C5-C4	2.35	118.49	113.00
4	B	489	MAN	C6-C5-C4	2.35	118.49	113.00
4	B	483	MAN	C2-C3-C4	2.36	114.99	110.88
4	B	489	MAN	C2-C3-C4	2.36	114.99	110.88
5	B	476	BMA	O5-C1-C2	2.36	114.49	110.79
3	A	476	BMA	O5-C1-C2	2.36	114.49	110.79
5	B	475	NAG	C1-O5-C5	2.38	115.45	112.17
3	A	475	NAG	C1-O5-C5	2.38	115.45	112.17
5	B	476	BMA	C1-C2-C3	2.44	112.74	109.65
3	A	476	BMA	C1-C2-C3	2.44	112.74	109.65
4	B	482	MAN	C1-O5-C5	2.55	115.69	112.17
4	B	488	MAN	C1-O5-C5	2.55	115.69	112.17
2	B	473	NAG	C2-N2-C7	2.58	126.71	122.94
2	A	473	NAG	C2-N2-C7	2.58	126.71	122.94
4	B	481	MAN	O2-C2-C3	2.60	115.29	110.17
4	B	487	MAN	O2-C2-C3	2.60	115.29	110.17
2	B	473	NAG	C4-C3-C2	2.62	114.86	111.02
2	A	473	NAG	C4-C3-C2	2.62	114.86	111.02
2	A	485	NAG	C6-C5-C4	2.72	119.36	113.00
2	B	491	NAG	C6-C5-C4	2.72	119.36	113.00
4	B	482	MAN	C1-C2-C3	2.78	113.18	109.65
4	B	488	MAN	C1-C2-C3	2.78	113.18	109.65
4	B	484	NAG	O4-C4-C5	2.86	116.49	109.28
4	B	478	NAG	O4-C4-C5	2.86	116.49	109.28
5	B	475	NAG	O4-C4-C3	2.88	116.61	110.36
3	A	475	NAG	O4-C4-C3	2.88	116.61	110.36
2	A	472	NAG	O4-C4-C5	2.89	116.57	109.28
2	B	472	NAG	O4-C4-C5	2.89	116.57	109.28
4	B	479	NAG	C8-C7-N2	2.94	121.42	116.11
4	B	485	NAG	C8-C7-N2	2.94	121.42	116.11
5	B	475	NAG	O3-C3-C4	2.95	116.77	110.36
3	A	475	NAG	O3-C3-C4	2.95	116.77	110.36
2	B	490	NAG	C6-C5-C4	3.29	120.69	113.00
2	A	484	NAG	C6-C5-C4	3.29	120.69	113.00
4	B	483	MAN	C1-O5-C5	3.44	116.90	112.17
4	B	489	MAN	C1-O5-C5	3.44	116.90	112.17
5	B	476	BMA	C1-O5-C5	3.45	116.92	112.17
3	A	476	BMA	C1-O5-C5	3.45	116.92	112.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	484	NAG	C1-O5-C5	3.54	117.05	112.17
4	B	478	NAG	C1-O5-C5	3.54	117.05	112.17
4	B	479	NAG	C6-C5-C4	3.75	121.77	113.00
4	B	485	NAG	C6-C5-C4	3.75	121.77	113.00
2	B	473	NAG	C1-C2-N2	4.19	117.65	110.49
2	A	473	NAG	C1-C2-N2	4.19	117.65	110.49
4	B	486	BMA	C1-O5-C5	4.35	118.16	112.17
4	B	480	BMA	C1-O5-C5	4.35	118.16	112.17
2	B	473	NAG	C8-C7-N2	4.55	124.33	116.11
2	A	473	NAG	C8-C7-N2	4.55	124.33	116.11
4	B	481	MAN	O5-C1-C2	4.81	118.32	110.79
4	B	487	MAN	O5-C1-C2	4.81	118.32	110.79
4	B	481	MAN	C1-O5-C5	5.61	119.90	112.17
4	B	487	MAN	C1-O5-C5	5.61	119.90	112.17
4	B	484	NAG	C2-N2-C7	6.21	132.00	122.94
4	B	478	NAG	C2-N2-C7	6.21	132.00	122.94
2	B	473	NAG	C1-O5-C5	6.22	120.74	112.17
2	A	473	NAG	C1-O5-C5	6.22	120.74	112.17
2	A	472	NAG	C6-C5-C4	6.40	127.98	113.00
2	B	472	NAG	C6-C5-C4	6.40	127.98	113.00
4	B	479	NAG	C1-O5-C5	6.52	121.15	112.17
4	B	485	NAG	C1-O5-C5	6.52	121.15	112.17
5	B	475	NAG	C1-C2-N2	6.70	121.93	110.49
3	A	475	NAG	C1-C2-N2	6.70	121.93	110.49
4	B	483	MAN	O2-C2-C1	7.34	124.10	109.18
4	B	489	MAN	O2-C2-C1	7.34	124.10	109.18
4	B	479	NAG	C2-N2-C7	7.57	133.98	122.94
4	B	485	NAG	C2-N2-C7	7.57	133.98	122.94
4	B	479	NAG	C1-C2-N2	7.90	123.98	110.49
4	B	485	NAG	C1-C2-N2	7.90	123.98	110.49
5	B	474	NAG	C1-O5-C5	8.12	123.36	112.17
3	A	474	NAG	C1-O5-C5	8.12	123.36	112.17
5	B	474	NAG	C1-C2-N2	10.02	127.61	110.49
3	A	474	NAG	C1-C2-N2	10.02	127.61	110.49
4	B	484	NAG	C6-C5-C4	10.62	137.86	113.00
4	B	478	NAG	C6-C5-C4	10.62	137.86	113.00

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	B	477	FUC	C1

There are no torsion outliers.

All (2) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	486	BMA	C1-C2-C3-C4-C5-O5
4	B	480	BMA	C1-C2-C3-C4-C5-O5

19 monomers are involved in 132 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	472	NAG	2	0
2	A	473	NAG	2	0
3	A	477	FUL	1	0
2	A	485	NAG	1	0
2	B	472	NAG	2	0
2	B	473	NAG	2	0
4	B	478	NAG	20	0
4	B	479	NAG	2	0
4	B	480	BMA	4	0
4	B	481	MAN	12	0
4	B	482	MAN	4	0
4	B	483	MAN	9	0
4	B	484	NAG	1	3
4	B	485	NAG	2	0
4	B	486	BMA	1	0
4	B	487	MAN	0	28
4	B	488	MAN	0	41
4	B	489	MAN	0	2
2	B	491	NAG	1	0

5.6 Ligand geometry

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	388/388 (100%)	-0.10	5 (1%) 77 80	8, 21, 34, 44	0
1	B	388/388 (100%)	-0.10	4 (1%) 82 84	8, 21, 34, 44	0
All	All	776/776 (100%)	-0.10	9 (1%) 79 82	8, 21, 34, 44	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	352	TRP	5.7
1	B	467	MET	3.5
1	A	352	TRP	3.2
1	A	392	ILE	2.5
1	A	333	SER	2.4
1	A	308	GLU	2.3
1	B	308	GLU	2.2
1	B	82	VAL	2.2
1	A	467	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](#)

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, 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	LLDF	B-factors(\AA^2)	Q<0.9
4	MAN	B	483	11/12	0.76	0.36	12.16	15,15,48,50	0
4	MAN	B	489	11/12	0.74	0.36	10.62	15,15,48,50	0
4	MAN	B	482	11/12	0.91	0.41	8.59	15,15,57,61	0
4	NAG	B	478	14/15	0.75	0.21	7.55	15,27,40,45	0
4	MAN	B	487	11/12	0.91	0.36	6.81	15,15,51,53	0
4	MAN	B	481	11/12	0.92	0.31	5.51	15,15,51,53	0
4	MAN	B	488	11/12	0.93	0.28	5.15	15,15,57,61	0
4	NAG	B	484	14/15	0.78	0.15	1.86	15,27,40,45	0
3	NAG	A	474	14/15	0.86	0.14	1.40	15,29,42,47	0
2	NAG	A	484	14/15	0.81	0.14	0.08	15,15,46,47	0
2	NAG	B	490	14/15	0.78	0.13	-0.03	15,15,46,47	0
5	NAG	B	474	14/15	0.86	0.13	-0.35	15,29,42,47	0
3	FUL	A	477	10/11	0.70	0.32	-	15,15,49,52	0
5	BMA	B	476	11/12	0.65	0.36	-	15,15,55,57	0
4	BMA	B	480	11/12	0.54	0.51	-	15,15,50,51	0
4	NAG	B	479	14/15	0.49	0.26	-	15,15,49,49	0
2	NAG	A	485	14/15	0.63	0.24	-	15,15,50,52	0
4	BMA	B	486	11/12	0.50	0.56	-	15,15,50,51	0
2	NAG	B	472	14/15	0.82	0.16	-	15,15,36,37	0
5	FUC	B	477	10/11	0.70	0.37	-	15,15,49,52	0
4	NAG	B	485	14/15	0.42	0.28	-	15,15,49,49	0
2	NAG	A	472	14/15	0.84	0.14	-	15,15,36,37	0
2	NAG	A	473	14/15	0.67	0.26	-	15,15,52,53	0
2	NAG	B	473	14/15	0.57	0.35	-	15,15,52,53	0
2	NAG	B	491	14/15	0.82	0.20	-	15,15,50,52	0
5	NAG	B	475	14/15	0.60	0.29	-	15,15,49,50	0
3	NAG	A	475	14/15	0.70	0.24	-	15,15,49,50	0
3	BMA	A	476	11/12	0.67	0.32	-	15,15,55,57	0

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. 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, 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	LLDF	B-factors(\AA^2)	Q<0.9
6	CA	A	470	1/1	0.99	0.06	-1.56	29,29,29,29	0
6	CA	B	470	1/1	0.95	0.05	-1.94	29,29,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.