



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 11:11 AM GMT

PDB ID : 2B4X
Title : Crystal Structure of Antithrombin-III
Authors : Adams, T.E.; Huntington, J.A.; Johnson, D.J.D.; Bock, S.C.
Deposited on : 2005-09-27
Resolution : 2.80 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

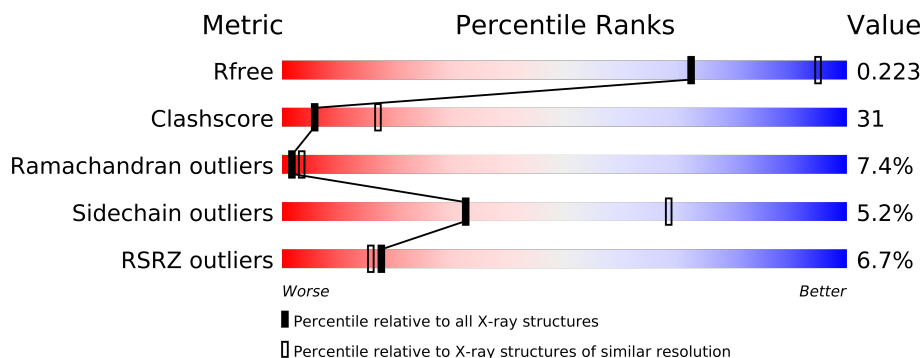
The following versions of software and data (see [references](#)) were used in the production of this report:

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

1 Overall quality at a glance

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	1799 (2.80-2.80)
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (2.80-2.80)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	I	427	
2	L	427	

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

Mol	Type	Chain	Res	Geometry	Electron density
3	NAG	I	801	-	X
3	NAG	I	861	-	X
3	NAG	L	801	-	X
3	NAG	L	842	-	X

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 6560 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Antithrombin-III.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	I	413	Total	C	N	O	S	0	0	0
			3203	2044	528	615	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	135	ALA	ASN	ENGINEERED	UNP P01008
I	220	LEU	TYR	ENGINEERED	UNP P01008

- Molecule 2 is a protein called Antithrombin-III.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	410	Total	C	N	O	S	0	0	0
			3078	1969	497	595	17			

- Molecule 3 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	I	1	Total C N O 14 8 1 5	0	0
3	I	1	Total C N O 14 8 1 5	0	0
3	I	1	Total C N O 14 8 1 5	0	0
3	L	1	Total C N O 14 8 1 5	0	0
3	L	1	Total C N O 14 8 1 5	0	0
3	L	1	Total C N O 14 8 1 5	0	0

- Molecule 4 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	I	2	Total C N O 28 16 2 10	0	0
4	L	2	Total C N O 27 16 2 9	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	135	ALA	ASN	ENGINEERED	UNP P01008
I	220	LEU	TYR	ENGINEERED	UNP P01008

- Molecule 5 is water.

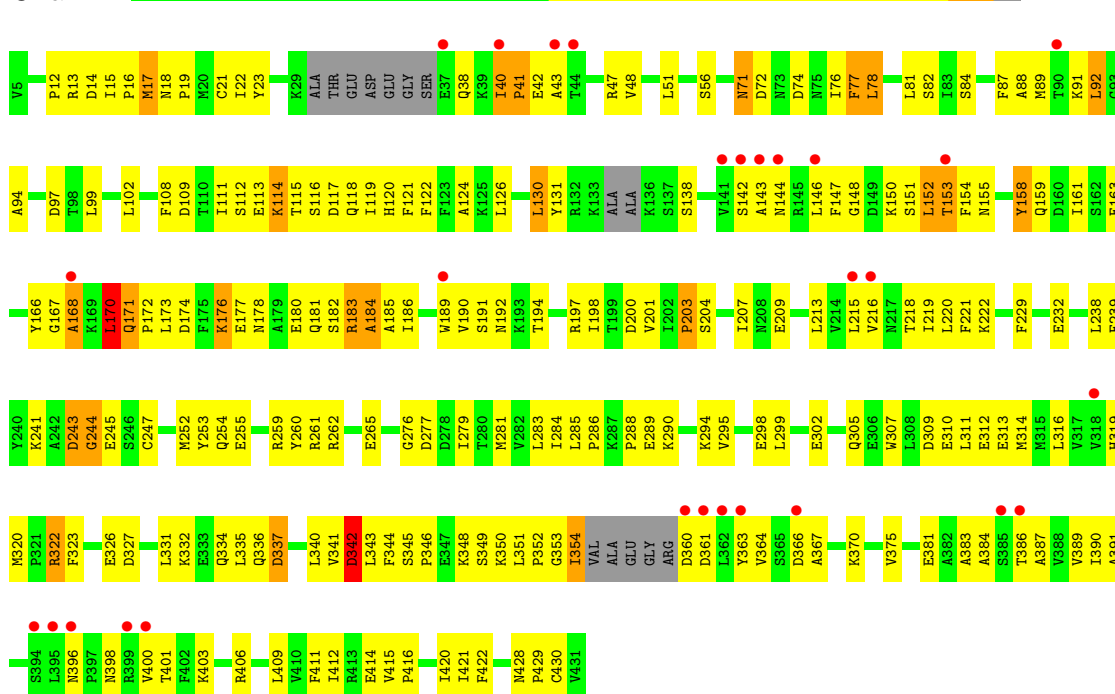
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	I	84	Total O 84 84	0	0
5	L	56	Total O 56 56	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

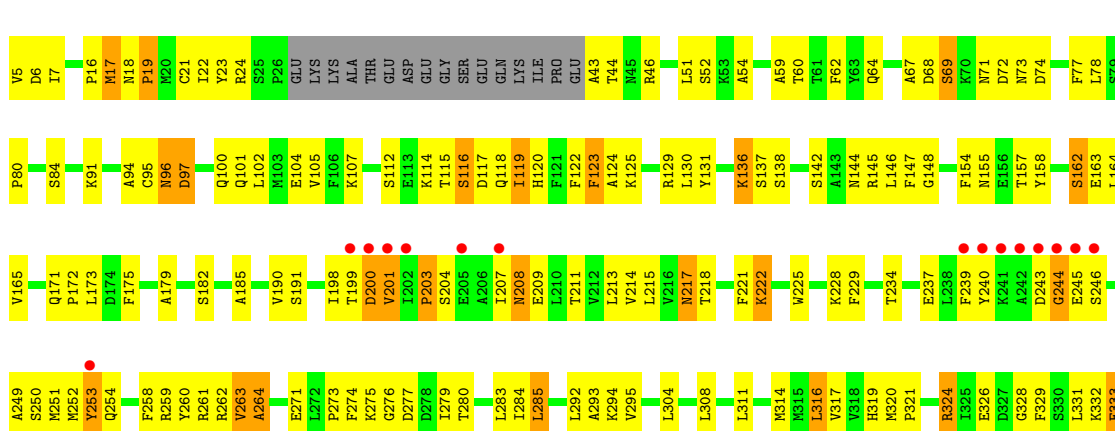
• Molecule 1: Antithrombin-III

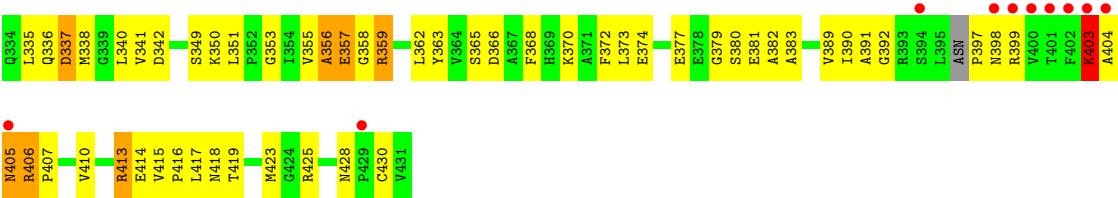
Chain I:



• Molecule 2: Antithrombin-III

Chain L:





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	68.47Å 98.67Å 89.24Å 90.00° 105.00° 90.00°	Depositor
Resolution (Å)	19.98 – 2.80 54.94 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (19.98-2.80) 99.9 (54.94-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 2.81Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.228 , 0.283 0.231 , 0.223	Depositor DCC
R_{free} test set	1438 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	79.5	Xtriage
Anisotropy	0.455	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 53.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 28330 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6560	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.17% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	I	0.40	0/3265	0.67	0/4423
2	L	0.40	0/3141	0.66	0/4277
All	All	0.40	0/6406	0.66	0/8700

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
4	I	1	0

There are no bond length outliers.

There are no bond angle outliers.

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	I	841	NAG	C1

There are no planarity outliers.

5.2 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	I	3203	0	3086	199	0
2	L	3078	0	2857	196	0
3	I	42	0	39	9	0
3	L	42	0	39	5	0
4	I	28	0	25	1	0
4	L	27	0	25	0	0
5	I	84	0	0	7	0
5	L	56	0	0	3	0
All	All	6560	0	6071	393	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 31.

All (393) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:150:LYS:HA	1:I:172:PRO:HB2	1.42	0.99
2:L:404:ALA:HB2	2:L:428:ASN:ND2	1.78	0.98
2:L:7:ILE:H	2:L:7:ILE:HD12	1.30	0.96
2:L:404:ALA:HB2	2:L:428:ASN:HD22	1.32	0.94
2:L:94:ALA:HA	2:L:351:LEU:HD23	1.52	0.91
1:I:190:VAL:HB	1:I:201:VAL:HG21	1.51	0.91
2:L:292:LEU:HD23	2:L:407:PRO:HG2	1.54	0.89
1:I:143:ALA:HB3	1:I:218:THR:HG23	1.51	0.89
2:L:124:ALA:HB2	2:L:165:VAL:HG13	1.58	0.84
2:L:211:THR:HA	2:L:391:ALA:O	1.78	0.83
2:L:365:SER:HB3	2:L:392:GLY:H	1.46	0.81
1:I:319:HIS:HB2	1:I:403:LYS:HA	1.64	0.79
1:I:415:VAL:HB	1:I:416:PRO:HD3	1.64	0.79
2:L:215:LEU:HD12	2:L:215:LEU:H	1.48	0.79
2:L:285:LEU:HD23	2:L:285:LEU:N	1.99	0.78
1:I:232:GLU:OE1	2:L:262:ARG:HD2	1.84	0.78
2:L:73:ASN:HB3	2:L:404:ALA:O	1.84	0.77
2:L:239:PHE:HE1	2:L:404:ALA:HB1	1.50	0.77
2:L:71:ASN:HB2	5:L:652:HOH:O	1.82	0.77
1:I:192:ASN:HD21	3:I:861:NAG:H2	1.49	0.77
1:I:183:ARG:HB2	1:I:207:ILE:HD12	1.66	0.76
2:L:145:ARG:HG2	2:L:147:PHE:CE1	2.22	0.74
1:I:176:LYS:HA	1:I:209:GLU:HG3	1.69	0.74
2:L:215:LEU:HD12	2:L:215:LEU:N	2.04	0.73
2:L:62:PHE:HA	2:L:338:MET:HE1	1.70	0.73
2:L:213:LEU:HB3	2:L:390:ILE:HD12	1.71	0.73
2:L:244:GLY:HA2	5:L:537:HOH:O	1.90	0.71
2:L:91:LYS:NZ	2:L:120:HIS:NE2	2.37	0.71

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:261:ARG:HA	1:I:261:ARG:NH1	2.05	0.71
1:I:192:ASN:HD21	3:I:861:NAG:C2	2.03	0.71
1:I:161:ILE:C	1:I:163:GLU:H	1.92	0.70
1:I:154:PHE:HA	1:I:354:ILE:HG23	1.74	0.70
1:I:213:LEU:HD23	1:I:364:VAL:HG22	1.74	0.70
2:L:252:MET:HE2	2:L:320:MET:HG2	1.73	0.69
1:I:354:ILE:HD12	1:I:354:ILE:N	2.08	0.69
2:L:190:VAL:HG21	2:L:201:VAL:HG21	1.73	0.69
1:I:192:ASN:HD21	3:I:861:NAG:C1	2.06	0.68
2:L:59:ALA:HB1	2:L:423:MET:CE	2.23	0.68
2:L:91:LYS:HE2	2:L:120:HIS:NE2	2.07	0.68
1:I:183:ARG:HE	1:I:204:SER:HA	1.59	0.68
1:I:336:GLN:HA	1:I:340:LEU:O	1.93	0.68
1:I:262:ARG:HG3	1:I:262:ARG:HH11	1.59	0.67
2:L:405:ASN:CG	2:L:406:ARG:H	1.97	0.66
1:I:332:LYS:O	1:I:336:GLN:HG3	1.95	0.66
2:L:7:ILE:HG12	2:L:164:LEU:O	1.96	0.66
2:L:77:PHE:CZ	2:L:373:LEU:HB2	2.30	0.66
1:I:91:LYS:NZ	1:I:120:HIS:NE2	2.38	0.66
3:L:841:NAG:O4	3:L:842:NAG:C1	2.44	0.66
1:I:146:LEU:HD13	1:I:215:LEU:HD13	1.78	0.66
2:L:77:PHE:CE2	2:L:373:LEU:HB2	2.32	0.65
2:L:324:ARG:HG3	2:L:324:ARG:O	1.96	0.65
1:I:259:ARG:NH2	1:I:311:LEU:HB2	2.11	0.65
2:L:292:LEU:CD2	2:L:407:PRO:HG2	2.26	0.65
2:L:355:VAL:HG23	2:L:362:LEU:HD11	1.77	0.65
1:I:143:ALA:HB3	1:I:218:THR:CG2	2.23	0.64
2:L:91:LYS:CE	2:L:120:HIS:NE2	2.60	0.64
2:L:59:ALA:HB1	2:L:423:MET:HE3	1.78	0.64
2:L:341:VAL:HG23	2:L:342:ASP:H	1.61	0.64
2:L:413:ARG:NH2	2:L:415:VAL:HG13	2.12	0.64
1:I:71:ASN:HD22	1:I:72:ASP:H	1.45	0.64
2:L:182:SER:O	2:L:185:ALA:HB3	1.97	0.64
1:I:91:LYS:HB2	1:I:102:LEU:HD13	1.79	0.64
2:L:237:GLU:HB2	5:L:656:HOH:O	1.97	0.64
1:I:294:LYS:O	1:I:298:GLU:HG3	1.96	0.64
2:L:365:SER:CB	2:L:392:GLY:H	2.12	0.63
2:L:239:PHE:CE1	2:L:404:ALA:HB1	2.32	0.63
2:L:261:ARG:HG3	2:L:263:VAL:CG1	2.29	0.63
1:I:182:SER:O	1:I:184:ALA:N	2.25	0.63
1:I:174:ASP:HB3	1:I:178:ASN:HB2	1.79	0.63
1:I:312:GLU:HB2	5:I:588:HOH:O	1.99	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:71:ASN:ND2	2:L:73:ASN:HB2	2.14	0.62
1:I:183:ARG:HG2	1:I:183:ARG:O	2.00	0.62
2:L:341:VAL:HG23	2:L:342:ASP:N	2.14	0.62
1:I:78:LEU:HD23	1:I:78:LEU:N	2.15	0.62
1:I:190:VAL:HG21	1:I:201:VAL:HG11	1.80	0.62
1:I:343:LEU:HD11	1:I:364:VAL:HG23	1.83	0.61
1:I:144:ASN:HB2	5:I:519:HOH:O	1.99	0.61
2:L:125:LYS:O	2:L:129:ARG:HG2	2.00	0.61
2:L:331:LEU:O	2:L:335:LEU:HB2	1.99	0.61
2:L:229:PHE:HB2	2:L:377:GLU:HA	1.82	0.61
2:L:175:PHE:O	2:L:209:GLU:HA	2.01	0.61
1:I:331:LEU:HB2	1:I:367:ALA:HB3	1.82	0.61
2:L:191:SER:OG	2:L:199:THR:HG22	2.01	0.61
1:I:326:GLU:HG2	1:I:370:LYS:HE3	1.82	0.60
1:I:190:VAL:CB	1:I:201:VAL:HG21	2.27	0.60
2:L:222:LYS:HG2	2:L:381:GLU:HG3	1.82	0.60
1:I:51:LEU:HA	1:I:111:ILE:HD11	1.83	0.60
1:I:170:LEU:O	1:I:171:GLN:HB3	2.02	0.60
2:L:261:ARG:O	2:L:263:VAL:HG13	2.02	0.60
1:I:194:THR:O	1:I:197:ARG:HG2	2.02	0.60
2:L:91:LYS:HZ3	2:L:120:HIS:CD2	2.19	0.59
2:L:274:PHE:HD2	2:L:279:ILE:O	1.84	0.59
1:I:283:LEU:HD11	1:I:320:MET:HE1	1.84	0.59
2:L:119:ILE:HG22	2:L:120:HIS:N	2.16	0.59
1:I:18:ASN:O	1:I:21:CYS:HB2	2.03	0.59
5:I:609:HOH:O	2:L:249:ALA:HB2	2.02	0.59
2:L:102:LEU:HD23	2:L:340:LEU:HD11	1.84	0.59
1:I:159:GLN:CB	1:I:170:LEU:HD21	2.32	0.59
1:I:111:ILE:O	1:I:114:LYS:HB2	2.03	0.58
1:I:182:SER:O	1:I:185:ALA:N	2.36	0.58
1:I:108:PHE:HB3	1:I:119:ILE:HG12	1.85	0.58
1:I:183:ARG:HG2	1:I:183:ARG:HH11	1.67	0.58
2:L:221:PHE:CE1	2:L:279:ILE:HG21	2.39	0.58
2:L:144:ASN:HD22	2:L:217:ASN:HA	1.68	0.58
2:L:252:MET:CE	2:L:320:MET:HG2	2.34	0.58
2:L:428:ASN:OD1	2:L:430:CYS:HB2	2.03	0.58
1:I:183:ARG:CB	1:I:207:ILE:HD12	2.34	0.58
1:I:391:ALA:O	2:L:321:PRO:HD3	2.02	0.58
2:L:7:ILE:H	2:L:7:ILE:CD1	2.07	0.58
2:L:24:ARG:HA	2:L:114:LYS:O	2.04	0.58
2:L:54:ALA:HB1	2:L:107:LYS:O	2.03	0.58
1:I:289:GLU:CD	1:I:289:GLU:H	2.08	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:292:LEU:HD21	2:L:407:PRO:O	2.05	0.57
1:I:47:ARG:HB3	1:I:122:PHE:CD1	2.39	0.57
2:L:254:GLN:NE2	2:L:258:PHE:HZ	2.02	0.57
1:I:115:THR:H	1:I:118:GLN:NE2	2.03	0.57
1:I:286:PRO:HB3	1:I:295:VAL:HG21	1.86	0.57
2:L:207:ILE:HG23	2:L:211:THR:HG21	1.85	0.56
2:L:208:ASN:C	2:L:208:ASN:HD22	2.08	0.56
1:I:115:THR:H	1:I:118:GLN:HE21	1.51	0.56
1:I:161:ILE:C	1:I:163:GLU:N	2.59	0.56
1:I:331:LEU:HB3	1:I:335:LEU:HD11	1.87	0.56
2:L:52:SER:HB2	2:L:419:THR:HG22	1.87	0.56
1:I:260:TYR:CD2	1:I:261:ARG:N	2.73	0.56
2:L:261:ARG:HG3	2:L:263:VAL:HG13	1.87	0.56
2:L:131:TYR:CE1	2:L:142:SER:HB2	2.39	0.56
1:I:396:ASN:OD1	1:I:398:ASN:N	2.36	0.56
1:I:71:ASN:HD22	1:I:72:ASP:N	2.03	0.56
1:I:387:ALA:O	2:L:316:LEU:HB2	2.06	0.56
1:I:284:ILE:HD13	1:I:307:TRP:HZ3	1.70	0.56
1:I:183:ARG:HG2	1:I:183:ARG:NH1	2.21	0.56
2:L:116:SER:O	2:L:118:GLN:N	2.40	0.55
1:I:285:LEU:HD11	1:I:406:ARG:HG3	1.87	0.55
2:L:7:ILE:HD12	2:L:7:ILE:N	2.11	0.55
1:I:51:LEU:HD22	1:I:122:PHE:CB	2.36	0.55
2:L:222:LYS:CG	2:L:381:GLU:HG3	2.37	0.55
2:L:96:ASN:ND2	3:L:801:NAG:C1	2.70	0.55
2:L:366:ASP:OD2	2:L:368:PHE:CZ	2.60	0.55
1:I:192:ASN:ND2	3:I:861:NAG:H2	2.18	0.55
2:L:332:LYS:O	2:L:336:GLN:HG3	2.07	0.55
1:I:191:SER:HA	1:I:198:ILE:O	2.07	0.54
1:I:342:ASP:O	1:I:349:SER:HA	2.06	0.54
2:L:148:GLY:HA3	2:L:154:PHE:CE2	2.42	0.54
2:L:198:ILE:HG23	2:L:370:LYS:HG2	1.88	0.54
1:I:177:GLU:C	1:I:178:ASN:HD22	2.11	0.54
1:I:17:MET:HG2	1:I:117:ASP:HB3	1.88	0.54
1:I:76:ILE:O	1:I:77:PHE:HB2	2.08	0.54
1:I:412:ILE:O	1:I:421:ILE:HB	2.08	0.54
1:I:221:PHE:CE1	1:I:279:ILE:HG21	2.43	0.54
1:I:262:ARG:HG3	1:I:262:ARG:NH1	2.23	0.54
2:L:234:THR:HG23	2:L:251:MET:O	2.07	0.54
2:L:115:THR:HG23	2:L:118:GLN:CB	2.38	0.54
2:L:328:GLY:O	2:L:329:PHE:HB3	2.09	0.53
2:L:372:PHE:O	2:L:382:ALA:HA	2.08	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:405:ASN:ND2	2:L:406:ARG:N	2.55	0.53
1:I:389:VAL:HA	5:I:655:HOH:O	2.07	0.53
2:L:199:THR:O	2:L:200:ASP:HB3	2.08	0.53
1:I:238:LEU:HA	1:I:247:CYS:O	2.08	0.53
1:I:331:LEU:HB3	1:I:335:LEU:CD1	2.39	0.53
2:L:403:LYS:HD3	2:L:404:ALA:N	2.24	0.53
2:L:18:ASN:O	2:L:19:PRO:O	2.26	0.53
1:I:183:ARG:HB2	1:I:207:ILE:CD1	2.37	0.53
1:I:183:ARG:CA	1:I:207:ILE:HD12	2.39	0.53
2:L:221:PHE:CG	2:L:222:LYS:N	2.77	0.53
2:L:105:VAL:HG21	2:L:340:LEU:HB2	1.91	0.52
1:I:345:SER:HB3	1:I:348:LYS:HB2	1.91	0.52
1:I:232:GLU:CD	2:L:262:ARG:HD2	2.30	0.52
1:I:260:TYR:CG	1:I:261:ARG:N	2.78	0.52
2:L:283:LEU:HD11	2:L:320:MET:CE	2.39	0.52
1:I:182:SER:C	1:I:184:ALA:H	2.12	0.52
2:L:71:ASN:C	2:L:73:ASN:H	2.13	0.52
2:L:71:ASN:HD21	2:L:73:ASN:HB2	1.74	0.52
2:L:22:ILE:N	2:L:22:ILE:HD12	2.24	0.52
2:L:130:LEU:HD23	2:L:417:LEU:HD13	1.92	0.52
2:L:283:LEU:O	2:L:284:ILE:HG13	2.08	0.52
1:I:213:LEU:HB3	1:I:364:VAL:HA	1.92	0.52
2:L:259:ARG:HD3	2:L:271:GLU:OE1	2.09	0.52
1:I:261:ARG:HD3	1:I:262:ARG:N	2.25	0.51
1:I:302:GLU:CD	1:I:302:GLU:H	2.13	0.51
2:L:91:LYS:HE2	2:L:120:HIS:CE1	2.45	0.51
1:I:89:MET:HA	1:I:166:TYR:CD2	2.45	0.51
1:I:170:LEU:HD23	1:I:170:LEU:H	1.76	0.51
2:L:67:ALA:O	2:L:69:SER:N	2.40	0.51
1:I:155:ASN:HB3	1:I:353:GLY:O	2.11	0.51
2:L:404:ALA:CB	2:L:428:ASN:HB2	2.41	0.51
2:L:179:ALA:HB1	2:L:207:ILE:O	2.11	0.51
2:L:74:ASP:O	2:L:425:ARG:HD2	2.10	0.51
1:I:153:THR:HG22	1:I:354:ILE:C	2.32	0.50
1:I:154:PHE:HA	1:I:354:ILE:CG2	2.40	0.50
1:I:15:ILE:N	1:I:16:PRO:HD3	2.26	0.50
2:L:213:LEU:HG	2:L:214:VAL:N	2.25	0.50
3:I:862:NAG:H3	3:I:862:NAG:O7	2.09	0.50
2:L:349:SER:OG	2:L:363:TYR:HA	2.11	0.50
1:I:19:PRO:C	1:I:21:CYS:H	2.13	0.50
1:I:148:GLY:O	1:I:172:PRO:HA	2.11	0.50
2:L:60:THR:O	2:L:64:GLN:HG3	2.12	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:304:LEU:O	2:L:308:LEU:HG	2.12	0.50
1:I:71:ASN:ND2	1:I:72:ASP:N	2.58	0.50
2:L:78:LEU:N	2:L:78:LEU:HD23	2.27	0.50
1:I:241:LYS:HD3	1:I:428:ASN:HD22	1.77	0.50
2:L:405:ASN:ND2	2:L:406:ARG:H	2.10	0.49
1:I:84:SER:O	1:I:88:ALA:HB2	2.12	0.49
2:L:292:LEU:O	2:L:295:VAL:N	2.41	0.49
1:I:313:GLU:O	1:I:314:MET:HG3	2.12	0.49
2:L:410:VAL:O	2:L:423:MET:HA	2.13	0.49
2:L:155:ASN:ND2	3:L:841:NAG:C7	2.76	0.49
1:I:182:SER:C	1:I:184:ALA:N	2.66	0.49
2:L:366:ASP:OD2	2:L:368:PHE:CE1	2.66	0.49
2:L:23:TYR:HB3	2:L:116:SER:OG	2.12	0.49
2:L:21:CYS:C	2:L:22:ILE:HD12	2.33	0.49
2:L:285:LEU:CD2	2:L:285:LEU:N	2.70	0.49
1:I:284:ILE:HD13	1:I:307:TRP:CZ3	2.47	0.49
2:L:374:GLU:O	2:L:380:SER:HA	2.13	0.49
1:I:91:LYS:HE3	1:I:120:HIS:CE1	2.47	0.49
2:L:23:TYR:CE1	2:L:100:GLN:HG3	2.47	0.49
2:L:355:VAL:O	2:L:355:VAL:HG12	2.12	0.49
2:L:100:GLN:O	2:L:104:GLU:HG3	2.13	0.49
1:I:92:LEU:HB3	1:I:158:TYR:CE1	2.48	0.49
2:L:5:VAL:HG12	2:L:6:ASP:N	2.28	0.49
2:L:284:ILE:C	2:L:285:LEU:HD23	2.33	0.48
2:L:51:LEU:CD2	2:L:123:PHE:HA	2.43	0.48
1:I:186:ILE:O	1:I:189:TRP:N	2.46	0.48
1:I:71:ASN:ND2	1:I:72:ASP:H	2.11	0.48
1:I:170:LEU:N	1:I:170:LEU:HD23	2.28	0.48
2:L:414:GLU:OE1	2:L:416:PRO:HG2	2.14	0.48
1:I:147:PHE:CE2	1:I:186:ILE:HG12	2.49	0.48
1:I:92:LEU:HB3	1:I:158:TYR:HE1	1.79	0.48
1:I:19:PRO:C	1:I:21:CYS:N	2.64	0.48
1:I:261:ARG:HD3	1:I:262:ARG:H	1.78	0.48
1:I:400:VAL:HG12	1:I:401:THR:N	2.28	0.48
1:I:89:MET:HA	1:I:166:TYR:CE2	2.49	0.48
1:I:178:ASN:HB3	1:I:181:GLN:HG3	1.96	0.48
1:I:76:ILE:HG22	1:I:77:PHE:N	2.29	0.48
2:L:326:GLU:HG2	2:L:372:PHE:HD2	1.79	0.47
1:I:346:PRO:HG3	1:I:363:TYR:CE2	2.49	0.47
2:L:158:TYR:HB2	2:L:353:GLY:O	2.13	0.47
1:I:396:ASN:HB2	2:L:240:TYR:CZ	2.49	0.47
1:I:260:TYR:CZ	1:I:400:VAL:HG11	2.50	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:214:VAL:HG22	2:L:389:VAL:HG22	1.97	0.47
2:L:215:LEU:CD1	2:L:215:LEU:H	2.23	0.47
2:L:96:ASN:O	2:L:97:ASP:HB2	2.14	0.47
1:I:261:ARG:NH1	5:I:658:HOH:O	2.46	0.47
2:L:355:VAL:O	2:L:357:GLU:N	2.48	0.47
2:L:355:VAL:CG2	2:L:362:LEU:HD11	2.45	0.47
2:L:308:LEU:HD13	2:L:413:ARG:NH2	2.30	0.47
1:I:87:PHE:CE1	1:I:335:LEU:HD13	2.49	0.47
1:I:22:ILE:HG22	1:I:23:TYR:N	2.30	0.47
2:L:119:ILE:O	2:L:122:PHE:N	2.48	0.47
1:I:261:ARG:CZ	1:I:261:ARG:HA	2.44	0.47
2:L:217:ASN:ND2	2:L:218:THR:N	2.63	0.47
1:I:390:ILE:HA	2:L:319:HIS:HB2	1.96	0.47
3:I:801:NAG:O7	3:I:801:NAG:H3	2.14	0.47
2:L:276:GLY:O	2:L:277:ASP:HB2	2.14	0.47
1:I:281:MET:HA	1:I:411:PHE:O	2.14	0.47
1:I:71:ASN:HB3	1:I:74:ASP:OD2	2.15	0.47
2:L:146:LEU:HD11	2:L:162:SER:HB3	1.97	0.47
1:I:51:LEU:HD22	1:I:122:PHE:HB3	1.97	0.47
1:I:121:PHE:O	1:I:124:ALA:HB3	2.15	0.47
2:L:404:ALA:HB2	2:L:428:ASN:CG	2.33	0.46
1:I:91:LYS:CB	1:I:102:LEU:HD13	2.42	0.46
1:I:326:GLU:CG	1:I:370:LYS:HE3	2.44	0.46
1:I:40:ILE:HG23	1:I:40:ILE:O	2.14	0.46
2:L:163:GLU:HG2	2:L:163:GLU:O	2.15	0.46
1:I:261:ARG:HH11	1:I:261:ARG:HA	1.80	0.46
2:L:215:LEU:CD1	2:L:215:LEU:N	2.76	0.46
2:L:131:TYR:CZ	2:L:142:SER:HB2	2.50	0.46
1:I:346:PRO:HA	1:I:363:TYR:CD1	2.50	0.46
1:I:155:ASN:O	1:I:158:TYR:HB3	2.16	0.46
2:L:405:ASN:O	2:L:406:ARG:C	2.54	0.46
2:L:398:ASN:CG	2:L:399:ARG:H	2.19	0.46
1:I:351:LEU:C	1:I:353:GLY:H	2.18	0.46
1:I:87:PHE:HD2	1:I:87:PHE:H	1.63	0.46
1:I:229:PHE:CE2	1:I:254:GLN:HG2	2.51	0.46
1:I:265:GLU:CD	1:I:290:LYS:HZ2	2.18	0.46
1:I:151:SER:O	1:I:152:LEU:O	2.34	0.46
3:I:861:NAG:H5	3:I:862:NAG:C1	2.46	0.46
2:L:95:CYS:O	2:L:350:LYS:HB2	2.16	0.46
1:I:51:LEU:HD22	1:I:122:PHE:HB2	1.97	0.46
2:L:71:ASN:O	2:L:73:ASN:N	2.49	0.45
1:I:190:VAL:HB	1:I:201:VAL:CG2	2.35	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:101:GLN:O	2:L:105:VAL:HG23	2.15	0.45
1:I:351:LEU:O	1:I:353:GLY:N	2.49	0.45
2:L:221:PHE:CE1	2:L:279:ILE:HD13	2.52	0.45
2:L:333:GLU:O	2:L:337:ASP:OD1	2.34	0.45
1:I:319:HIS:CB	1:I:403:LYS:HG3	2.46	0.45
2:L:326:GLU:HG2	2:L:372:PHE:CD2	2.52	0.45
2:L:16:PRO:O	2:L:17:MET:C	2.55	0.45
2:L:172:PRO:O	2:L:173:LEU:HG	2.16	0.45
1:I:327:ASP:O	1:I:370:LYS:HD2	2.16	0.45
2:L:43:ALA:HB3	2:L:46:ARG:CB	2.47	0.45
1:I:56:SER:OG	1:I:420:ILE:HD12	2.17	0.45
1:I:89:MET:HE2	1:I:166:TYR:HB3	1.99	0.45
1:I:253:TYR:OH	1:I:255:GLU:OE2	2.34	0.45
5:I:609:HOH:O	2:L:239:PHE:HB2	2.17	0.45
1:I:331:LEU:C	1:I:335:LEU:HD12	2.38	0.45
1:I:219:ILE:HG22	1:I:220:LEU:N	2.32	0.45
2:L:145:ARG:HG2	2:L:147:PHE:CZ	2.51	0.44
2:L:200:ASP:O	2:L:201:VAL:C	2.56	0.44
1:I:276:GLY:O	1:I:277:ASP:HB2	2.17	0.44
2:L:404:ALA:HB2	2:L:428:ASN:CB	2.47	0.44
2:L:418:ASN:O	2:L:418:ASN:OD1	2.35	0.44
2:L:294:LYS:O	2:L:294:LYS:HG2	2.18	0.44
2:L:71:ASN:N	2:L:74:ASP:OD2	2.46	0.44
2:L:199:THR:O	2:L:200:ASP:CB	2.64	0.44
1:I:78:LEU:H	1:I:78:LEU:HD23	1.80	0.44
2:L:258:PHE:O	2:L:314:MET:N	2.49	0.44
2:L:263:VAL:O	2:L:264:ALA:C	2.57	0.44
2:L:258:PHE:HB2	2:L:316:LEU:HD21	2.00	0.44
1:I:180:GLU:C	1:I:182:SER:H	2.22	0.43
1:I:346:PRO:HG3	1:I:363:TYR:CZ	2.53	0.43
2:L:404:ALA:HB2	2:L:428:ASN:HB2	1.99	0.43
2:L:91:LYS:HZ3	2:L:120:HIS:CE1	2.34	0.43
1:I:332:LYS:HG3	1:I:344:PHE:CD1	2.52	0.43
1:I:285:LEU:HD12	1:I:286:PRO:HD2	2.00	0.43
1:I:219:ILE:HB	1:I:422:PHE:CZ	2.53	0.43
2:L:245:GLU:HG2	2:L:246:SER:N	2.34	0.43
1:I:360:ASP:OD2	1:I:361:ASP:N	2.52	0.43
2:L:358:GLY:O	2:L:359:ARG:C	2.57	0.43
1:I:239:PHE:CD2	1:I:429:PRO:HG3	2.53	0.43
1:I:198:ILE:HG23	1:I:370:LYS:HG2	2.01	0.43
1:I:151:SER:HB3	1:I:152:LEU:H	1.54	0.43
1:I:192:ASN:ND2	3:I:861:NAG:C1	2.80	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:157:THR:OG1	3:L:841:NAG:H5	2.19	0.43
2:L:258:PHE:HB2	2:L:316:LEU:CD2	2.48	0.43
2:L:157:THR:CB	3:L:841:NAG:H5	2.49	0.43
1:I:288:PRO:HB2	1:I:289:GLU:OE2	2.18	0.43
2:L:171:GLN:HA	2:L:172:PRO:HD3	1.81	0.43
1:I:192:ASN:ND2	3:I:861:NAG:C2	2.77	0.43
1:I:153:THR:O	1:I:354:ILE:HG22	2.18	0.43
1:I:241:LYS:CD	1:I:428:ASN:HD22	2.32	0.43
1:I:89:MET:O	1:I:92:LEU:HB2	2.19	0.42
1:I:284:ILE:HB	1:I:409:LEU:HB2	2.01	0.42
1:I:14:ASP:C	1:I:16:PRO:HD3	2.40	0.42
1:I:154:PHE:O	1:I:155:ASN:C	2.57	0.42
1:I:22:ILE:CG2	1:I:23:TYR:N	2.81	0.42
2:L:356:ALA:C	2:L:358:GLY:N	2.73	0.42
1:I:167:GLY:O	1:I:168:ALA:HB2	2.19	0.42
1:I:94:ALA:HB3	1:I:99:LEU:HD13	2.01	0.42
1:I:341:VAL:O	1:I:343:LEU:N	2.52	0.42
2:L:80:PRO:HG3	2:L:423:MET:HE2	2.02	0.42
1:I:421:ILE:HG22	1:I:422:PHE:CD2	2.54	0.42
2:L:259:ARG:NH2	2:L:311:LEU:HB2	2.33	0.42
2:L:365:SER:OG	2:L:389:VAL:HG12	2.20	0.42
1:I:155:ASN:HA	4:I:841:NAG:C8	2.50	0.42
1:I:113:GLU:O	1:I:114:LYS:C	2.57	0.42
1:I:414:GLU:HB2	1:I:421:ILE:HD11	2.01	0.42
2:L:356:ALA:O	2:L:358:GLY:N	2.53	0.42
2:L:136:LYS:O	2:L:137:SER:HB2	2.20	0.42
1:I:189:TRP:O	1:I:192:ASN:HB3	2.20	0.42
2:L:245:GLU:HG2	2:L:246:SER:H	1.85	0.42
2:L:356:ALA:C	2:L:358:GLY:H	2.22	0.42
2:L:138:SER:HB3	2:L:279:ILE:HD12	2.02	0.42
2:L:225:TRP:CD1	2:L:379:GLY:HA2	2.55	0.42
1:I:334:GLN:O	1:I:335:LEU:C	2.58	0.42
1:I:299:LEU:C	1:I:299:LEU:HD13	2.41	0.42
1:I:337:ASP:OD2	1:I:337:ASP:O	2.37	0.42
1:I:81:LEU:HD21	1:I:130:LEU:CD1	2.50	0.42
1:I:183:ARG:NE	1:I:203:PRO:O	2.52	0.41
1:I:354:ILE:CD1	1:I:354:ILE:N	2.77	0.41
1:I:341:VAL:C	1:I:343:LEU:H	2.24	0.41
2:L:217:ASN:C	2:L:217:ASN:ND2	2.74	0.41
1:I:222:LYS:HB2	1:I:381:GLU:HG2	2.01	0.41
1:I:343:LEU:CD1	1:I:364:VAL:HG23	2.49	0.41
2:L:273:PRO:HA	2:L:280:THR:HG22	2.02	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:243:ASP:OD1	1:I:244:GLY:N	2.42	0.41
2:L:341:VAL:CG2	2:L:342:ASP:N	2.83	0.41
1:I:138:SER:OG	1:I:138:SER:O	2.37	0.41
1:I:305:GLN:HG2	5:I:586:HOH:O	2.21	0.41
2:L:372:PHE:CD1	2:L:372:PHE:C	2.93	0.41
1:I:48:VAL:HG13	1:I:126:LEU:HD13	2.03	0.41
1:I:131:TYR:CZ	1:I:142:SER:HB2	2.56	0.41
2:L:102:LEU:CD2	2:L:340:LEU:HD11	2.51	0.41
2:L:91:LYS:HE2	2:L:91:LYS:HB3	1.80	0.41
1:I:114:LYS:HD2	1:I:114:LYS:N	2.34	0.41
1:I:241:LYS:HD3	1:I:428:ASN:ND2	2.35	0.41
2:L:221:PHE:HE1	2:L:279:ILE:HD13	1.86	0.41
1:I:81:LEU:HD21	1:I:130:LEU:HD11	2.02	0.41
1:I:323:PHE:CZ	1:I:375:VAL:HG21	2.56	0.41
1:I:252:MET:HG3	1:I:322:ARG:HG2	2.01	0.41
1:I:361:ASP:O	1:I:361:ASP:OD1	2.39	0.40
2:L:105:VAL:HG13	2:L:338:MET:HB3	2.02	0.40
1:I:200:ASP:OD1	1:I:370:LYS:HE2	2.21	0.40
1:I:40:ILE:HA	1:I:41:PRO:HD2	2.01	0.40
1:I:150:LYS:N	1:I:173:LEU:O	2.54	0.40
2:L:260:TYR:O	2:L:261:ARG:HB2	2.21	0.40
2:L:253:TYR:CD1	2:L:253:TYR:C	2.94	0.40
2:L:253:TYR:CE2	2:L:317:VAL:CG1	3.05	0.40
2:L:283:LEU:HD11	2:L:320:MET:HE3	2.02	0.40
2:L:373:LEU:HD12	2:L:374:GLU:H	1.86	0.40
1:I:119:ILE:HG22	1:I:120:HIS:N	2.35	0.40
2:L:251:MET:CE	2:L:319:HIS:HB3	2.52	0.40
2:L:372:PHE:N	2:L:383:ALA:O	2.54	0.40
2:L:203:PRO:HB2	2:L:204:SER:H	1.69	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	I	405/427 (95%)	309 (76%)	65 (16%)	31 (8%)	1	3
2	L	404/427 (95%)	328 (81%)	47 (12%)	29 (7%)	2	4
All	All	809/854 (95%)	637 (79%)	112 (14%)	60 (7%)	2	3

All (60) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	I	42	GLU
1	I	152	LEU
1	I	170	LEU
1	I	176	LYS
1	I	430	CYS
2	L	19	PRO
2	L	96	ASN
2	L	112	SER
2	L	116	SER
2	L	117	ASP
2	L	136	LYS
2	L	200	ASP
2	L	203	PRO
2	L	228	LYS
2	L	250	SER
2	L	264	ALA
2	L	356	ALA
2	L	403	LYS
1	I	184	ALA
1	I	245	GLU
1	I	350	LYS
1	I	383	ALA
1	I	384	ALA
2	L	44	THR
2	L	69	SER
2	L	72	ASP
2	L	243	ASP
1	I	77	PHE
1	I	168	ALA
1	I	183	ARG
1	I	243	ASP
1	I	244	GLY
1	I	337	ASP
1	I	342	ASP
2	L	17	MET

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Mol	Chain	Res	Type
2	L	68	ASP
2	L	293	ALA
2	L	359	ARG
1	I	38	GLN
1	I	82	SER
1	I	158	TYR
1	I	203	PRO
1	I	352	PRO
1	I	386	THR
2	L	244	GLY
2	L	357	GLU
1	I	12	PRO
1	I	43	ALA
1	I	116	SER
1	I	130	LEU
1	I	171	GLN
2	L	222	LYS
2	L	406	ARG
1	I	40	ILE
1	I	112	SER
2	L	275	LYS
2	L	333	GLU
1	I	41	PRO
2	L	263	VAL
2	L	119	ILE

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	I	341/378 (90%)	323 (95%)	18 (5%)	32	67
2	L	312/379 (82%)	296 (95%)	16 (5%)	33	69
All	All	653/757 (86%)	619 (95%)	34 (5%)	32	68

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	13	ARG
1	I	17	MET
1	I	71	ASN
1	I	78	LEU
1	I	92	LEU
1	I	97	ASP
1	I	109	ASP
1	I	114	LYS
1	I	153	THR
1	I	170	LEU
1	I	216	VAL
1	I	309	ASP
1	I	310	GLU
1	I	316	LEU
1	I	322	ARG
1	I	342	ASP
1	I	354	ILE
1	I	366	ASP
2	L	84	SER
2	L	97	ASP
2	L	123	PHE
2	L	162	SER
2	L	201	VAL
2	L	208	ASN
2	L	217	ASN
2	L	253	TYR
2	L	285	LEU
2	L	316	LEU
2	L	324	ARG
2	L	337	ASP
2	L	397	PRO
2	L	403	LYS
2	L	405	ASN
2	L	413	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	I	55	ASN
1	I	71	ASN
1	I	118	GLN
1	I	178	ASN
1	I	192	ASN

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Mol	Chain	Res	Type
1	I	334	GLN
1	I	428	ASN
2	L	55	ASN
2	L	65	HIS
2	L	71	ASN
2	L	96	ASN
2	L	144	ASN
2	L	208	ASN
2	L	217	ASN
2	L	254	GLN
2	L	336	GLN
2	L	405	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

4 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	$\# Z > 2$
4	NAG	I	841	1,4	12,14,15	0.59	0	15,19,21	0.84	1 (6%)
4	NAG	I	842	4	12,14,15	0.55	0	15,19,21	0.67	0
4	NAG	L	861	2,4	12,14,15	0.71	0	15,19,21	0.93	0
4	NAG	L	862	4	9,13,15	0.30	0	11,17,21	0.75	1 (9%)

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
4	NAG	I	841	1,4	1/1/5/7	0/6/23/26	0/1/1/1
4	NAG	I	842	4	-	0/6/23/26	0/1/1/1
4	NAG	L	861	2,4	-	0/6/23/26	0/1/1/1
4	NAG	L	862	4	-	0/4/22/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	L	862	NAG	C3-C2-N2	-2.20	108.95	112.26
4	I	841	NAG	C2-N2-C7	-2.13	119.51	123.09

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	I	841	NAG	C1

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry ⓘ

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	I	801	1	12,14,15	0.47	0	15,19,21	0.65	0
3	NAG	I	861	-	12,14,15	0.45	0	15,19,21	0.71	0
3	NAG	I	862	-	12,14,15	0.45	0	15,19,21	0.50	0
3	NAG	L	801	-	12,14,15	0.53	0	15,19,21	0.59	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	L	841	2	12,14,15	0.72	0	15,19,21	0.96	1 (6%)
3	NAG	L	842	-	12,14,15	0.45	0	15,19,21	0.57	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	I	801	1	-	0/6/23/26	0/1/1/1
3	NAG	I	861	-	-	0/6/23/26	0/1/1/1
3	NAG	I	862	-	-	0/6/23/26	0/1/1/1
3	NAG	L	801	-	-	1/6/23/26	0/1/1/1
3	NAG	L	841	2	-	0/6/23/26	0/1/1/1
3	NAG	L	842	-	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	841	NAG	C3-C4-C5	2.19	114.11	110.20

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	L	801	NAG	C1-C2-N2-C7

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	I	413/427 (96%)	0.23	28 (6%) 17 15	36, 79, 115, 151	0
2	L	410/427 (96%)	0.06	25 (6%) 21 20	42, 77, 121, 141	0
All	All	823/854 (96%)	0.15	53 (6%) 17 17	36, 78, 120, 151	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	240	TYR	6.7
2	L	398	ASN	5.5
2	L	242	ALA	5.2
1	I	395	LEU	5.0
1	I	144	ASN	4.7
2	L	239	PHE	4.6
1	I	360	ASP	4.4
1	I	361	ASP	4.2
2	L	394	SER	4.2
2	L	207	ILE	4.0
1	I	189	TRP	3.9
2	L	403	LYS	3.9
2	L	401	THR	3.9
2	L	246	SER	3.7
2	L	241	LYS	3.7
2	L	399	ARG	3.6
1	I	394	SER	3.6
1	I	153	THR	3.5
2	L	405	ASN	3.3
1	I	43	ALA	3.2
2	L	402	PHE	3.2
1	I	168	ALA	3.2
1	I	363	TYR	3.0
1	I	143	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
2	L	205	GLU	2.9
2	L	253	TYR	2.8
1	I	362	LEU	2.8
2	L	202	ILE	2.8
2	L	429	PRO	2.6
1	I	40	ILE	2.6
1	I	366	ASP	2.5
2	L	244	GLY	2.5
1	I	386	THR	2.4
2	L	243	ASP	2.4
1	I	90	THR	2.4
1	I	216	VAL	2.3
2	L	245	GLU	2.3
2	L	404	ALA	2.3
1	I	396	ASN	2.3
1	I	385	SER	2.3
2	L	200	ASP	2.3
1	I	44	THR	2.3
1	I	399	ARG	2.3
1	I	146	LEU	2.2
1	I	141	VAL	2.2
2	L	201	VAL	2.2
1	I	215	LEU	2.1
2	L	400	VAL	2.1
1	I	318	VAL	2.1
2	L	199	THR	2.1
1	I	142	SER	2.1
1	I	37	GLU	2.1
1	I	400	VAL	2.0

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

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

6.3 Carbohydrates ⓘ

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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	NAG	I	841	14/15	0.54	2.58	106,112,116,118	0
4	NAG	L	861	14/15	0.24	1.51	115,119,122,124	0
4	NAG	I	842	14/15	0.30	-	120,122,122,122	0
4	NAG	L	862	13/15	0.12	-	127,128,129,129	0

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	NAG	I	801	14/15	0.17	4.76	125,128,130,130	0
3	NAG	L	842	14/15	0.24	3.46	148,150,151,152	0
3	NAG	I	861	14/15	0.67	2.90	139,143,145,147	0
3	NAG	L	801	14/15	0.18	2.28	126,128,129,129	0
3	NAG	L	841	14/15	0.14	-0.34	86,93,96,97	0
3	NAG	I	862	14/15	0.60	-	160,162,163,163	0

6.5 Other polymers

There are no such residues in this entry.