



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:12 PM GMT

PDB ID : 1J5Q
Title : The Structure and Evolution of the Major Capsid Protein of a Large, Lipid-containing, DNA virus.
Authors : Nandhagopal, N.; Simpson, A.A.; Gurnon, J.R.; Yan, X.; Baker, T.S.; Graves, M.V.; Van Etten, J.L.; Rossmann, M.G.
Deposited on : 2002-06-27
Resolution : 2.55 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

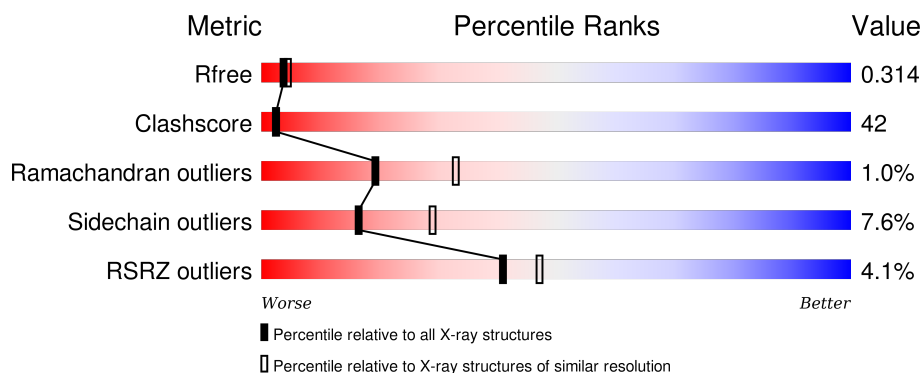
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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}	91344	4549 (2.58-2.50)
Clashscore	102246	5292 (2.58-2.50)
Ramachandran outliers	100387	5194 (2.58-2.50)
Sidechain outliers	100360	5196 (2.58-2.50)
RSRZ outliers	91569	4561 (2.58-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	437	<div> <div>4%</div> <div>48%</div> <div>41%</div> <div>6%</div> <div>5%</div> </div>
1	B	437	<div> <div>3%</div> <div>52%</div> <div>37%</div> <div>5%</div> <div>5%</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
2	NAG	A	438	-	-	X	X
2	NAG	A	444	-	-	X	-
2	NAG	B	538	-	-	X	X
2	NAG	B	549	-	-	X	X
3	MAN	A	439	X	-	-	-
3	MAN	A	440	-	-	X	-
3	MAN	A	441	-	-	X	-
3	MAN	A	442	X	-	-	-
3	MAN	A	443	-	-	X	-
3	MAN	A	445	-	-	X	-
3	MAN	A	447	X	-	X	-
3	MAN	A	448	-	-	X	-
3	MAN	A	451	X	-	-	-
3	MAN	A	452	-	-	X	-
3	MAN	A	454	X	-	-	-
3	MAN	B	540	-	-	X	-
3	MAN	B	541	-	-	X	-
3	MAN	B	542	-	-	X	-
3	MAN	B	543	X	-	X	-
3	MAN	B	547	X	-	X	-
3	MAN	B	548	X	-	X	-
3	MAN	B	551	X	-	-	-
3	MAN	B	553	-	-	X	-
3	MAN	B	554	X	-	-	-
4	NDG	A	446	-	-	X	X
4	NDG	A	449	-	-	X	X
4	NDG	B	544	-	-	X	-
4	NDG	B	550	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6956 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 Major capsid protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	413	Total	C	N	O	S	0	0	0
			3232	2052	549	623	8			
1	B	413	Total	C	N	O	S	0	0	0
			3232	2052	549	623	8			

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



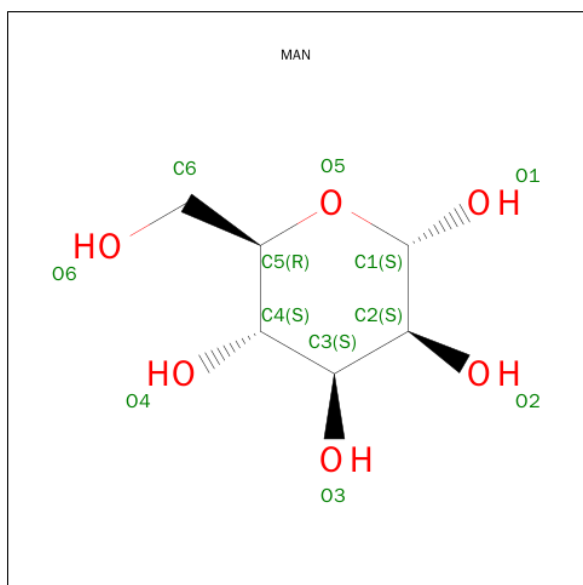
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			15	8	1	6		
2	A	1	Total	C	N	O	0	0
			15	8	1	6		
2	A	1	Total	C	N	O	0	0
			15	8	1	6		
2	B	1	Total	C	N	O	0	0
			15	8	1	6		

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

- Molecule 3 is SUGAR (ALPHA-D-MANNOSE) (three-letter code: MAN) (formula: C₆H₁₂O₆).



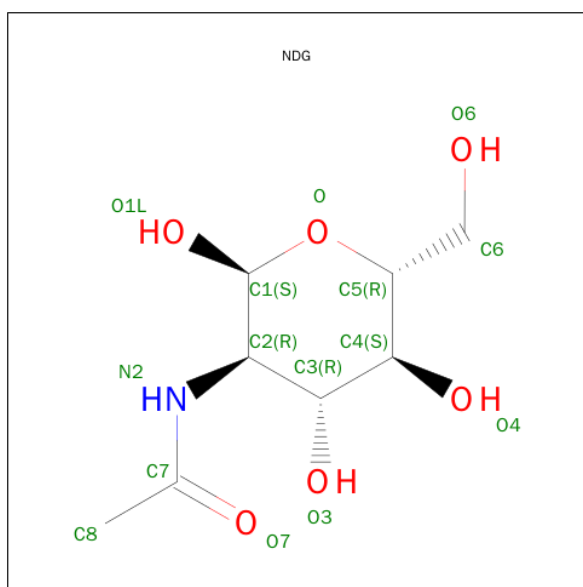
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	6	6		
3	A	1	Total	C	O	0	0
			12	6	6		
3	A	1	Total	C	O	0	0
			12	6	6		
3	A	1	Total	C	O	0	0
			12	6	6		
3	A	1	Total	C	O	0	0
			12	6	6		
3	A	1	Total	C	O	0	0
			12	6	6		
3	A	1	Total	C	O	0	0
			12	6	6		
3	A	1	Total	C	O	0	0
			12	6	6		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	6	6		
3	A	1	Total	C	O	0	0
			12	6	6		
3	A	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		

- Molecule 4 is SUGAR (2-(ACETYLAMINO)-2-DEOXY-A-D-GLUCOPYRANOSE) (three-letter code: NDG) (formula: C₈H₁₅NO₆).



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

- Molecule 5 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total	Hg	0	0
			2	2		
5	A	2	Total	Hg	0	0
			2	2		

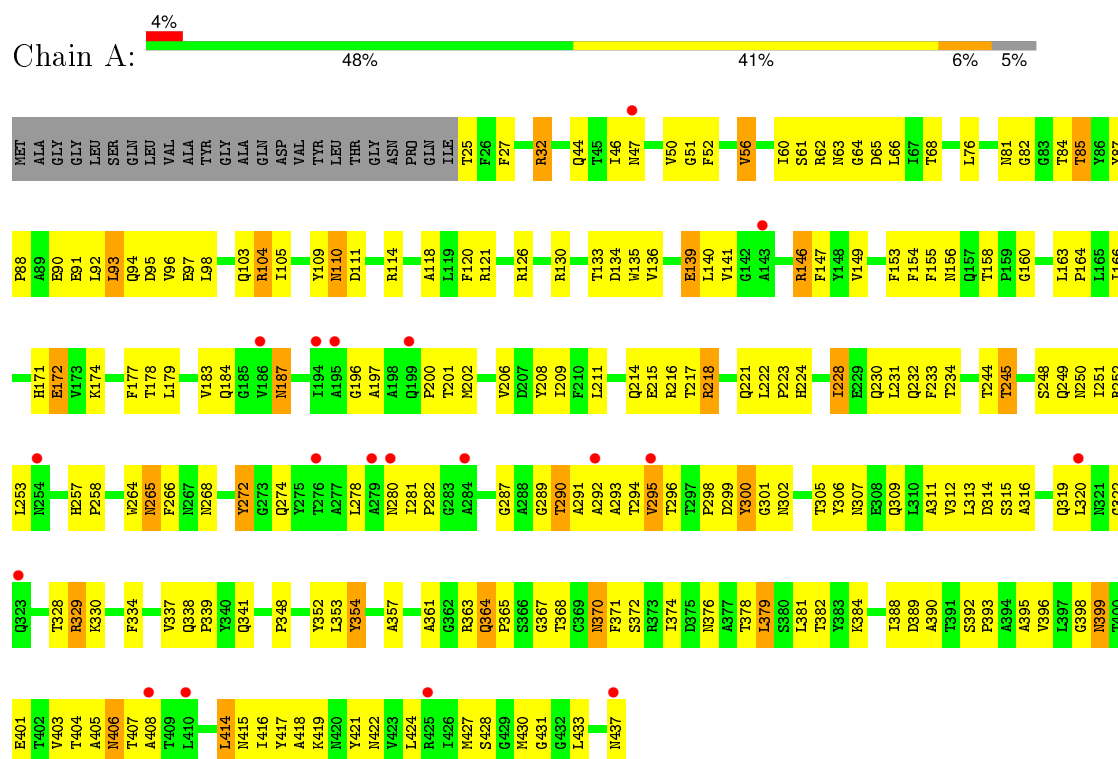
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	23	Total	O	0	0
			23	23		
6	B	27	Total	O	0	0
			27	27		

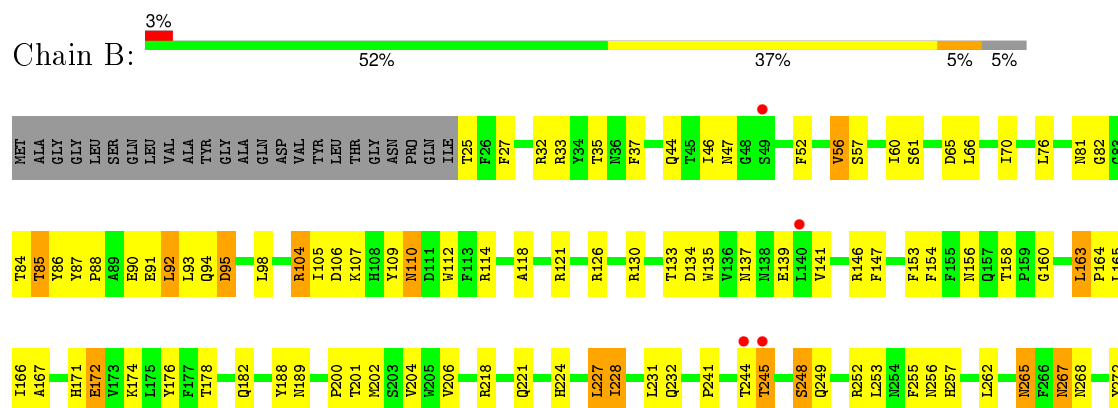
3 Residue-property plots [i](#)

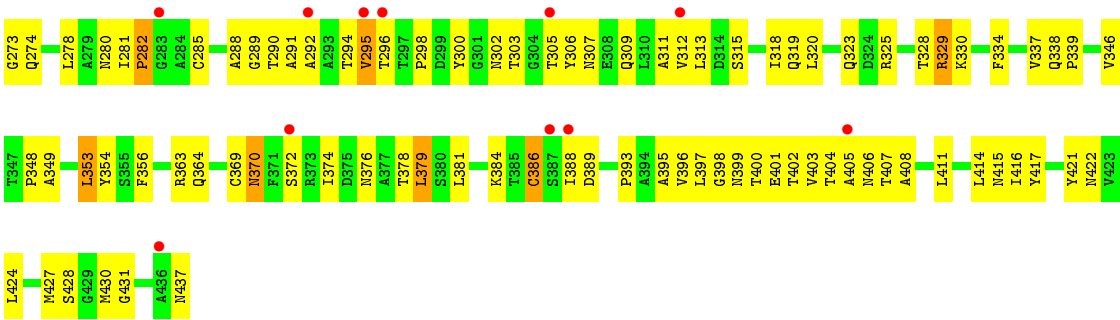
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ($RSRZ > 2$). 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: Major capsid protein



• Molecule 1: Major capsid protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 41 3 2	Depositor
Cell constants a, b, c, α , β , γ	188.76 Å 188.76 Å 188.76 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	84.42 – 2.55 84.42 – 2.54	Depositor EDS
% Data completeness (in resolution range)	99.4 (84.42-2.55) 99.6 (84.42-2.54)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.87 (at 2.55 Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.269 , 0.309 0.272 , 0.314	Depositor DCC
R_{free} test set	1514 reflections (4.18%)	DCC
Wilson B-factor (Å ²)	36.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 33.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 71863 reflections	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	6956	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.79% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HG, NAG, NDG, 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	0.39	0/3307	0.65	0/4506
1	B	0.40	0/3307	0.65	1/4506 (0.0%)
All	All	0.40	0/6614	0.65	1/9012 (0.0%)

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

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

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	189	ASN	N-CA-C	-5.28	96.74	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	354	TYR	Sidechain

5.2 Too-close contacts [i](#)

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

the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3232	0	3127	237	0
1	B	3232	0	3129	211	0
2	A	45	0	45	45	0
2	B	45	0	45	55	0
3	A	144	0	144	58	0
3	B	144	0	144	66	0
4	A	30	0	30	30	0
4	B	30	0	30	17	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
6	A	23	0	0	1	0
6	B	27	0	0	1	0
All	All	6956	0	6694	570	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 42.

All (570) 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:389:ASP:H	4:A:449:NDG:H8C3	1.03	1.16
1:A:393:PRO:HG2	2:A:438:NAG:H82	1.19	1.15
2:B:538:NAG:H3	3:B:540:MAN:H2	1.26	1.13
1:B:389:ASP:H	2:B:549:NAG:H81	1.10	1.11
3:A:443:MAN:H1	3:A:447:MAN:O2	1.53	1.08
1:B:302:ASN:ND2	2:B:546:NAG:H61	1.67	1.08
2:B:538:NAG:C3	3:B:540:MAN:H2	1.85	1.07
1:A:280:ASN:HD21	2:A:438:NAG:C1	1.69	1.05
2:A:438:NAG:O3	3:A:440:MAN:H2	1.59	1.01
3:A:443:MAN:H1	2:A:444:NAG:O4	1.62	1.00
1:B:395:ALA:HB1	2:B:549:NAG:H82	1.44	0.99
1:A:399:ASN:HD21	4:A:449:NDG:H1	1.28	0.99
1:A:294:THR:HG22	1:A:295:VAL:H	1.27	0.97
1:A:280:ASN:HD21	2:A:438:NAG:H1	1.25	0.96
3:A:441:MAN:O4	3:A:442:MAN:H1	1.66	0.95
2:A:444:NAG:H61	4:A:446:NDG:O4	1.64	0.95
1:B:135:TRP:CD1	1:B:139:GLU:HG3	2.01	0.94
1:B:290:THR:HG22	1:B:292:ALA:H	1.28	0.94
3:A:447:MAN:H4	3:A:448:MAN:O1	1.68	0.94
1:B:302:ASN:HD21	2:B:546:NAG:H61	1.31	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:399:ASN:HD21	4:A:449:NDG:C1	1.81	0.93
2:A:438:NAG:HO3	3:A:440:MAN:H2	1.32	0.92
1:B:406:ASN:HD21	4:B:550:NDG:HA	0.95	0.90
1:A:389:ASP:N	4:A:449:NDG:H8C3	1.85	0.90
3:A:443:MAN:H1	3:A:447:MAN:HO2	1.30	0.89
1:A:389:ASP:H	4:A:449:NDG:C8	1.87	0.88
2:B:538:NAG:H3	3:B:540:MAN:C2	2.03	0.88
3:B:541:MAN:O4	3:B:542:MAN:H5	1.73	0.88
2:B:549:NAG:C6	3:B:553:MAN:H5	2.04	0.88
1:A:135:TRP:HB3	1:A:139:GLU:HG3	1.56	0.87
1:A:171:HIS:HD2	1:A:431:GLY:H	1.22	0.87
1:B:82:GLY:H	3:B:543:MAN:C6	1.88	0.86
1:A:290:THR:HG23	1:A:292:ALA:H	1.38	0.86
3:A:447:MAN:H61	3:A:448:MAN:H5	1.57	0.86
1:B:389:ASP:N	2:B:549:NAG:H81	1.91	0.86
1:B:81:ASN:HB3	3:B:543:MAN:H61	1.56	0.85
1:B:294:THR:HG22	1:B:295:VAL:H	1.38	0.85
1:B:406:ASN:ND2	4:B:550:NDG:HA	1.74	0.84
2:B:538:NAG:HO6	3:B:541:MAN:H2	1.41	0.84
1:B:353:LEU:HD22	1:B:354:TYR:N	1.92	0.84
1:A:406:ASN:HD22	1:A:407:THR:N	1.76	0.83
1:A:393:PRO:CG	2:A:438:NAG:H82	2.05	0.82
1:B:289:GLY:HA2	2:B:538:NAG:O7	1.79	0.82
1:A:158:THR:HG22	1:A:160:GLY:H	1.43	0.82
2:A:438:NAG:O3	3:A:440:MAN:C2	2.27	0.82
1:B:294:THR:HG22	1:B:296:THR:H	1.45	0.81
1:B:46:ILE:HG23	1:B:56:VAL:HG21	1.62	0.81
1:A:393:PRO:HG2	2:A:438:NAG:C8	2.07	0.81
1:A:183:VAL:H	1:A:187:ASN:HD21	1.28	0.81
1:A:244:THR:O	1:A:245:THR:HG22	1.81	0.81
1:A:25:THR:HG22	1:A:27:PHE:H	1.44	0.81
1:A:90:GLU:HB2	1:A:133:THR:HG21	1.63	0.81
1:A:217:THR:O	1:A:221:GLN:HG2	1.80	0.80
1:A:44:GLN:HE22	1:A:61:SER:H	1.28	0.80
1:B:389:ASP:H	2:B:549:NAG:C8	1.92	0.79
2:A:444:NAG:H4	3:A:447:MAN:O1	1.83	0.79
3:A:443:MAN:C1	2:A:444:NAG:O4	2.29	0.79
3:B:547:MAN:O4	3:B:548:MAN:H1	1.81	0.79
3:B:543:MAN:H2	3:B:545:MAN:H1	1.64	0.79
1:A:384:LYS:HD2	1:A:404:THR:HG23	1.65	0.79
1:B:384:LYS:HD2	1:B:404:THR:HG23	1.65	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:82:GLY:H	3:B:543:MAN:H61	1.46	0.78
1:B:46:ILE:CG2	1:B:56:VAL:HG21	2.13	0.78
1:A:146:ARG:HH11	1:A:146:ARG:HG2	1.49	0.78
2:B:538:NAG:C4	3:B:540:MAN:H2	2.13	0.78
1:B:171:HIS:CD2	1:B:431:GLY:H	2.01	0.78
1:B:305:THR:HG22	1:B:307:ASN:H	1.48	0.77
1:A:305:THR:HG22	1:A:307:ASN:H	1.48	0.77
1:B:91:GLU:HA	1:B:110:ASN:OD1	1.84	0.77
1:B:302:ASN:HD21	2:B:546:NAG:C6	1.98	0.77
1:B:46:ILE:HD13	1:B:202:MET:HE2	1.67	0.77
1:A:319:GLN:HE21	1:A:322:GLY:HA2	1.49	0.77
1:B:312:VAL:HG12	1:B:408:ALA:HB1	1.68	0.76
1:B:56:VAL:HG11	1:B:202:MET:HE1	1.67	0.76
2:B:538:NAG:O3	3:B:541:MAN:H2	1.85	0.76
2:B:538:NAG:C8	3:B:552:MAN:H5	2.14	0.76
1:A:278:LEU:HD21	1:A:298:PRO:HB3	1.68	0.76
3:A:443:MAN:C1	2:A:444:NAG:HO4	1.99	0.75
1:A:65:ASP:HB3	1:A:166:ILE:HD13	1.69	0.75
1:B:384:LYS:HD2	1:B:404:THR:CG2	2.17	0.75
4:B:544:NDG:O3	3:B:547:MAN:H4	1.87	0.75
3:B:543:MAN:HO3	3:B:545:MAN:C5	1.98	0.75
1:B:171:HIS:HD2	1:B:431:GLY:H	1.34	0.75
1:A:280:ASN:ND2	2:A:438:NAG:C1	2.49	0.75
1:A:257:HIS:CE1	1:A:437:ASN:HB2	2.23	0.74
3:B:539:MAN:H62	3:B:541:MAN:H3	1.68	0.74
1:B:174:LYS:HB3	1:B:174:LYS:HZ3	1.53	0.74
1:B:65:ASP:HA	1:B:166:ILE:HG23	1.70	0.74
1:A:311:ALA:O	1:A:404:THR:HG21	1.88	0.73
2:B:549:NAG:H61	3:B:553:MAN:H61	1.69	0.73
1:A:330:LYS:HE2	1:A:401:GLU:OE1	1.89	0.73
1:B:257:HIS:NE2	1:B:437:ASN:HB2	2.03	0.73
1:A:294:THR:HG22	1:A:296:THR:H	1.52	0.73
3:B:541:MAN:H61	3:B:542:MAN:O6	1.89	0.72
1:A:171:HIS:CD2	1:A:431:GLY:H	2.05	0.72
1:B:406:ASN:ND2	4:B:550:NDG:H1	2.04	0.72
2:A:438:NAG:H81	3:A:452:MAN:H5	1.71	0.72
1:B:280:ASN:ND2	2:B:538:NAG:C1	2.52	0.72
1:B:265:ASN:HD21	1:B:415:ASN:HD22	1.38	0.72
1:A:47:ASN:O	1:A:56:VAL:HG22	1.89	0.72
1:A:388:ILE:HG23	4:A:449:NDG:H8C1	1.71	0.72
1:B:90:GLU:HB2	1:B:133:THR:HG21	1.70	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:280:ASN:ND2	2:B:538:NAG:H1	2.05	0.71
1:B:153:PHE:H	1:B:156:ASN:HD22	1.38	0.71
1:A:290:THR:CG2	1:A:292:ALA:H	2.02	0.71
2:A:438:NAG:C8	3:A:452:MAN:H5	2.20	0.71
1:B:47:ASN:O	1:B:56:VAL:HG23	1.91	0.71
1:A:399:ASN:ND2	4:A:449:NDG:C1	2.53	0.70
1:B:46:ILE:HG23	1:B:56:VAL:CG2	2.21	0.70
2:A:444:NAG:C6	4:A:446:NDG:O4	2.38	0.70
2:A:438:NAG:H4	3:A:441:MAN:H2	1.72	0.70
1:A:278:LEU:HB2	1:A:407:THR:HG21	1.74	0.70
1:B:174:LYS:NZ	1:B:174:LYS:HB3	2.07	0.70
1:B:66:LEU:HD23	1:B:164:PRO:HA	1.72	0.70
1:B:280:ASN:HD21	2:B:538:NAG:C1	2.05	0.70
3:B:543:MAN:H2	3:B:545:MAN:C1	2.21	0.69
1:B:88:PRO:HD2	1:B:134:ASP:CB	2.23	0.69
3:B:541:MAN:HO4	3:B:542:MAN:H5	1.56	0.69
3:A:452:MAN:O6	3:A:452:MAN:H1	1.92	0.68
1:A:218:ARG:HH21	1:A:224:HIS:CE1	2.12	0.68
2:A:438:NAG:O3	3:A:441:MAN:O2	2.06	0.68
3:A:443:MAN:C1	3:A:447:MAN:O2	2.37	0.68
1:B:104:ARG:HG3	1:B:104:ARG:O	1.93	0.68
1:A:312:VAL:HG12	1:A:408:ALA:HB1	1.75	0.67
1:B:280:ASN:HD22	1:B:393:PRO:HG3	1.58	0.67
1:A:153:PHE:H	1:A:156:ASN:HD22	1.42	0.67
1:A:52:PHE:CE2	1:A:200:PRO:HD3	2.29	0.67
1:B:98:LEU:HD12	1:B:105:ILE:HB	1.75	0.67
1:B:88:PRO:HD2	1:B:134:ASP:HB3	1.77	0.67
1:A:289:GLY:HA2	2:A:438:NAG:O1	1.95	0.67
4:B:544:NDG:O	2:B:546:NAG:O4	2.12	0.67
1:A:25:THR:HG22	1:A:27:PHE:N	2.10	0.67
2:A:438:NAG:C4	3:A:441:MAN:H2	2.25	0.67
2:B:549:NAG:H61	3:B:553:MAN:H5	1.74	0.67
1:B:163:LEU:HD22	1:B:164:PRO:HD2	1.76	0.66
2:B:538:NAG:H3	3:B:541:MAN:O2	1.95	0.66
2:B:549:NAG:H62	3:B:553:MAN:H5	1.76	0.66
1:A:294:THR:O	1:A:295:VAL:HG23	1.96	0.66
1:A:329:ARG:HB3	1:A:329:ARG:HH11	1.61	0.66
1:B:320:LEU:HD11	1:B:325:ARG:HD3	1.77	0.66
1:B:320:LEU:HD22	1:B:374:ILE:HD13	1.76	0.66
1:B:414:LEU:HD22	1:B:416:ILE:HG13	1.78	0.66
1:B:44:GLN:NE2	1:B:60:ILE:HA	2.11	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:399:ASN:HD21	2:B:549:NAG:C1	2.09	0.66
1:A:379:LEU:HD13	1:A:381:LEU:HD21	1.78	0.66
3:A:443:MAN:H2	3:A:445:MAN:H1	1.79	0.65
1:B:82:GLY:H	3:B:543:MAN:H62	1.58	0.65
1:A:68:THR:HG23	1:A:209:ILE:HD13	1.78	0.65
1:A:294:THR:HG22	1:A:295:VAL:N	2.08	0.65
1:A:312:VAL:CG1	1:A:408:ALA:HB1	2.26	0.65
1:A:329:ARG:NH1	1:A:329:ARG:HB3	2.12	0.65
1:B:244:THR:O	1:B:245:THR:HG22	1.97	0.65
1:A:399:ASN:ND2	4:A:449:NDG:H1	2.08	0.64
1:A:251:ILE:HD12	1:A:251:ILE:N	2.12	0.64
1:B:353:LEU:HD22	1:B:354:TYR:H	1.59	0.64
3:A:447:MAN:C4	3:A:448:MAN:O1	2.45	0.64
3:A:447:MAN:H4	3:A:448:MAN:HO1	1.63	0.63
1:B:46:ILE:HD13	1:B:202:MET:CE	2.28	0.63
1:B:338:GLN:HB3	1:B:353:LEU:HD13	1.81	0.63
3:B:540:MAN:H3	3:B:541:MAN:O2	1.98	0.63
1:A:187:ASN:H	1:A:187:ASN:HD22	1.46	0.63
1:A:302:ASN:ND2	4:A:446:NDG:O	2.31	0.63
1:A:395:ALA:HB1	4:A:449:NDG:C8	2.29	0.62
1:B:44:GLN:HE21	1:B:60:ILE:HA	1.64	0.62
3:B:541:MAN:C6	3:B:542:MAN:HO6	2.13	0.62
1:B:388:ILE:HA	2:B:549:NAG:O7	2.00	0.62
1:A:309:GLN:O	1:A:404:THR:HG22	2.00	0.62
1:B:329:ARG:NH1	1:B:329:ARG:HB3	2.13	0.62
1:A:135:TRP:CD1	1:A:139:GLU:HG3	2.35	0.61
2:B:538:NAG:C3	3:B:541:MAN:O2	2.48	0.61
1:B:44:GLN:HE22	1:B:61:SER:H	1.48	0.61
1:B:290:THR:HG22	1:B:292:ALA:N	2.09	0.61
4:A:449:NDG:O3	3:A:454:MAN:O5	2.17	0.61
4:B:544:NDG:HB	3:B:547:MAN:C5	2.13	0.61
1:A:153:PHE:H	1:A:156:ASN:ND2	1.98	0.61
1:B:320:LEU:CD1	1:B:325:ARG:HD3	2.31	0.61
1:B:400:THR:HG22	1:B:402:THR:H	1.65	0.61
1:B:25:THR:HG22	1:B:27:PHE:H	1.65	0.61
3:A:447:MAN:C6	3:A:448:MAN:O1	2.49	0.60
1:B:311:ALA:O	1:B:404:THR:HG21	2.00	0.60
1:B:267:ASN:HB2	1:B:274:GLN:O	2.01	0.60
1:B:135:TRP:HD1	1:B:139:GLU:HG3	1.64	0.60
1:A:118:ALA:HB1	1:A:348:PRO:HD2	1.83	0.60
1:B:232:GLN:HE22	1:B:255:PHE:HA	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:ASN:O	1:A:393:PRO:HG3	2.01	0.60
1:B:281:ILE:HG23	1:B:282:PRO:HD2	1.82	0.60
1:B:395:ALA:CB	2:B:549:NAG:H82	2.26	0.60
1:A:395:ALA:HB1	4:A:449:NDG:H8C1	1.82	0.60
2:B:538:NAG:H81	3:B:552:MAN:H5	1.84	0.60
1:A:135:TRP:CB	1:A:139:GLU:HG3	2.30	0.60
1:A:25:THR:CG2	1:A:27:PHE:H	2.15	0.60
1:A:46:ILE:HD13	1:A:202:MET:HG2	1.84	0.60
1:B:280:ASN:HD21	2:B:538:NAG:H1	1.66	0.60
2:A:444:NAG:O5	4:A:446:NDG:O4	2.20	0.60
1:A:257:HIS:NE2	1:A:437:ASN:HB2	2.16	0.60
4:A:449:NDG:O4	3:A:453:MAN:O1	2.14	0.60
1:A:337:VAL:O	1:A:341:GLN:HG3	2.01	0.60
1:A:44:GLN:NE2	1:A:60:ILE:HA	2.16	0.59
1:B:87:TYR:N	1:B:88:PRO:HD3	2.17	0.59
1:A:82:GLY:H	3:A:443:MAN:C6	2.15	0.59
1:B:280:ASN:O	1:B:280:ASN:ND2	2.35	0.59
2:B:549:NAG:O1	2:B:549:NAG:C7	2.50	0.59
1:B:218:ARG:HH21	1:B:224:HIS:CE1	2.20	0.59
1:A:258:PRO:HB3	1:A:368:THR:HG21	1.85	0.59
1:A:299:ASP:O	1:A:301:GLY:N	2.36	0.59
1:A:312:VAL:O	1:A:384:LYS:HG3	2.03	0.59
1:A:265:ASN:C	1:A:265:ASN:HD22	2.05	0.59
1:A:257:HIS:NE2	1:A:437:ASN:C	2.56	0.59
3:A:439:MAN:O5	3:A:441:MAN:O3	2.21	0.58
1:A:62:ARG:NH2	1:A:166:ILE:HA	2.18	0.58
2:B:538:NAG:O4	3:B:540:MAN:H2	2.02	0.58
1:B:98:LEU:CD1	1:B:105:ILE:HB	2.33	0.58
1:B:121:ARG:HG2	1:B:126:ARG:HA	1.84	0.58
1:A:231:LEU:HD13	1:A:232:GLN:N	2.18	0.58
1:A:274:GLN:NE2	1:A:278:LEU:HG	2.18	0.58
3:B:547:MAN:O6	3:B:548:MAN:H1	2.03	0.58
1:A:94:GLN:HA	1:A:110:ASN:ND2	2.19	0.58
3:A:452:MAN:O6	3:A:452:MAN:C1	2.50	0.58
1:B:81:ASN:OD1	1:B:141:VAL:HG23	2.03	0.58
1:A:95:ASP:OD1	1:A:178:THR:HB	2.04	0.58
1:B:399:ASN:ND2	2:B:549:NAG:O1	2.27	0.58
1:A:98:LEU:HD12	1:A:105:ILE:HB	1.84	0.58
1:B:232:GLN:NE2	1:B:255:PHE:HA	2.18	0.57
2:B:549:NAG:H61	3:B:553:MAN:C6	2.34	0.57
1:A:65:ASP:HB3	1:A:166:ILE:CD1	2.33	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:THR:HG23	1:A:88:PRO:HD3	1.84	0.57
1:B:330:LYS:HE2	1:B:401:GLU:OE1	2.03	0.57
1:A:314:ASP:HB3	1:A:382:THR:O	2.04	0.57
3:A:439:MAN:C1	3:A:441:MAN:O3	2.51	0.57
1:A:120:PHE:HD2	1:A:361:ALA:HB2	1.69	0.57
1:B:302:ASN:HD21	2:B:546:NAG:C5	2.18	0.57
1:B:338:GLN:HB2	1:B:339:PRO:HD3	1.86	0.57
1:A:281:ILE:HG23	1:A:282:PRO:HD2	1.87	0.57
1:B:82:GLY:N	3:B:543:MAN:H62	2.20	0.57
1:A:130:ARG:HG2	1:A:136:VAL:CG1	2.35	0.57
1:B:370:ASN:ND2	1:B:372:SER:H	2.03	0.57
1:A:218:ARG:NH2	1:A:224:HIS:CE1	2.73	0.57
1:B:329:ARG:HH11	1:B:329:ARG:HB3	1.69	0.57
1:B:290:THR:CG2	1:B:291:ALA:N	2.68	0.56
1:A:280:ASN:HD22	1:A:393:PRO:HD3	1.70	0.56
3:B:541:MAN:H61	3:B:542:MAN:HO6	1.70	0.56
1:B:52:PHE:CE2	1:B:200:PRO:HD3	2.40	0.56
4:A:449:NDG:HC	3:A:453:MAN:HO1	1.49	0.56
2:B:538:NAG:O3	3:B:541:MAN:C2	2.52	0.56
1:A:302:ASN:HD21	4:A:446:NDG:C5	2.18	0.56
1:B:406:ASN:CG	4:B:550:NDG:H1	2.26	0.56
1:B:281:ILE:HD13	1:B:397:LEU:HD21	1.87	0.56
2:B:538:NAG:H82	3:B:552:MAN:H5	1.86	0.56
3:B:547:MAN:O4	3:B:548:MAN:C1	2.51	0.56
1:A:93:LEU:HD21	1:A:177:PHE:HD2	1.70	0.56
1:A:172:GLU:OE1	1:A:174:LYS:HD2	2.05	0.56
2:B:538:NAG:H3	3:B:540:MAN:C1	2.35	0.56
1:A:183:VAL:H	1:A:187:ASN:ND2	2.00	0.56
2:B:538:NAG:O6	3:B:541:MAN:H2	2.02	0.56
1:A:81:ASN:HB3	3:A:443:MAN:H61	1.86	0.56
1:B:406:ASN:ND2	4:B:550:NDG:C1	2.68	0.56
1:A:88:PRO:HD2	1:A:134:ASP:CB	2.36	0.56
1:A:32:ARG:H	1:A:32:ARG:HD2	1.71	0.56
1:B:320:LEU:HG	1:B:325:ARG:HB2	1.88	0.55
1:A:305:THR:HG22	1:A:306:TYR:N	2.22	0.55
1:B:85:THR:HG23	1:B:88:PRO:HD3	1.87	0.55
1:B:257:HIS:CE1	1:B:437:ASN:HB2	2.42	0.55
1:A:66:LEU:HD23	1:A:164:PRO:HA	1.88	0.55
1:A:370:ASN:ND2	1:A:372:SER:H	2.05	0.55
1:A:290:THR:HG23	1:A:291:ALA:N	2.22	0.55
1:B:84:THR:HG22	1:B:84:THR:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:543:MAN:HO3	3:B:545:MAN:C6	2.20	0.55
1:B:295:VAL:O	1:B:295:VAL:HG12	2.07	0.55
1:A:133:THR:HG22	1:A:133:THR:O	2.06	0.55
2:B:549:NAG:H61	3:B:553:MAN:C5	2.37	0.54
2:A:444:NAG:O1	4:A:446:NDG:H5	2.07	0.54
2:A:444:NAG:C5	4:A:446:NDG:HC	2.19	0.54
3:A:440:MAN:H2	3:A:441:MAN:O2	2.08	0.54
1:B:398:GLY:O	2:B:549:NAG:H62	2.06	0.54
1:A:171:HIS:HD2	1:A:431:GLY:N	2.00	0.54
1:A:363:ARG:HA	1:A:363:ARG:HE	1.73	0.54
4:A:449:NDG:O7	3:A:454:MAN:H61	2.07	0.54
1:B:406:ASN:ND2	4:B:550:NDG:N2	2.44	0.54
1:B:294:THR:HG22	1:B:295:VAL:N	2.16	0.54
1:B:280:ASN:ND2	2:B:538:NAG:O1	2.40	0.54
1:B:66:LEU:CD2	1:B:164:PRO:HA	2.38	0.54
1:B:86:TYR:HE2	1:B:273:GLY:HA3	1.73	0.54
3:B:541:MAN:C6	3:B:542:MAN:O6	2.56	0.54
1:B:312:VAL:O	1:B:384:LYS:HG3	2.07	0.54
1:B:281:ILE:HD11	1:B:396:VAL:HB	1.90	0.54
1:B:82:GLY:HA3	3:B:543:MAN:H62	1.89	0.54
1:A:302:ASN:ND2	4:A:446:NDG:H6C2	2.22	0.54
1:A:320:LEU:CD2	1:A:374:ILE:HD13	2.38	0.53
1:B:280:ASN:OD1	2:B:538:NAG:O1	2.18	0.53
1:A:140:LEU:HD23	3:A:443:MAN:H61	1.90	0.53
1:A:184:GLN:HG2	1:A:300:TYR:CD2	2.43	0.53
1:A:384:LYS:HD2	1:A:404:THR:CG2	2.36	0.53
1:A:60:ILE:HD12	1:A:206:VAL:HG21	1.89	0.53
1:A:111:ASP:HB3	1:A:417:TYR:OH	2.08	0.53
1:B:319:GLN:HG3	1:B:323:GLN:O	2.08	0.53
1:A:287:GLY:O	1:A:293:ALA:HB2	2.07	0.53
1:B:158:THR:HG22	1:B:160:GLY:H	1.72	0.53
1:B:278:LEU:HD21	1:B:298:PRO:HB3	1.91	0.53
1:A:62:ARG:HH21	1:A:166:ILE:HA	1.74	0.53
1:B:379:LEU:HD12	1:B:416:ILE:CD1	2.39	0.53
2:B:549:NAG:HO3	3:B:554:MAN:C5	2.21	0.52
1:A:134:ASP:HA	1:A:147:PHE:CE2	2.44	0.52
1:B:379:LEU:HD12	1:B:416:ILE:HD11	1.90	0.52
1:A:250:ASN:C	1:A:251:ILE:HD12	2.30	0.52
1:A:230:GLN:O	1:A:419:LYS:HD2	2.09	0.52
1:B:338:GLN:HB3	1:B:353:LEU:CD1	2.39	0.52
1:A:171:HIS:CD2	1:A:430:MET:HA	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:543:MAN:O3	3:B:545:MAN:O6	2.26	0.52
1:A:305:THR:HG22	1:A:307:ASN:N	2.21	0.52
1:A:82:GLY:H	3:A:443:MAN:H62	1.74	0.52
1:A:94:GLN:HG3	1:A:110:ASN:ND2	2.25	0.52
1:B:94:GLN:HB3	1:B:178:THR:O	2.08	0.52
1:B:244:THR:O	1:B:245:THR:O	2.28	0.52
1:B:252:ARG:NE	1:B:376:ASN:ND2	2.57	0.52
2:A:438:NAG:H83	3:A:452:MAN:C5	2.40	0.52
1:B:65:ASP:HB3	1:B:166:ILE:HD13	1.91	0.52
4:B:544:NDG:O3	3:B:547:MAN:C4	2.57	0.52
1:B:118:ALA:HB1	1:B:348:PRO:HD2	1.91	0.52
1:B:265:ASN:HD22	1:B:265:ASN:C	2.14	0.52
1:B:278:LEU:HB2	1:B:407:THR:HG21	1.92	0.52
1:B:305:THR:HG22	1:B:307:ASN:N	2.22	0.51
1:A:135:TRP:CG	1:A:139:GLU:HG3	2.46	0.51
1:A:87:TYR:N	1:A:88:PRO:HD3	2.26	0.51
1:B:95:ASP:OD2	1:B:178:THR:HB	2.10	0.51
1:A:223:PRO:HB3	1:A:427:MET:HE2	1.92	0.51
1:A:146:ARG:NH1	1:A:146:ARG:HG2	2.22	0.51
1:A:364:GLN:NE2	1:A:365:PRO:O	2.44	0.51
1:B:285:CYS:SG	1:B:288:ALA:HB2	2.50	0.51
2:B:538:NAG:O6	3:B:539:MAN:H5	2.10	0.51
1:A:221:GLN:HE21	1:A:221:GLN:HA	1.75	0.51
1:A:265:ASN:C	1:A:265:ASN:ND2	2.63	0.51
4:B:544:NDG:HA	3:B:547:MAN:C6	2.24	0.51
1:B:298:PRO:HG2	1:B:300:TYR:CE2	2.45	0.51
1:A:91:GLU:HA	1:A:110:ASN:OD1	2.10	0.51
1:B:46:ILE:HD11	1:B:204:VAL:HG23	1.93	0.50
1:B:114:ARG:HD2	1:B:272:TYR:CE2	2.45	0.50
1:B:227:LEU:HD13	1:B:421:TYR:CD1	2.46	0.50
1:A:135:TRP:HB3	1:A:139:GLU:CG	2.36	0.50
1:A:81:ASN:OD1	1:A:141:VAL:HG23	2.11	0.50
1:A:46:ILE:CD1	1:A:202:MET:HG2	2.41	0.50
1:B:253:LEU:HD22	1:B:255:PHE:CE1	2.47	0.50
1:B:370:ASN:HD22	1:B:370:ASN:C	2.14	0.50
3:B:547:MAN:O6	3:B:548:MAN:C1	2.59	0.50
1:B:274:GLN:N	1:B:303:THR:HG21	2.27	0.50
1:B:94:GLN:HG2	1:B:95:ASP:OD1	2.12	0.50
1:B:137:ASN:N	1:B:139:GLU:OE2	2.40	0.50
2:A:444:NAG:C5	4:A:446:NDG:O4	2.59	0.49
1:A:130:ARG:HG2	1:A:136:VAL:HG11	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:395:ALA:O	1:A:399:ASN:ND2	2.44	0.49
1:B:86:TYR:CE2	1:B:303:THR:HB	2.46	0.49
1:A:281:ILE:HD11	1:A:396:VAL:HB	1.93	0.49
2:B:538:NAG:O4	2:B:538:NAG:N2	2.41	0.49
1:B:174:LYS:NZ	1:B:176:TYR:CZ	2.80	0.49
2:A:438:NAG:HO3	3:A:441:MAN:C2	2.19	0.49
2:A:438:NAG:C8	3:A:452:MAN:C5	2.88	0.49
1:B:338:GLN:HG3	1:B:353:LEU:O	2.12	0.49
1:A:163:LEU:HD23	1:A:164:PRO:HD2	1.94	0.49
1:A:232:GLN:HB2	1:A:418:ALA:HB3	1.95	0.49
1:A:370:ASN:HD22	1:A:370:ASN:C	2.14	0.49
1:A:103:GLN:HE22	1:A:433:LEU:HD12	1.78	0.49
1:A:44:GLN:HE21	1:A:60:ILE:HA	1.77	0.49
1:B:414:LEU:CD2	1:B:416:ILE:HG13	2.41	0.49
1:B:274:GLN:NE2	1:B:278:LEU:HG	2.27	0.49
3:A:445:MAN:H1	3:A:448:MAN:H2	1.93	0.49
1:B:221:GLN:OE1	1:B:221:GLN:HA	2.13	0.49
1:B:386:CYS:HB2	1:B:406:ASN:HA	1.93	0.48
1:A:76:LEU:CD2	1:A:200:PRO:HB3	2.43	0.48
1:A:251:ILE:O	1:A:378:THR:HG23	2.13	0.48
1:B:329:ARG:HB2	1:B:334:PHE:CE1	2.48	0.48
1:A:398:GLY:O	4:A:449:NDG:H6C2	2.13	0.48
1:B:280:ASN:CG	2:B:538:NAG:HO1	2.14	0.48
1:B:312:VAL:HG23	1:B:313:LEU:N	2.27	0.48
1:A:389:ASP:CB	4:A:449:NDG:H8C3	2.44	0.48
1:A:280:ASN:ND2	1:A:393:PRO:HD3	2.29	0.48
2:A:438:NAG:H83	3:A:452:MAN:H5	1.94	0.48
1:B:312:VAL:CG1	1:B:408:ALA:HB1	2.40	0.48
1:A:87:TYR:OH	1:A:130:ARG:HD2	2.13	0.48
1:A:97:GLU:CD	1:A:104:ARG:HH11	2.16	0.48
3:A:443:MAN:C3	3:A:445:MAN:H1	2.43	0.48
1:A:82:GLY:H	3:A:443:MAN:H61	1.79	0.48
1:B:378:THR:HG22	1:B:379:LEU:N	2.29	0.48
1:A:309:GLN:NE2	1:A:403:VAL:O	2.46	0.48
1:A:312:VAL:HG23	1:A:313:LEU:N	2.29	0.48
3:B:540:MAN:C3	3:B:541:MAN:O2	2.61	0.48
1:B:130:ARG:HG2	1:B:134:ASP:OD2	2.14	0.48
2:A:438:NAG:O4	3:A:441:MAN:C2	2.62	0.48
1:A:154:PHE:CD1	1:A:155:PHE:N	2.82	0.48
1:A:114:ARG:NE	1:A:114:ARG:HA	2.29	0.48
1:B:309:GLN:O	1:B:404:THR:HG22	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:280:ASN:ND2	1:B:393:PRO:HG3	2.27	0.47
2:B:549:NAG:O3	3:B:554:MAN:C1	2.62	0.47
2:A:444:NAG:O1	4:A:446:NDG:H6C1	2.14	0.47
1:A:353:LEU:HD13	1:A:354:TYR:N	2.28	0.47
1:B:379:LEU:HD13	1:B:381:LEU:HD11	1.97	0.47
1:B:82:GLY:CA	3:B:543:MAN:H62	2.44	0.47
1:B:82:GLY:HA3	4:B:544:NDG:H6C1	1.96	0.47
1:A:85:THR:CG2	1:A:134:ASP:HB2	2.45	0.47
1:A:65:ASP:CG	1:A:216:ARG:HD3	2.35	0.47
1:B:87:TYR:CE2	1:B:130:ARG:HG3	2.50	0.47
1:A:370:ASN:HD22	1:A:372:SER:H	1.63	0.47
1:A:51:GLY:O	1:A:179:LEU:HD12	2.13	0.47
1:A:253:LEU:HD21	1:A:416:ILE:HD13	1.95	0.47
1:B:85:THR:HG23	1:B:88:PRO:CD	2.45	0.47
1:A:388:ILE:HD11	1:A:403:VAL:HG23	1.96	0.47
1:A:65:ASP:O	1:A:166:ILE:HG12	2.15	0.47
1:A:320:LEU:HD21	1:A:374:ILE:HD13	1.96	0.47
1:A:218:ARG:HH21	1:A:224:HIS:HE1	1.57	0.46
1:B:86:TYR:CE2	1:B:273:GLY:HA3	2.51	0.46
1:B:384:LYS:CD	1:B:404:THR:HG23	2.41	0.46
3:A:443:MAN:H3	3:A:445:MAN:H1	1.97	0.46
3:A:447:MAN:H61	3:A:448:MAN:O1	2.15	0.46
1:A:320:LEU:HD22	1:A:374:ILE:HG21	1.96	0.46
1:B:231:LEU:HD11	1:B:417:TYR:HD2	1.80	0.46
2:A:444:NAG:C4	3:A:447:MAN:O1	2.61	0.46
1:A:114:ARG:HD2	1:A:272:TYR:CE2	2.50	0.46
1:A:392:SER:O	1:A:396:VAL:HG23	2.15	0.46
1:A:302:ASN:HD21	4:A:446:NDG:H6C2	1.79	0.46
1:B:280:ASN:O	1:B:281:ILE:HG13	2.16	0.46
3:B:547:MAN:O4	3:B:548:MAN:C2	2.64	0.46
1:A:158:THR:HG22	1:A:160:GLY:N	2.20	0.46
1:A:251:ILE:CD1	1:A:251:ILE:N	2.78	0.46
1:B:388:ILE:HB	1:B:404:THR:O	2.15	0.46
2:B:538:NAG:O4	3:B:540:MAN:C2	2.64	0.46
1:A:87:TYR:HB3	1:A:90:GLU:HB3	1.97	0.46
1:B:232:GLN:HE22	1:B:256:ASN:H	1.64	0.46
1:A:93:LEU:HD12	1:A:149:VAL:HG11	1.98	0.46
1:B:88:PRO:HD2	1:B:134:ASP:HB2	1.97	0.45
1:A:338:GLN:HB2	1:A:339:PRO:HD3	1.98	0.45
2:B:538:NAG:H3	3:B:540:MAN:O1	2.16	0.45
1:A:294:THR:CG2	1:A:295:VAL:H	2.06	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:406:ASN:CG	2:A:450:NAG:O1	2.54	0.45
1:A:88:PRO:HD2	1:A:134:ASP:HB3	1.98	0.45
4:B:550:NDG:H8C2	3:B:551:MAN:O2	2.17	0.45
1:A:309:GLN:O	1:A:309:GLN:HG3	2.16	0.45
1:B:281:ILE:HD13	1:B:397:LEU:CD2	2.46	0.45
1:A:221:GLN:NE2	1:A:221:GLN:HA	2.32	0.45
1:B:172:GLU:OE1	1:B:174:LYS:HD2	2.16	0.45
1:A:50:VAL:HG13	1:A:202:MET:HE2	1.99	0.45
2:A:438:NAG:C4	3:A:441:MAN:C2	2.94	0.45
1:A:266:PHE:CE2	1:A:414:LEU:HG	2.52	0.45
2:B:538:NAG:H3	3:B:540:MAN:HO1	1.81	0.45
1:B:133:THR:O	1:B:133:THR:HG22	2.17	0.45
1:A:280:ASN:ND2	2:A:438:NAG:O5	2.50	0.45
1:A:357:ALA:HB2	1:A:367:GLY:O	2.16	0.45
1:B:309:GLN:NE2	1:B:403:VAL:O	2.50	0.44
1:A:56:VAL:HG11	1:A:202:MET:CE	2.47	0.44
1:A:265:ASN:HD21	1:A:415:ASN:HD22	1.65	0.44
1:A:338:GLN:HG3	1:A:353:LEU:O	2.17	0.44
1:B:109:TYR:O	1:B:112:TRP:HB3	2.17	0.44
1:B:305:THR:O	1:B:349:ALA:N	2.47	0.44
1:A:65:ASP:O	1:A:66:LEU:HD23	2.17	0.44
1:B:88:PRO:CD	1:B:134:ASP:CB	2.95	0.44
1:B:228:ILE:HD11	1:B:422:ASN:HB3	1.98	0.44
1:B:337:VAL:O	1:B:338:GLN:C	2.55	0.44
1:A:334:PHE:HB2	1:A:352:TYR:HB3	1.98	0.44
1:A:295:VAL:O	1:A:295:VAL:HG12	2.17	0.44
2:A:438:NAG:H83	3:A:452:MAN:O5	2.17	0.44
1:A:61:SER:OG	1:A:63:ASN:ND2	2.51	0.44
4:B:544:NDG:HA	3:B:547:MAN:H62	1.83	0.44
2:A:450:NAG:H83	3:A:451:MAN:O3	2.17	0.44
1:A:389:ASP:HB3	4:A:449:NDG:H8C3	1.99	0.44
1:B:265:ASN:ND2	1:B:265:ASN:C	2.70	0.44
1:A:121:ARG:HG2	1:A:126:ARG:HA	1.99	0.44
1:A:406:ASN:OD1	2:A:450:NAG:C1	2.66	0.44
1:A:319:GLN:HE21	1:A:322:GLY:CA	2.26	0.44
1:A:84:THR:HG22	1:A:84:THR:O	2.18	0.44
1:B:312:VAL:CG1	1:B:411:LEU:HD12	2.47	0.43
1:B:395:ALA:O	1:B:399:ASN:ND2	2.46	0.43
1:A:379:LEU:CD1	1:A:381:LEU:HD21	2.47	0.43
1:B:158:THR:OG1	1:B:363:ARG:HG3	2.18	0.43
1:A:233:PHE:CD1	1:A:234:THR:N	2.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:ASN:HB3	1:A:272:TYR:HE1	1.84	0.43
3:A:443:MAN:C2	3:A:445:MAN:H1	2.46	0.43
1:A:228:ILE:HD13	1:A:422:ASN:O	2.19	0.43
3:A:445:MAN:C1	3:A:448:MAN:H2	2.48	0.43
1:B:241:PRO:HB2	1:B:386:CYS:SG	2.58	0.43
1:B:52:PHE:CZ	1:B:200:PRO:HD3	2.54	0.43
1:A:365:PRO:HB3	1:A:421:TYR:CE2	2.54	0.43
1:A:390:ALA:HB2	1:A:405:ALA:HB1	2.00	0.43
1:A:293:ALA:HB3	2:A:438:NAG:O6	2.19	0.43
1:B:172:GLU:HG3	1:B:172:GLU:H	1.61	0.43
1:B:106:ASP:OD1	1:B:107:LYS:N	2.51	0.43
1:B:182:GLN:NE2	1:B:188:TYR:CZ	2.87	0.43
1:A:295:VAL:CG1	1:A:295:VAL:O	2.67	0.43
3:A:447:MAN:C5	3:A:448:MAN:O1	2.67	0.43
1:A:406:ASN:HD22	1:A:407:THR:H	1.58	0.43
1:A:218:ARG:HG2	1:A:222:LEU:CD1	2.49	0.43
1:A:427:MET:HG2	1:A:428:SER:H	1.84	0.43
1:A:315:SER:HB2	1:A:328:THR:CG2	2.48	0.43
1:B:56:VAL:CG2	1:B:57:SER:N	2.82	0.43
1:B:248:SER:O	1:B:249:GLN:HB2	2.19	0.43
3:A:452:MAN:O6	3:A:452:MAN:O2	2.15	0.42
3:B:547:MAN:O4	3:B:548:MAN:H2	2.19	0.42
2:B:538:NAG:C3	3:B:540:MAN:C2	2.75	0.42
1:A:299:ASP:CG	1:A:302:ASN:HD22	2.22	0.42
4:B:550:NDG:O7	3:B:551:MAN:O2	2.36	0.42
1:A:392:SER:O	1:A:395:ALA:HB3	2.18	0.42
1:A:155:PHE:CE2	1:A:163:LEU:HD12	2.54	0.42
1:A:252:ARG:NE	1:A:376:ASN:ND2	2.67	0.42
1:A:61:SER:CB	1:A:63:ASN:HD21	2.32	0.42
1:A:211:LEU:HD22	1:A:215:GLU:OE1	2.20	0.42
1:A:52:PHE:CZ	1:A:200:PRO:HD3	2.54	0.42
1:A:315:SER:O	1:A:316:ALA:HB2	2.18	0.42
1:A:264:TRP:HA	1:A:415:ASN:O	2.19	0.42
1:B:35:THR:O	1:B:37:PHE:HD2	2.02	0.42
1:A:196:GLY:O	1:A:197:ALA:C	2.57	0.42
4:B:550:NDG:O7	3:B:551:MAN:H4	2.19	0.42
1:B:329:ARG:HH11	1:B:329:ARG:CB	2.32	0.42
1:B:232:GLN:HE22	1:B:256:ASN:N	2.17	0.42
1:B:319:GLN:HG3	1:B:323:GLN:C	2.40	0.42
1:B:33:ARG:HH11	1:B:33:ARG:HG3	1.84	0.42
2:A:438:NAG:HO4	3:A:441:MAN:C2	2.33	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:549:NAG:C4	3:B:553:MAN:HO1	2.30	0.42
1:A:406:ASN:C	1:A:406:ASN:HD22	2.16	0.42
1:B:154:PHE:CB	1:B:228:ILE:HB	2.50	0.42
1:A:172:GLU:H	1:A:172:GLU:HG3	1.59	0.42
1:B:290:THR:CG2	1:B:292:ALA:H	2.15	0.42
1:A:158:THR:OG1	1:A:363:ARG:HG3	2.18	0.42
2:A:438:NAG:O1	2:A:438:NAG:C7	2.68	0.42
1:B:291:ALA:HA	3:B:541:MAN:O1	2.18	0.42
1:A:163:LEU:HD23	1:A:163:LEU:HA	1.91	0.41
1:A:94:GLN:HB3	1:A:178:THR:O	2.20	0.41
1:B:262:LEU:HD12	1:B:318:ILE:CD1	2.50	0.41
1:A:221:GLN:NE2	1:A:221:GLN:CA	2.81	0.41
1:A:280:ASN:ND2	2:A:438:NAG:H1	2.10	0.41
1:A:187:ASN:N	1:A:187:ASN:HD22	2.10	0.41
1:A:287:GLY:O	1:A:290:THR:HG22	2.20	0.41
1:A:155:PHE:CD2	1:A:163:LEU:HD12	2.56	0.41
1:B:25:THR:HG22	1:B:27:PHE:N	2.34	0.41
2:A:438:NAG:C3	3:A:440:MAN:O2	2.69	0.41
1:A:302:ASN:HD21	4:A:446:NDG:C6	2.33	0.41
1:B:46:ILE:CG2	1:B:56:VAL:CG2	2.89	0.41
1:B:76:LEU:HB2	1:B:147:PHE:HE1	1.84	0.41
1:B:406:ASN:HD21	4:B:550:NDG:C1	2.34	0.41
1:A:133:THR:O	1:A:133:THR:CG2	2.69	0.41
1:A:146:ARG:NH1	1:A:146:ARG:CG	2.82	0.41
1:A:66:LEU:HD11	1:A:211:LEU:HD11	2.03	0.41
1:B:427:MET:HG2	1:B:428:SER:N	2.36	0.41
1:B:165:LEU:C	1:B:167:ALA:H	2.23	0.41
1:B:280:ASN:HA	1:B:285:CYS:SG	2.61	0.41
1:B:171:HIS:NE2	1:B:430:MET:HB3	2.36	0.41
1:B:405:ALA:C	1:B:407:THR:N	2.74	0.41
1:B:280:ASN:CG	2:B:538:NAG:O1	2.59	0.41
2:B:549:NAG:O3	3:B:554:MAN:O5	2.32	0.41
1:A:120:PHE:CD2	1:A:361:ALA:HB2	2.51	0.41
1:B:306:TYR:CD2	1:B:348:PRO:HA	2.56	0.41
1:A:103:GLN:HG2	6:A:466:HOH:O	2.21	0.41
1:B:33:ARG:NH1	1:B:33:ARG:HG3	2.36	0.41
1:A:64:GLY:HA3	1:A:208:TYR:HB3	2.03	0.41
1:A:395:ALA:CB	4:A:449:NDG:H8C2	2.51	0.41
1:B:290:THR:HG22	1:B:291:ALA:N	2.36	0.41
1:B:46:ILE:O	1:B:46:ILE:HG22	2.21	0.41
1:A:370:ASN:HD22	1:A:371:PHE:N	2.18	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:ILE:HG12	1:B:206:VAL:HG22	2.03	0.41
1:A:98:LEU:CD1	1:A:105:ILE:HB	2.49	0.40
1:B:158:THR:OG1	1:B:363:ARG:HD2	2.21	0.40
1:A:389:ASP:OD1	1:A:392:SER:N	2.54	0.40
2:A:438:NAG:HO3	3:A:440:MAN:C2	2.15	0.40
1:B:85:THR:HG23	1:B:88:PRO:CG	2.52	0.40
1:B:407:THR:O	1:B:407:THR:HG22	2.21	0.40
1:B:76:LEU:HD11	1:B:92:LEU:HD12	2.03	0.40
1:A:221:GLN:HE21	1:A:221:GLN:CA	2.34	0.40
1:B:315:SER:HB2	1:B:328:THR:CG2	2.52	0.40
1:B:356:PHE:HA	6:B:581:HOH:O	2.20	0.40
2:A:438:NAG:H3	3:A:440:MAN:O2	2.22	0.40
1:B:146:ARG:HH11	1:B:146:ARG:HG3	1.85	0.40
1:A:109:TYR:CD1	1:A:109:TYR:N	2.89	0.40
1:A:248:SER:O	1:A:249:GLN:HB2	2.21	0.40
1:A:280:ASN:ND2	1:A:393:PRO:CG	2.84	0.40
3:A:447:MAN:O4	3:A:448:MAN:H3	2.21	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	411/437 (94%)	361 (88%)	45 (11%)	5 (1%)	16	27
1	B	411/437 (94%)	371 (90%)	37 (9%)	3 (1%)	26	44
All	All	822/874 (94%)	732 (89%)	82 (10%)	8 (1%)	19	33

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	295	VAL

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Mol	Chain	Res	Type
1	A	300	TYR
1	B	245	THR
1	A	214	GLN
1	A	272	TYR
1	B	295	VAL
1	B	110	ASN
1	A	110	ASN

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	341/358 (95%)	315 (92%)	26 (8%)	16	29
1	B	341/358 (95%)	315 (92%)	26 (8%)	16	29
All	All	682/716 (95%)	630 (92%)	52 (8%)	16	29

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ARG
1	A	56	VAL
1	A	85	THR
1	A	92	LEU
1	A	93	LEU
1	A	96	VAL
1	A	104	ARG
1	A	139	GLU
1	A	146	ARG
1	A	172	GLU
1	A	187	ASN
1	A	201	THR
1	A	218	ARG
1	A	228	ILE
1	A	245	THR
1	A	265	ASN

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Mol	Chain	Res	Type
1	A	268	ASN
1	A	290	THR
1	A	329	ARG
1	A	364	GLN
1	A	370	ASN
1	A	379	LEU
1	A	399	ASN
1	A	406	ASN
1	A	414	LEU
1	A	424	LEU
1	B	32	ARG
1	B	56	VAL
1	B	85	THR
1	B	92	LEU
1	B	93	LEU
1	B	95	ASP
1	B	104	ARG
1	B	163	LEU
1	B	172	GLU
1	B	201	THR
1	B	227	LEU
1	B	228	ILE
1	B	248	SER
1	B	265	ASN
1	B	267	ASN
1	B	268	ASN
1	B	282	PRO
1	B	329	ARG
1	B	346	VAL
1	B	353	LEU
1	B	364	GLN
1	B	369	CYS
1	B	370	ASN
1	B	379	LEU
1	B	386	CYS
1	B	424	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	44	GLN

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Mol	Chain	Res	Type
1	A	63	ASN
1	A	103	GLN
1	A	110	ASN
1	A	128	ASN
1	A	156	ASN
1	A	157	GLN
1	A	171	HIS
1	A	182	GLN
1	A	187	ASN
1	A	221	GLN
1	A	232	GLN
1	A	246	GLN
1	A	265	ASN
1	A	268	ASN
1	A	280	ASN
1	A	302	ASN
1	A	319	GLN
1	A	364	GLN
1	A	370	ASN
1	A	376	ASN
1	A	399	ASN
1	A	406	ASN
1	A	415	ASN
1	B	43	GLN
1	B	44	GLN
1	B	54	ASN
1	B	63	ASN
1	B	108	HIS
1	B	128	ASN
1	B	156	ASN
1	B	171	HIS
1	B	182	GLN
1	B	187	ASN
1	B	232	GLN
1	B	265	ASN
1	B	268	ASN
1	B	302	ASN
1	B	364	GLN
1	B	370	ASN
1	B	376	ASN
1	B	399	ASN
1	B	406	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 38 ligands modelled in this entry, 4 are monoatomic - leaving 34 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	438	-	15,15,15	0.37	0	17,21,21	1.15	2 (11%)
3	MAN	A	439	-	12,12,12	0.54	0	17,17,17	0.57	0
3	MAN	A	440	-	12,12,12	0.67	0	17,17,17	1.35	3 (17%)
3	MAN	A	441	-	12,12,12	0.52	0	17,17,17	1.09	2 (11%)
3	MAN	A	442	-	12,12,12	0.36	0	17,17,17	0.90	1 (5%)
3	MAN	A	443	-	12,12,12	0.56	0	17,17,17	1.07	2 (11%)
2	NAG	A	444	-	15,15,15	0.73	0	17,21,21	0.88	0
3	MAN	A	445	-	12,12,12	0.49	0	17,17,17	0.56	0
4	NDG	A	446	-	15,15,15	0.76	0	17,21,21	1.43	2 (11%)
3	MAN	A	447	-	12,12,12	0.58	0	17,17,17	0.93	1 (5%)
3	MAN	A	448	-	12,12,12	0.54	0	17,17,17	0.79	0
4	NDG	A	449	-	15,15,15	0.53	0	17,21,21	0.82	1 (5%)
2	NAG	A	450	-	15,15,15	0.70	0	17,21,21	0.71	0
3	MAN	A	451	-	12,12,12	0.41	0	17,17,17	0.66	0
3	MAN	A	452	-	12,12,12	0.43	0	17,17,17	0.66	0
3	MAN	A	453	-	12,12,12	0.68	0	17,17,17	0.84	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	MAN	A	454	-	12,12,12	0.51	0	17,17,17	0.59	0
2	NAG	B	538	-	15,15,15	0.61	0	17,21,21	1.12	2 (11%)
3	MAN	B	539	-	12,12,12	0.62	0	17,17,17	0.94	1 (5%)
3	MAN	B	540	-	12,12,12	0.33	0	17,17,17	0.59	0
3	MAN	B	541	-	12,12,12	0.60	0	17,17,17	0.97	1 (5%)
3	MAN	B	542	-	12,12,12	0.48	0	17,17,17	0.66	0
3	MAN	B	543	-	12,12,12	0.47	0	17,17,17	0.78	0
4	NDG	B	544	-	15,15,15	0.45	0	17,21,21	0.76	0
3	MAN	B	545	-	12,12,12	0.67	0	17,17,17	0.45	0
2	NAG	B	546	-	15,15,15	0.74	0	17,21,21	0.76	0
3	MAN	B	547	-	12,12,12	0.57	0	17,17,17	0.48	0
3	MAN	B	548	-	12,12,12	0.41	0	17,17,17	0.53	0
2	NAG	B	549	-	15,15,15	0.44	0	17,21,21	0.71	0
4	NDG	B	550	-	15,15,15	0.60	0	17,21,21	0.50	0
3	MAN	B	551	-	12,12,12	0.56	0	17,17,17	0.61	0
3	MAN	B	552	-	12,12,12	0.69	0	17,17,17	0.65	0
3	MAN	B	553	-	12,12,12	0.51	0	17,17,17	0.66	0
3	MAN	B	554	-	12,12,12	0.61	0	17,17,17	0.99	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
2	NAG	A	438	-	-	0/6/26/26	0/1/1/1
3	MAN	A	439	-	1/1/5/5	0/2/22/22	0/1/1/1
3	MAN	A	440	-	-	0/2/22/22	0/1/1/1
3	MAN	A	441	-	-	0/2/22/22	0/1/1/1
3	MAN	A	442	-	1/1/5/5	0/2/22/22	0/1/1/1
3	MAN	A	443	-	-	0/2/22/22	0/1/1/1
2	NAG	A	444	-	-	0/6/26/26	0/1/1/1
3	MAN	A	445	-	-	0/2/22/22	0/1/1/1
4	NDG	A	446	-	-	0/6/26/26	0/1/1/1
3	MAN	A	447	-	1/1/5/5	0/2/22/22	0/1/1/1
3	MAN	A	448	-	-	0/2/22/22	0/1/1/1
4	NDG	A	449	-	-	0/6/26/26	0/1/1/1
2	NAG	A	450	-	-	0/6/26/26	0/1/1/1
3	MAN	A	451	-	1/1/5/5	0/2/22/22	0/1/1/1
3	MAN	A	452	-	-	0/2/22/22	0/1/1/1
3	MAN	A	453	-	-	0/2/22/22	0/1/1/1
3	MAN	A	454	-	1/1/5/5	0/2/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	538	-	-	0/6/26/26	0/1/1/1
3	MAN	B	539	-	-	0/2/22/22	0/1/1/1
3	MAN	B	540	-	-	0/2/22/22	0/1/1/1
3	MAN	B	541	-	-	0/2/22/22	0/1/1/1
3	MAN	B	542	-	-	0/2/22/22	0/1/1/1
3	MAN	B	543	-	1/1/5/5	0/2/22/22	0/1/1/1
4	NDG	B	544	-	-	0/6/26/26	0/1/1/1
3	MAN	B	545	-	-	0/2/22/22	0/1/1/1
2	NAG	B	546	-	-	0/6/26/26	0/1/1/1
3	MAN	B	547	-	1/1/5/5	0/2/22/22	0/1/1/1
3	MAN	B	548	-	1/1/5/5	0/2/22/22	0/1/1/1
2	NAG	B	549	-	-	0/6/26/26	0/1/1/1
4	NDG	B	550	-	-	0/6/26/26	0/1/1/1
3	MAN	B	551	-	1/1/5/5	0/2/22/22	0/1/1/1
3	MAN	B	552	-	-	0/2/22/22	0/1/1/1
3	MAN	B	553	-	-	0/2/22/22	0/1/1/1
3	MAN	B	554	-	1/1/5/5	0/2/22/22	0/1/1/1

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	446	NDG	C4-C3-C2	-4.41	104.31	110.43
3	A	441	MAN	C4-C3-C2	-2.81	105.55	110.79
3	B	539	MAN	C3-C4-C5	-2.59	105.68	110.20
2	B	538	NAG	C4-C3-C2	-2.58	106.85	110.43
3	A	447	MAN	C4-C3-C2	-2.51	106.10	110.79
4	A	446	NDG	C3-C4-C5	-2.48	105.87	110.20
2	B	538	NAG	C6-C5-C4	-2.46	106.96	113.02
3	A	441	MAN	C3-C4-C5	-2.26	106.26	110.20
4	A	449	NDG	C4-C3-C2	-2.24	107.32	110.43
3	B	541	MAN	C3-C4-C5	-2.13	106.49	110.20
3	A	440	MAN	C3-C4-C5	-2.01	106.69	110.20
2	A	438	NAG	C3-C4-C5	2.03	113.74	110.20
3	A	442	MAN	C3-C4-C5	2.24	114.09	110.20
3	A	443	MAN	C4-C3-C2	2.44	115.35	110.79
3	A	440	MAN	O5-C1-C2	2.54	113.85	109.80
3	B	554	MAN	C1-C2-C3	2.73	114.50	110.43
2	A	438	NAG	C4-C3-C2	2.74	114.22	110.43
3	A	443	MAN	C3-C4-C5	2.88	115.21	110.20
3	A	440	MAN	C1-C2-C3	4.18	116.64	110.43

All (10) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	554	MAN	C1
3	A	451	MAN	C1
3	A	447	MAN	C1
3	A	439	MAN	C1
3	B	548	MAN	C1
3	A	442	MAN	C1
3	A	454	MAN	C1
3	B	543	MAN	C1
3	B	551	MAN	C1
3	B	547	MAN	C1

There are no torsion outliers.

There are no ring outliers.

34 monomers are involved in 196 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	438	NAG	30	0
3	A	439	MAN	2	0
3	A	440	MAN	7	0
3	A	441	MAN	11	0
3	A	442	MAN	1	0
3	A	443	MAN	15	0
2	A	444	NAG	12	0
3	A	445	MAN	6	0
4	A	446	NDG	12	0
3	A	447	MAN	13	0
3	A	448	MAN	10	0
4	A	449	NDG	18	0
2	A	450	NAG	3	0
3	A	451	MAN	1	0
3	A	452	MAN	9	0
3	A	453	MAN	2	0
3	A	454	MAN	2	0
2	B	538	NAG	30	0
3	B	539	MAN	2	0
3	B	540	MAN	12	0
3	B	541	MAN	16	0
3	B	542	MAN	6	0
3	B	543	MAN	12	0
4	B	544	NDG	7	0
3	B	545	MAN	5	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	546	NAG	5	0
3	B	547	MAN	11	0
3	B	548	MAN	6	0
2	B	549	NAG	20	0
4	B	550	NDG	10	0
3	B	551	MAN	3	0
3	B	552	MAN	3	0
3	B	553	MAN	7	0
3	B	554	MAN	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	413/437 (94%)	0.41	19 (4%) 36 42	22, 34, 48, 52	0
1	B	413/437 (94%)	0.46	15 (3%) 46 53	21, 34, 48, 55	0
All	All	826/874 (94%)	0.43	34 (4%) 41 47	21, 34, 48, 55	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	194	ILE	3.2
1	A	410	LEU	3.2
1	B	244	THR	3.2
1	A	295	VAL	3.1
1	A	320	LEU	2.9
1	B	292	ALA	2.9
1	A	276	THR	2.7
1	B	283	GLY	2.7
1	B	312	VAL	2.7
1	A	195	ALA	2.6
1	B	296	THR	2.6
1	A	186	VAL	2.5
1	A	437	ASN	2.5
1	A	425	ARG	2.5
1	A	279	ALA	2.4
1	A	47	ASN	2.3
1	B	372	SER	2.3
1	B	405	ALA	2.3
1	B	305	THR	2.3
1	B	388	ILE	2.3
1	A	254	ASN	2.3
1	A	199	GLN	2.2
1	B	387	SER	2.2
1	A	284	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	436	ALA	2.2
1	A	143	ALA	2.1
1	A	280	ASN	2.1
1	A	292	ALA	2.1
1	A	408	ALA	2.1
1	A	323	GLN	2.1
1	B	140	LEU	2.1
1	B	49	SER	2.1
1	B	295	VAL	2.0
1	B	245	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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(Å ²)	Q<0.9
4	NDG	A	446	15/15	0.75	0.31	9.72	55,56,59,59	0
4	NDG	A	449	15/15	0.74	0.32	3.91	45,48,51,51	0
2	NAG	A	438	15/15	0.87	0.35	3.73	56,56,58,59	0
2	NAG	B	538	15/15	0.72	0.34	3.13	55,57,61,63	0
2	NAG	B	549	15/15	0.77	0.35	3.11	43,49,51,53	0
2	NAG	A	444	15/15	0.67	0.20	1.68	38,43,46,48	0
3	MAN	A	443	12/12	0.83	0.25	1.34	51,51,52,53	0
3	MAN	B	553	12/12	0.84	0.24	1.16	37,40,43,43	0
3	MAN	A	453	12/12	0.87	0.23	0.73	39,41,42,42	0
4	NDG	B	550	15/15	0.75	0.27	0.65	53,55,56,56	0
2	NAG	A	450	15/15	0.81	0.25	0.39	55,55,56,56	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	NDG	B	544	15/15	0.71	0.20	-0.02	38,44,45,46	0
5	HG	A	455	1/1	0.99	0.10	-1.43	61,61,61,61	1
5	HG	B	457	1/1	0.99	0.10	-1.63	69,69,69,69	1
5	HG	B	458	1/1	0.92	0.10	-2.76	86,86,86,86	1
5	HG	A	456	1/1	1.00	0.11	-4.38	71,71,71,71	0
3	MAN	A	452	12/12	0.73	0.31	-	50,53,55,55	0
3	MAN	A	448	12/12	0.87	0.16	-	53,54,56,57	0
3	MAN	B	542	12/12	0.57	0.28	-	64,66,67,69	0
3	MAN	A	447	12/12	0.79	0.23	-	58,58,59,59	0
3	MAN	B	548	12/12	0.81	0.28	-	55,56,58,60	0
3	MAN	A	442	12/12	0.72	0.39	-	64,65,66,66	0
3	MAN	A	439	12/12	0.85	0.30	-	52,53,54,56	0
2	NAG	B	546	15/15	0.67	0.30	-	50,55,60,60	0
3	MAN	B	539	12/12	0.79	0.27	-	52,53,54,55	0
3	MAN	A	440	12/12	0.84	0.32	-	54,55,58,59	0
3	MAN	A	454	12/12	0.88	0.27	-	48,50,53,54	0
3	MAN	B	541	12/12	0.71	0.27	-	56,57,59,60	0
3	MAN	B	540	12/12	0.67	0.38	-	54,55,58,58	0
3	MAN	B	543	12/12	0.83	0.27	-	50,50,52,52	0
3	MAN	B	552	12/12	0.49	0.28	-	49,53,54,54	0
3	MAN	B	551	12/12	0.86	0.26	-	59,61,62,63	0
3	MAN	B	545	12/12	0.44	0.47	-	64,66,67,67	0
3	MAN	A	451	12/12	0.79	0.19	-	60,62,63,63	0
3	MAN	A	445	12/12	0.75	0.41	-	64,66,67,69	0
3	MAN	B	547	12/12	0.73	0.34	-	58,59,61,62	0
3	MAN	B	554	12/12	0.82	0.37	-	50,53,54,55	0
3	MAN	A	441	12/12	0.71	0.25	-	54,57,58,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.