



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:39 PM GMT

PDB ID : 4PJ6
Title : Crystal Structure of Human Insulin Regulated Aminopeptidase with Lysine in Active Site
Authors : Hermans, S.J.; Ascher, D.B.; Hancock, N.C.; Holien, J.K.; Michell, B.; Morton, C.J.; Parker, M.W.
Deposited on : 2014-05-12
Resolution : 2.96 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

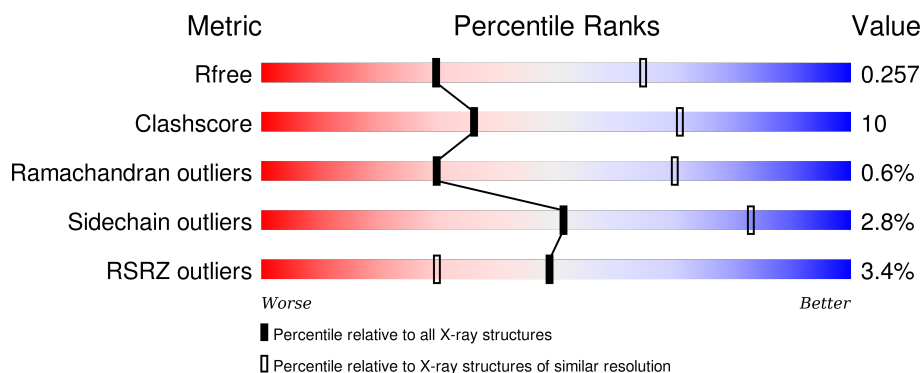
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.96 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	2184 (3.00-2.92)
Clashscore	102246	2552 (3.00-2.92)
Ramachandran outliers	100387	2468 (3.00-2.92)
Sidechain outliers	100360	2471 (3.00-2.92)
RSRZ outliers	91569	2201 (3.00-2.92)

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	872	 2% 74% 24% ..
1	B	872	 4% 74% 23% ..

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
3	LYS	A	1102	-	-	-	X
3	LYS	B	1102	-	-	-	X
4	NAG	B	1106	-	-	-	X
4	NAG	B	1109	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 14244 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 Leucyl-cystinyl aminopeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	860	Total	C	N	O	S	0	2	0
			6967	4502	1132	1306	27			
1	B	849	Total	C	N	O	S	0	0	0
			6861	4442	1107	1287	25			

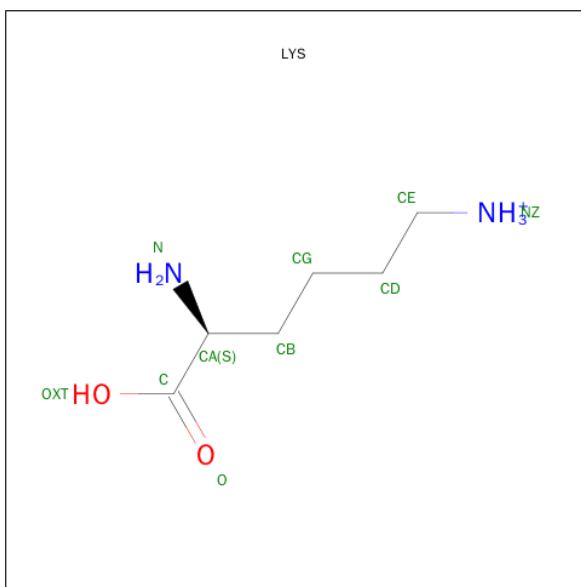
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	154	SER	-	expression tag	UNP Q9UIQ6
B	154	SER	-	expression tag	UNP Q9UIQ6

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is LYSINE (three-letter code: LYS) (formula: C₆H₁₅N₂O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			9	6	2	1		
3	B	1	Total	C	N	O	0	0
			9	6	2	1		

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

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

- Molecule 5 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	101	Total	O	0	0
			101	101		
6	B	57	Total	O	0	0
			57	57		

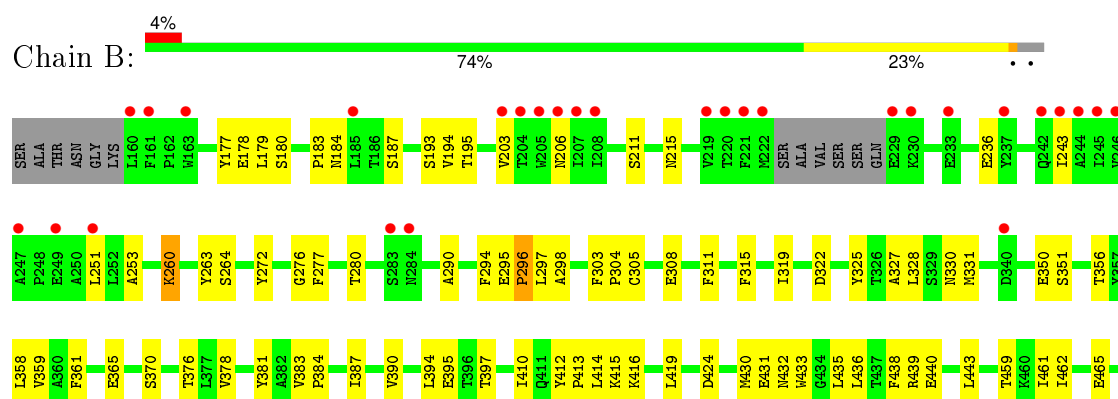
3 Residue-property plots

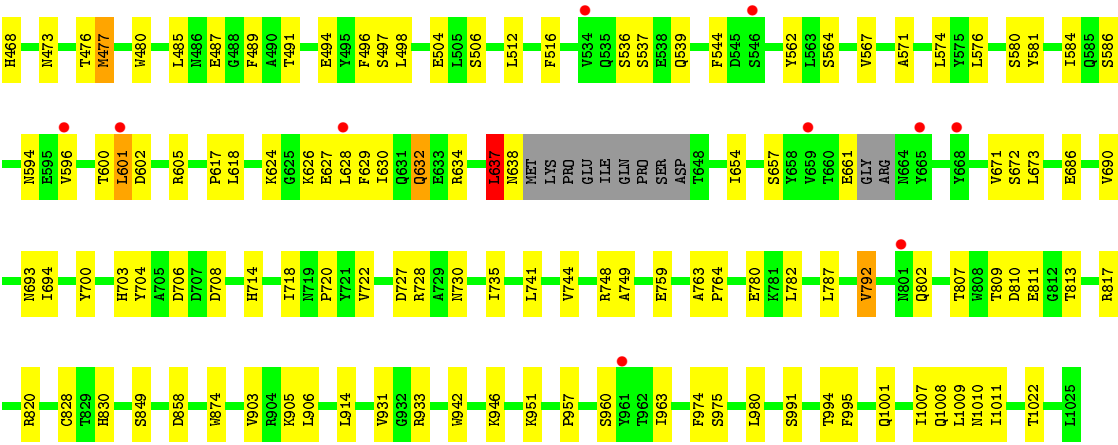
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Leucyl-cystinyl aminopeptidase



• Molecule 1: Leucyl-cystinyl aminopeptidase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	68.03Å 256.62Å 71.31Å 90.00° 114.35° 90.00°	Depositor
Resolution (Å)	48.32 – 2.96 48.27 – 2.96	Depositor EDS
% Data completeness (in resolution range)	99.6 (48.32-2.96) 99.6 (48.27-2.96)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.91 (at 2.96Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.185 , 0.257 0.185 , 0.257	Depositor DCC
R_{free} test set	2328 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	67.3	Xtriage
Anisotropy	0.119	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 73.1	EDS
Estimated twinning fraction	0.034 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 46198 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14244	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% 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, 0.375 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: ZN, NAG

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.67	1/7143 (0.0%)	0.76	1/9691 (0.0%)
1	B	0.65	1/7029 (0.0%)	0.72	0/9541
All	All	0.66	2/14172 (0.0%)	0.74	1/19232 (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	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	637	LEU	CG-CD1	7.31	1.78	1.51
1	A	230	LYS	CD-CE	5.40	1.64	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1025	LEU	CA-CB-CG	6.84	131.03	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	632	GLN	Sidechain

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	6967	0	6817	137	0
1	B	6861	0	6701	134	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	9	0	12	0	0
3	B	9	0	12	1	0
4	A	98	0	91	0	0
4	B	112	0	104	1	0
5	B	28	0	25	0	0
6	A	101	0	0	3	0
6	B	57	0	0	1	0
All	All	14244	0	13762	268	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 10.

All (268) 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:637:LEU:CG	1:B:637:LEU:CD1	1.78	1.56
1:B:637:LEU:HD13	1:B:1009:LEU:CD2	1.80	1.11
1:B:637:LEU:HD13	1:B:1009:LEU:HD21	1.29	1.10
1:A:594:ASN:OD1	1:A:602:ASP:O	1.79	1.00
1:B:637:LEU:CD1	1:B:1009:LEU:HD21	2.03	0.87
1:A:597:THR:O	1:A:600:THR:HG22	1.79	0.83
1:A:307:ASP:HA	1:A:356:THR:HG21	1.61	0.81
1:B:397:THR:HG21	1:B:419:LEU:HD22	1.63	0.81
1:A:748:ARG:HG3	6:A:1202:HOH:O	1.81	0.79
1:A:600:THR:HG23	1:A:601:LEU:HD12	1.66	0.76
1:B:975:SER:HB3	1:B:1010:ASN:HB3	1.66	0.76
1:A:564:SER:HB3	1:A:567:VAL:HG23	1.67	0.75
1:B:637:LEU:HD12	1:B:638:ASN:N	2.02	0.75
1:A:600:THR:HG23	1:A:601:LEU:CD1	2.18	0.74
1:B:376:THR:HA	1:B:415:LYS:O	1.88	0.73
1:B:330:ASN:O	1:B:416:LYS:HE2	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:741:LEU:HD23	1:A:743:LYS:HG3	1.71	0.72
1:A:936:PRO:HA	1:B:903:VAL:HG21	1.72	0.71
1:A:794:ARG:HH11	1:A:1025:LEU:HA	1.55	0.71
1:A:741:LEU:CD2	1:A:743:LYS:HG3	2.22	0.70
1:A:432:ASN:HB2	1:A:435:LEU:O	1.93	0.69
1:B:295:GLU:HB3	1:B:430:MET:HE1	1.74	0.69
1:B:637:LEU:HD13	1:B:1009:LEU:HD22	1.73	0.68
1:B:637:LEU:CD1	1:B:637:LEU:CB	2.71	0.68
1:B:637:LEU:HA	1:B:741:LEU:HG	1.76	0.67
1:B:906:LEU:HD22	1:B:931:VAL:HG22	1.76	0.67
1:B:632:GLN:OE1	1:B:654:ILE:HD12	1.94	0.67
1:B:763:ALA:HB3	1:B:764:PRO:HD3	1.77	0.66
1:A:806:GLN:HA	6:A:1263:HOH:O	1.95	0.66
1:A:614:LYS:HE3	1:A:634:ARG:HH21	1.62	0.65
1:A:820:ARG:HG3	1:A:824:LEU:HD12	1.79	0.65
1:B:571:ALA:HB2	1:B:596:VAL:HG21	1.78	0.64
1:A:372:ASP:OD1	1:A:374:ASN:O	2.16	0.64
1:A:958:LEU:HG	1:A:997:LEU:HD11	1.80	0.63
1:A:571:ALA:HB2	1:A:596:VAL:HG21	1.80	0.62
1:B:632:GLN:OE1	1:B:654:ILE:CD1	2.47	0.62
1:B:203:VAL:HG22	1:B:253:ALA:H	1.65	0.62
1:A:904:ARG:NH1	1:B:780:GLU:OE2	2.32	0.62
1:A:606:MET:CE	1:A:655:PRO:HG3	2.30	0.61
1:A:906:LEU:HD22	1:A:931:VAL:HG22	1.82	0.61
1:A:626:LYS:HE2	1:A:685:GLU:HA	1.81	0.61
1:B:628:LEU:HD22	1:B:690:VAL:HG21	1.83	0.61
1:B:184:ASN:O	1:B:187:SER:O	2.19	0.61
1:B:657:SER:HB2	1:B:694:ILE:HD12	1.83	0.60
1:B:628:LEU:CD1	1:B:630:ILE:HD11	2.31	0.60
1:A:658:TYR:CZ	1:A:670:SER:HB3	2.36	0.60
1:A:594:ASN:ND2	6:A:1285:HOH:O	2.29	0.60
1:B:624:LYS:O	1:B:627:GLU:HG3	2.01	0.59
1:A:477:MET:HG2	1:A:477:MET:O	2.02	0.59
1:B:319:ILE:HD11	1:B:327:ALA:HB1	1.85	0.59
1:A:233:GLU:HG2	1:A:246:VAL:HB	1.83	0.59
1:B:397:THR:CG2	1:B:419:LEU:HD22	2.33	0.59
1:A:205:TRP:CZ2	1:A:250:ALA:HB2	2.37	0.59
1:A:272:TYR:CZ	1:A:298:ALA:HB2	2.38	0.59
1:B:628:LEU:HD13	1:B:630:ILE:HD11	1.85	0.59
1:A:630:ILE:CD1	1:A:674:LEU:HD13	2.33	0.59
1:A:294:PHE:HA	1:A:298:ALA:HB3	1.83	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:431:GLU:HG3	1:A:468:HIS:HB3	1.84	0.58
1:A:380:ILE:HG21	1:A:390:VAL:HB	1.84	0.58
1:B:294:PHE:HA	1:B:298:ALA:HB3	1.85	0.58
1:A:515:ARG:NH2	1:A:696:MET:O	2.34	0.58
1:B:177:TYR:CD2	1:B:315:PHE:CE2	2.91	0.58
1:B:564:SER:HB3	1:B:567:VAL:H	1.69	0.58
1:B:960:SER:HB3	1:B:963:ILE:HG22	1.86	0.58
1:B:295:GLU:HB3	1:B:430:MET:CE	2.33	0.58
1:B:991:SER:HB2	1:B:994:THR:OG1	2.04	0.58
1:A:175:LEU:HD11	1:A:199:GLN:HB2	1.86	0.57
1:B:432:ASN:HB2	1:B:435:LEU:O	2.04	0.57
1:A:975:SER:HB3	1:A:1010:ASN:HB3	1.86	0.57
1:B:914:LEU:HD11	1:B:951:LYS:HG2	1.86	0.57
1:A:495:TYR:CD1	1:A:511:PHE:HB2	2.39	0.57
1:A:168:LEU:HD13	1:A:306:PHE:HD2	1.70	0.57
1:A:874:TRP:CE2	1:A:905:LYS:HD3	2.40	0.57
1:B:331:MET:SD	1:B:351:SER:HA	2.45	0.56
1:A:220:THR:OG1	1:A:229:GLU:HB3	2.05	0.56
1:A:634:ARG:HD3	1:A:636:PHE:HB2	1.86	0.56
1:B:476:THR:O	1:B:584:ILE:HG13	2.06	0.55
1:B:397:THR:HG21	1:B:419:LEU:CD2	2.36	0.55
1:B:319:ILE:CD1	1:B:327:ALA:HB1	2.36	0.55
1:A:720:PRO:HB2	1:A:728:ARG:HD3	1.88	0.55
1:B:487:GLU:OE2	3:B:1102:LYS:N	2.41	0.55
1:B:325:TYR:CD1	1:B:365:GLU:HG3	2.42	0.54
1:A:209:LEU:HD11	1:A:243:ILE:HD11	1.88	0.54
1:A:630:ILE:HD12	1:A:674:LEU:HD13	1.88	0.54
1:B:308:GLU:HB2	1:B:311:PHE:HD2	1.72	0.54
1:A:176:ARG:HG2	1:A:314:THR:OG1	2.07	0.54
1:B:383:VAL:HG23	1:B:384:PRO:HD2	1.89	0.54
1:A:1018:LEU:HA	1:A:1021:LEU:HD12	1.90	0.53
1:A:936:PRO:HA	1:B:903:VAL:CG2	2.38	0.53
1:A:515:ARG:O	1:A:519:MET:HG3	2.08	0.53
1:B:580:SER:HB3	1:B:581:TYR:CD2	2.43	0.53
1:A:205:TRP:CE2	1:A:250:ALA:HB2	2.43	0.53
1:A:282:GLU:N	1:A:385:GLU:OE1	2.41	0.53
1:B:735:ILE:HG12	1:B:749:ALA:HA	1.90	0.53
1:A:874:TRP:CD2	1:A:905:LYS:HD3	2.44	0.53
1:A:796:PHE:CG	1:A:832:LEU:HD13	2.44	0.52
1:B:637:LEU:CD2	1:B:637:LEU:CD1	2.79	0.52
1:A:458:VAL:HA	1:A:461:ILE:HG22	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:251:LEU:HD23	1:A:257:TYR:CD2	2.44	0.52
1:B:236:GLU:HG2	1:B:243:ILE:HG22	1.91	0.52
1:A:186:THR:HG22	1:A:288:TYR:CE1	2.45	0.52
1:A:405:GLN:HG2	1:A:411:GLN:HA	1.92	0.51
1:A:561:THR:HG21	1:A:697:ASN:ND2	2.24	0.51
1:A:358:LEU:CD2	1:A:473:ASN:OD1	2.58	0.51
1:A:537:SER:HA	1:A:540:ILE:HD12	1.93	0.51
1:A:318:LYS:HD3	1:A:347:GLU:HG3	1.93	0.51
1:A:761:HIS:O	1:A:764:PRO:HD2	2.10	0.51
1:A:584:ILE:HG23	1:A:588:ASP:HB2	1.91	0.51
1:A:762:THR:HA	1:A:819:LEU:HD13	1.93	0.51
1:A:656:LEU:O	1:A:671:VAL:HG23	2.11	0.50
1:A:789:SER:O	1:A:793:THR:HG23	2.11	0.50
1:A:421:ALA:HB1	1:A:440:GLU:HA	1.92	0.50
1:B:179:LEU:HD13	1:B:194:VAL:HG22	1.92	0.50
1:B:874:TRP:CD2	1:B:905:LYS:HD3	2.47	0.50
1:A:325:TYR:CD1	1:A:365:GLU:HG3	2.46	0.50
1:B:813:THR:O	1:B:817:ARG:HG2	2.11	0.50
1:A:310:ALA:HA	1:A:583:SER:OG	2.11	0.50
1:B:504:GLU:CD	1:B:504:GLU:H	2.14	0.50
1:B:506:SER:N	6:B:1253:HOH:O	2.45	0.50
1:A:617:PRO:HB3	1:A:652:TRP:CE3	2.46	0.50
1:B:296:PRO:HB2	1:B:297:LEU:HD13	1.94	0.49
1:B:626:LYS:HG3	1:B:686:GLU:HG3	1.94	0.49
1:A:355:SER:HB2	1:A:358:LEU:HD12	1.93	0.49
1:A:762:THR:HB	1:A:819:LEU:HB2	1.94	0.49
1:B:459:THR:HG22	1:B:498:LEU:HD11	1.94	0.49
1:A:796:PHE:CD1	1:A:832:LEU:HD13	2.48	0.49
1:A:360:ALA:HB1	1:A:435:LEU:HD23	1.95	0.49
1:B:600:THR:HB	1:B:601:LEU:HD13	1.94	0.49
1:A:219:VAL:HG13	1:A:232:ALA:O	2.12	0.48
1:A:613:GLN:OE1	1:A:650:TYR:HA	2.12	0.48
1:A:403:PHE:HB2	1:A:501:ILE:HD11	1.95	0.48
1:A:476:THR:HG22	1:A:477:MET:N	2.29	0.48
1:B:703:HIS:ND1	1:B:704:TYR:N	2.62	0.48
1:B:397:THR:CG2	1:B:419:LEU:CD2	2.92	0.48
1:B:438:PHE:CD1	1:B:443:LEU:HD11	2.49	0.48
1:B:178:GLU:HB3	1:B:195:THR:HB	1.95	0.48
1:B:671:VAL:HG22	1:B:672:SER:H	1.79	0.48
1:B:974:PHE:O	1:B:1007:ILE:HG23	2.14	0.48
1:B:215:ASN:HB3	1:B:264:SER:HB2	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:492:PHE:HZ	1:A:560:LYS:HD3	1.79	0.47
1:A:180:SER:HA	1:A:318:LYS:O	2.14	0.47
1:B:562:TYR:CZ	1:B:673:LEU:HD13	2.49	0.47
1:A:456:LYS:O	1:A:460:LYS:HG3	2.14	0.47
1:A:579:HIS:HB3	1:A:584:ILE:HD11	1.95	0.47
1:A:357:TYR:CE1	1:A:358:LEU:HG	2.50	0.47
1:B:322:ASP:HB2	1:B:325:TYR:CD2	2.50	0.47
1:B:296:PRO:HB2	1:B:297:LEU:CD1	2.44	0.47
1:A:606:MET:HE3	1:A:655:PRO:HG3	1.96	0.47
1:B:624:LYS:HB2	1:B:629:PHE:HE2	1.79	0.47
1:A:469:GLN:O	1:A:473:ASN:HB2	2.15	0.46
1:B:574:LEU:HD11	4:B:1109:NAG:H81	1.97	0.46
1:B:807:THR:O	1:B:820:ARG:HD3	2.15	0.46
1:A:790:ARG:NH1	1:A:1024:TRP:O	2.48	0.46
1:B:272:TYR:OH	1:B:296:PRO:HD2	2.15	0.46
1:B:358:LEU:HD22	1:B:473:ASN:OD1	2.15	0.46
1:B:693:ASN:HB2	1:B:700:TYR:CE2	2.51	0.46
1:A:713:ILE:HG23	1:A:752:LEU:HA	1.97	0.46
1:A:373:VAL:HG11	1:A:401:LEU:HD23	1.98	0.46
1:B:211:SER:HB2	1:B:304:PRO:HB3	1.97	0.46
1:B:810:ASP:HB3	1:B:817:ARG:CZ	2.46	0.46
1:A:790:ARG:HH22	1:A:1022:THR:HA	1.80	0.46
1:B:387:ILE:O	1:B:390:VAL:HG22	2.16	0.46
1:B:637:LEU:H	1:B:637:LEU:HG	1.46	0.45
1:B:476:THR:HG22	1:B:477:MET:O	2.16	0.45
1:B:703:HIS:CE1	1:B:704:TYR:O	2.70	0.45
1:A:376:THR:HA	1:A:415:LYS:O	2.17	0.45
1:A:300:ARG:HH21	1:A:307:ASP:HB3	1.82	0.45
1:A:422:ILE:O	1:A:440:GLU:HB2	2.16	0.45
1:B:370:SER:HA	1:B:378:VAL:O	2.17	0.45
1:B:1008:GLN:HA	1:B:1011:ILE:HD12	1.98	0.45
1:A:331:MET:SD	1:A:351:SER:HA	2.57	0.45
1:B:995:PHE:O	1:B:1001:GLN:NE2	2.48	0.45
1:A:454:ASP:O	1:A:458:VAL:HG12	2.17	0.45
1:B:744:VAL:HG13	1:B:748:ARG:HD2	1.97	0.45
1:A:502:PHE:O	1:A:503:LYS:C	2.55	0.45
1:B:975:SER:CB	1:B:1010:ASN:HB3	2.42	0.45
1:B:536:SER:HB3	1:B:539:GLN:HB2	1.99	0.44
1:B:328:LEU:HD21	1:B:381:TYR:CZ	2.52	0.44
1:B:544:PHE:CD1	1:B:544:PHE:N	2.85	0.44
1:A:984:GLN:HG2	1:A:988:GLU:OE2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:489:PHE:O	1:A:493:MET:HG2	2.17	0.44
1:B:809:THR:HB	1:B:810:ASP:H	1.63	0.44
1:A:440:GLU:HG2	1:A:444:LEU:HD11	2.00	0.44
1:B:809:THR:OG1	1:B:811:GLU:HG2	2.18	0.44
1:B:817:ARG:HB2	1:B:858:ASP:OD2	2.18	0.44
1:B:412:TYR:HA	1:B:413:PRO:HD3	1.91	0.44
1:B:303:PHE:CE2	1:B:305:CYS:HB3	2.53	0.44
1:B:439:ARG:O	1:B:440:GLU:C	2.56	0.44
1:A:270:SER:OG	1:A:271:TYR:N	2.50	0.44
1:A:984:GLN:HG3	1:A:995:PHE:HE1	1.82	0.43
1:B:394:LEU:O	1:B:395:GLU:C	2.57	0.43
1:A:451:SER:HB3	1:A:454:ASP:CG	2.38	0.43
1:B:617:PRO:HG2	1:B:700:TYR:HB3	2.00	0.43
1:A:914:LEU:HD13	1:A:924:LEU:HD21	2.01	0.43
1:A:383:VAL:HG22	1:A:384:PRO:HD2	2.01	0.43
1:A:405:GLN:HG3	1:A:470:TRP:HZ2	1.84	0.43
1:B:277:PHE:CE2	1:B:424:ASP:HB2	2.54	0.43
1:B:594:ASN:OD1	1:B:602:ASP:O	2.36	0.43
1:A:564:SER:HB3	1:A:567:VAL:CG2	2.44	0.43
1:B:957:PRO:O	1:B:960:SER:HB2	2.19	0.43
1:A:975:SER:CB	1:A:1010:ASN:HB3	2.49	0.43
1:B:975:SER:HB3	1:B:1010:ASN:CB	2.44	0.43
1:B:476:THR:HG22	1:B:477:MET:N	2.33	0.43
1:A:358:LEU:HD22	1:A:473:ASN:OD1	2.17	0.43
1:B:195:THR:OG1	1:B:260:LYS:HG3	2.18	0.43
1:B:491:THR:O	1:B:494:GLU:HB3	2.18	0.43
1:A:634:ARG:HG2	1:A:634:ARG:HH11	1.83	0.43
1:A:178:GLU:HB3	1:A:195:THR:HB	2.01	0.43
1:B:376:THR:HG22	1:B:414:LEU:O	2.19	0.43
1:A:410:ILE:HD11	1:A:577:HIS:HB2	2.00	0.43
1:A:796:PHE:HB2	1:A:832:LEU:HD13	2.01	0.42
1:B:496:PHE:O	1:B:497:SER:C	2.58	0.42
1:B:410:ILE:CD1	1:B:576:LEU:HB3	2.49	0.42
1:B:438:PHE:CG	1:B:443:LEU:HD11	2.54	0.42
1:B:438:PHE:HE2	1:B:461:ILE:HG12	1.84	0.42
1:A:295:GLU:HA	1:A:296:PRO:HA	1.87	0.42
1:B:792:VAL:HG11	1:B:830:HIS:CE1	2.54	0.42
1:A:600:THR:HG23	1:A:601:LEU:HD13	1.99	0.42
1:A:720:PRO:HG2	1:A:721:TYR:CE2	2.54	0.42
1:B:704:TYR:HB3	1:B:708:ASP:HB2	2.01	0.42
1:A:272:TYR:CE2	1:A:298:ALA:HB2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:439:ARG:O	1:A:440:GLU:C	2.58	0.42
1:B:436:LEU:HB3	1:B:438:PHE:CE1	2.55	0.42
1:A:252:LEU:O	1:A:255:HIS:HB2	2.18	0.42
1:A:656:LEU:HD11	1:A:674:LEU:HB2	2.02	0.42
1:A:161:PHE:HA	1:A:162:PRO:HD3	1.96	0.42
1:A:802:GLN:O	1:A:806:GLN:HG2	2.20	0.42
1:B:178:GLU:O	1:B:194:VAL:HA	2.20	0.42
1:B:180:SER:HB3	1:B:193:SER:HB3	2.02	0.42
1:A:694:ILE:O	1:A:695:ASN:HB2	2.19	0.42
1:A:964:GLN:HG2	1:A:999:CYS:HB3	2.01	0.42
1:A:461:ILE:O	1:A:465:GLU:HG2	2.19	0.42
1:B:759:GLU:O	1:B:802:GLN:NE2	2.42	0.42
1:A:762:THR:CG2	1:A:815:SER:OG	2.68	0.41
1:A:403:PHE:CB	1:A:501:ILE:HD11	2.50	0.41
1:B:720:PRO:HB2	1:B:728:ARG:HD3	2.01	0.41
1:B:431:GLU:HG3	1:B:468:HIS:HB3	2.01	0.41
1:A:479:TRP:CG	1:A:480:TRP:N	2.89	0.41
1:A:412:TYR:CZ	1:A:417:LEU:HD13	2.55	0.41
1:A:315:PHE:N	1:A:351:SER:OG	2.53	0.41
1:A:331:MET:HE3	1:A:351:SER:HB2	2.01	0.41
1:B:874:TRP:CE2	1:B:905:LYS:HD3	2.55	0.41
1:B:787:LEU:HD12	1:B:787:LEU:HA	1.93	0.41
1:B:295:GLU:HG3	1:B:480:TRP:CH2	2.56	0.41
1:A:277:PHE:HZ	1:A:423:PRO:HD2	1.86	0.41
1:B:727:ASP:O	1:B:730:ASN:HB3	2.21	0.41
1:A:746:LEU:O	1:A:749:ALA:HB3	2.21	0.41
1:B:465:GLU:O	1:B:468:HIS:HB2	2.20	0.41
1:B:356:THR:O	1:B:359:VAL:HG23	2.21	0.41
1:B:980:LEU:HD13	1:B:1007:ILE:HB	2.02	0.41
1:A:551:LYS:O	1:A:555:LEU:HG	2.20	0.41
1:B:203:VAL:HA	1:B:251:LEU:O	2.21	0.41
1:B:194:VAL:HG23	1:B:263:TYR:HE2	1.86	0.41
1:B:397:THR:HG22	1:B:462:ILE:CG2	2.50	0.41
1:B:618:LEU:HD11	1:B:703:HIS:HB2	2.03	0.41
1:B:942:TRP:O	1:B:946:LYS:HG3	2.21	0.41
1:A:438:PHE:HE1	1:A:461:ILE:HG12	1.86	0.41
1:A:422:ILE:HA	1:A:423:PRO:HD3	1.95	0.41
1:B:276:GLY:HA2	1:B:290:ALA:HA	2.03	0.40
1:B:512:LEU:HG	1:B:516:PHE:CE2	2.56	0.40
1:A:168:LEU:HD13	1:A:306:PHE:CD2	2.53	0.40
1:A:847:MET:HA	1:A:876:PHE:CE2	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:372:ASP:HA	1:A:377:LEU:HA	2.03	0.40
1:B:714:HIS:CE1	1:B:718:ILE:HG13	2.57	0.40
1:A:597:THR:HB	1:A:600:THR:HG22	2.03	0.40
1:B:485:LEU:HD22	1:B:586:SER:HA	2.04	0.40
1:B:661:GLU:OE1	1:B:722:VAL:HG21	2.22	0.40
1:A:534:VAL:HG13	1:A:539:GLN:HB3	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	856/872 (98%)	800 (94%)	48 (6%)	8 (1%)	21 61
1	B	841/872 (96%)	787 (94%)	52 (6%)	2 (0%)	52 86
All	All	1697/1744 (97%)	1587 (94%)	100 (6%)	10 (1%)	30 70

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	226	SER
1	A	638	ASN
1	A	664	ASN
1	A	282	GLU
1	A	227	SER
1	A	868	ALA
1	A	340	ASP
1	B	183	PRO
1	A	296	PRO
1	B	296	PRO

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	765/781 (98%)	743 (97%)	22 (3%)	50	82
1	B	751/781 (96%)	731 (97%)	20 (3%)	52	84
All	All	1516/1562 (97%)	1474 (97%)	42 (3%)	51	83

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	236	GLU
1	A	319	ILE
1	A	322	ASP
1	A	383	VAL
1	A	405	GLN
1	A	416	LYS
1	A	420	VAL
1	A	433	TRP
1	A	455	ARG
1	A	477	MET
1	A	481	ASN
1	A	489	PHE
1	A	563	LEU
1	A	634	ARG
1	A	665	TYR
1	A	684	THR
1	A	730	ASN
1	A	762	THR
1	A	785	MET
1	A	904	ARG
1	A	1016	LYS
1	A	1022	THR
1	B	206	ASN
1	B	260	LYS
1	B	280	THR
1	B	350	GLU
1	B	361	PHE

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Mol	Chain	Res	Type
1	B	433	TRP
1	B	477	MET
1	B	489	PHE
1	B	537	SER
1	B	601	LEU
1	B	605	ARG
1	B	634	ARG
1	B	637	LEU
1	B	706	ASP
1	B	782	LEU
1	B	792	VAL
1	B	828	CYS
1	B	849	SER
1	B	933	ARG
1	B	1022	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

17 non-standard protein/DNA/RNA residues 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	A	1103	1	14,14,15	0.76	1 (7%)	15,19,21	1.06	2 (13%)
4	NAG	A	1104	1	14,14,15	0.54	0	15,19,21	1.28	2 (13%)
4	NAG	A	1105	1	14,14,15	0.67	0	15,19,21	1.59	1 (6%)
4	NAG	A	1106	1	14,14,15	0.69	0	15,19,21	1.44	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	A	1107	1	14,14,15	0.74	0	15,19,21	2.03	1 (6%)
4	NAG	A	1108	1	14,14,15	0.68	0	15,19,21	1.60	4 (26%)
4	NAG	A	1109	1	14,14,15	1.00	1 (7%)	15,19,21	1.93	1 (6%)
4	NAG	B	1103	1	14,14,15	0.82	0	15,19,21	1.72	2 (13%)
5	NAG	B	1104	1,5	14,14,15	0.78	0	15,19,21	1.96	3 (20%)
5	NAG	B	1105	5	14,14,15	0.63	0	15,19,21	1.04	1 (6%)
4	NAG	B	1106	1	14,14,15	0.52	0	15,19,21	1.77	1 (6%)
4	NAG	B	1107	1	14,14,15	0.56	0	15,19,21	1.66	1 (6%)
4	NAG	B	1108	1	14,14,15	0.80	0	15,19,21	1.55	2 (13%)
4	NAG	B	1109	1	14,14,15	0.61	0	15,19,21	1.93	1 (6%)
4	NAG	B	1110	1	14,14,15	0.56	0	15,19,21	0.97	1 (6%)
4	NAG	B	1111	1	14,14,15	0.52	0	15,19,21	1.32	1 (6%)
4	NAG	B	1112	1	14,14,15	0.56	0	15,19,21	1.04	2 (13%)

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	A	1103	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1104	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1105	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1106	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1107	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1108	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1109	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1103	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1104	1,5	-	0/6/23/26	0/1/1/1
5	NAG	B	1105	5	-	0/6/23/26	0/1/1/1
4	NAG	B	1106	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1107	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1108	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1109	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1110	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1111	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1112	1	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1103	NAG	C1-C2	2.15	1.55	1.52
4	A	1109	NAG	C1-C2	2.86	1.56	1.52

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1108	NAG	C6-C5-C4	-2.63	106.52	113.02
4	A	1108	NAG	O7-C7-C8	-2.13	118.15	122.06
4	B	1103	NAG	O7-C7-C8	-2.08	118.24	122.06
5	B	1104	NAG	O7-C7-C8	-2.03	118.33	122.06
4	B	1112	NAG	O5-C5-C6	2.09	111.88	107.35
4	A	1103	NAG	C1-O5-C5	2.10	114.92	112.25
4	A	1103	NAG	C2-N2-C7	2.15	125.81	123.04
5	B	1104	NAG	O7-C7-N2	2.16	126.27	121.86
4	A	1108	NAG	C2-N2-C7	2.21	125.88	123.04
4	B	1112	NAG	C1-O5-C5	2.22	115.07	112.25
4	B	1108	NAG	C2-N2-C7	2.32	126.02	123.04
4	A	1104	NAG	C1-O5-C5	2.53	115.46	112.25
5	B	1105	NAG	C1-O5-C5	2.59	115.53	112.25
4	B	1110	NAG	C1-O5-C5	2.79	115.79	112.25
4	A	1104	NAG	C4-C3-C2	3.02	115.92	111.23
4	B	1111	NAG	C1-O5-C5	3.93	117.23	112.25
4	A	1108	NAG	C1-O5-C5	3.99	117.32	112.25
4	A	1106	NAG	C1-O5-C5	4.20	117.57	112.25
4	B	1108	NAG	C1-O5-C5	4.40	117.83	112.25
4	B	1103	NAG	C1-O5-C5	5.06	118.67	112.25
4	A	1105	NAG	C1-O5-C5	5.39	119.09	112.25
4	B	1107	NAG	C1-O5-C5	5.50	119.23	112.25
4	B	1106	NAG	C1-O5-C5	5.73	119.52	112.25
5	B	1104	NAG	C1-O5-C5	5.83	119.64	112.25
4	A	1109	NAG	C1-O5-C5	6.10	119.98	112.25
4	B	1109	NAG	C1-O5-C5	6.65	120.69	112.25
4	A	1107	NAG	C1-O5-C5	7.00	121.13	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1109	NAG	1	0

5.5 Carbohydrates ⓘ

2 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	$\# Z > 2$
5	NAG	B	1104	1,5	14,14,15	0.78	0	15,19,21	1.96	3 (20%)
5	NAG	B	1105	5	14,14,15	0.63	0	15,19,21	1.04	1 (6%)

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
5	NAG	B	1104	1,5	-	0/6/23/26	0/1/1/1
5	NAG	B	1105	5	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1104	NAG	O7-C7-C8	-2.03	118.33	122.06
5	B	1104	NAG	O7-C7-N2	2.16	126.27	121.86
5	B	1105	NAG	C1-O5-C5	2.59	115.53	112.25
5	B	1104	NAG	C1-O5-C5	5.83	119.64	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.6 Ligand geometry

Of 19 ligands modelled in this entry, 2 are monoatomic - leaving 17 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
3	LYS	A	1102	2	7,8,9	0.42	0	6,8,10	1.01	1 (16%)
4	NAG	A	1103	1	14,14,15	0.76	1 (7%)	15,19,21	1.06	2 (13%)
4	NAG	A	1104	1	14,14,15	0.54	0	15,19,21	1.28	2 (13%)
4	NAG	A	1105	1	14,14,15	0.67	0	15,19,21	1.59	1 (6%)
4	NAG	A	1106	1	14,14,15	0.69	0	15,19,21	1.44	1 (6%)
4	NAG	A	1107	1	14,14,15	0.74	0	15,19,21	2.03	1 (6%)
4	NAG	A	1108	1	14,14,15	0.68	0	15,19,21	1.60	4 (26%)
4	NAG	A	1109	1	14,14,15	1.00	1 (7%)	15,19,21	1.93	1 (6%)
3	LYS	B	1102	-	7,8,9	0.69	0	6,8,10	1.23	1 (16%)
4	NAG	B	1103	1	14,14,15	0.82	0	15,19,21	1.72	2 (13%)
4	NAG	B	1106	1	14,14,15	0.52	0	15,19,21	1.77	1 (6%)
4	NAG	B	1107	1	14,14,15	0.56	0	15,19,21	1.66	1 (6%)
4	NAG	B	1108	1	14,14,15	0.80	0	15,19,21	1.55	2 (13%)
4	NAG	B	1109	1	14,14,15	0.61	0	15,19,21	1.93	1 (6%)
4	NAG	B	1110	1	14,14,15	0.56	0	15,19,21	0.97	1 (6%)
4	NAG	B	1111	1	14,14,15	0.52	0	15,19,21	1.32	1 (6%)
4	NAG	B	1112	1	14,14,15	0.56	0	15,19,21	1.04	2 (13%)

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	LYS	A	1102	2	-	0/5/7/9	0/0/0/0
4	NAG	A	1103	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1104	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1105	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	1106	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1107	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1108	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1109	1	-	0/6/23/26	0/1/1/1
3	LYS	B	1102	-	-	0/5/7/9	0/0/0/0
4	NAG	B	1103	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1106	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1107	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1108	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1109	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1110	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1111	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1112	1	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1103	NAG	C1-C2	2.15	1.55	1.52
4	A	1109	NAG	C1-C2	2.86	1.56	1.52

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1102	LYS	O-C-CA	-2.84	118.09	125.49
4	A	1108	NAG	C6-C5-C4	-2.63	106.52	113.02
3	A	1102	LYS	O-C-CA	-2.36	119.33	125.49
4	A	1108	NAG	O7-C7-C8	-2.13	118.15	122.06
4	B	1103	NAG	O7-C7-C8	-2.08	118.24	122.06
4	B	1112	NAG	O5-C5-C6	2.09	111.88	107.35
4	A	1103	NAG	C1-O5-C5	2.10	114.92	112.25
4	A	1103	NAG	C2-N2-C7	2.15	125.81	123.04
4	A	1108	NAG	C2-N2-C7	2.21	125.88	123.04
4	B	1112	NAG	C1-O5-C5	2.22	115.07	112.25
4	B	1108	NAG	C2-N2-C7	2.32	126.02	123.04
4	A	1104	NAG	C1-O5-C5	2.53	115.46	112.25
4	B	1110	NAG	C1-O5-C5	2.79	115.79	112.25
4	A	1104	NAG	C4-C3-C2	3.02	115.92	111.23
4	B	1111	NAG	C1-O5-C5	3.93	117.23	112.25
4	A	1108	NAG	C1-O5-C5	3.99	117.32	112.25
4	A	1106	NAG	C1-O5-C5	4.20	117.57	112.25
4	B	1108	NAG	C1-O5-C5	4.40	117.83	112.25
4	B	1103	NAG	C1-O5-C5	5.06	118.67	112.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1105	NAG	C1-O5-C5	5.39	119.09	112.25
4	B	1107	NAG	C1-O5-C5	5.50	119.23	112.25
4	B	1106	NAG	C1-O5-C5	5.73	119.52	112.25
4	A	1109	NAG	C1-O5-C5	6.10	119.98	112.25
4	B	1109	NAG	C1-O5-C5	6.65	120.69	112.25
4	A	1107	NAG	C1-O5-C5	7.00	121.13	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1102	LYS	1	0
4	B	1109	NAG	1	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	860/872 (98%)	-0.05	19 (2%)	65 44	23, 60, 96, 122	0
1	B	849/872 (97%)	0.16	39 (4%)	36 21	29, 70, 105, 139	0
All	All	1709/1744 (97%)	0.05	58 (3%)	49 30	23, 65, 101, 139	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	648	THR	5.0
1	B	229	GLU	4.6
1	B	221	PHE	4.5
1	B	205	TRP	4.2
1	B	244	ALA	4.0
1	B	246	VAL	4.0
1	B	207	ILE	4.0
1	B	203	VAL	3.6
1	B	340	ASP	3.5
1	B	628	LEU	3.5
1	A	206	ASN	3.3
1	B	161	PHE	3.2
1	B	284	ASN	3.2
1	A	244	ALA	3.1
1	B	245	ILE	3.1
1	A	235	LEU	3.0
1	A	992	GLU	3.0
1	A	340	ASP	3.0
1	B	249	GLU	3.0
1	B	220	THR	2.9
1	B	208	ILE	2.8
1	B	230	LYS	2.8
1	B	601	LEU	2.8
1	B	801	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	219	VAL	2.7
1	B	546	SER	2.6
1	B	251	LEU	2.6
1	B	247	ALA	2.6
1	B	668	TYR	2.5
1	B	206	ASN	2.5
1	A	204	THR	2.5
1	B	534	VAL	2.5
1	B	204	THR	2.5
1	A	343	LEU	2.5
1	B	233	GLU	2.5
1	A	201	LEU	2.4
1	B	283	SER	2.4
1	B	160	LEU	2.4
1	A	245	ILE	2.4
1	B	185	LEU	2.4
1	A	221	PHE	2.4
1	A	597	THR	2.3
1	A	203	VAL	2.3
1	A	159	LYS	2.3
1	A	233	GLU	2.3
1	B	659	VAL	2.3
1	B	243	ILE	2.3
1	B	961	TYR	2.2
1	B	163	TRP	2.2
1	B	222	MET	2.2
1	A	202	GLN	2.2
1	A	286	LYS	2.1
1	B	237	TYR	2.1
1	B	242	GLN	2.1
1	A	251	LEU	2.1
1	B	596	VAL	2.1
1	A	247	ALA	2.0
1	B	665	TYR	2.0

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

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	NAG	B	1106	14/15	0.86	0.28	4.22	96,104,111,118	0
4	NAG	B	1109	14/15	0.72	0.42	3.71	118,128,135,135	0
4	NAG	B	1103	14/15	0.73	0.36	1.29	102,119,123,127	0
4	NAG	A	1105	14/15	0.68	0.26	0.92	110,128,154,156	0
4	NAG	A	1104	14/15	0.91	0.17	0.84	101,111,114,116	0
4	NAG	A	1103	14/15	0.83	0.29	0.64	101,116,119,122	0
5	NAG	B	1104	14/15	0.90	0.16	-0.66	84,96,104,113	0
4	NAG	A	1109	14/15	0.87	0.26	-	99,106,118,118	0
4	NAG	B	1107	14/15	0.85	0.23	-	96,101,107,107	0
4	NAG	A	1107	14/15	0.90	0.10	-	86,93,100,106	0
4	NAG	A	1108	14/15	0.82	0.31	-	116,130,143,146	0
5	NAG	B	1105	14/15	0.79	0.28	-	105,117,131,131	0
4	NAG	B	1110	14/15	0.82	0.21	-	97,108,119,123	0
4	NAG	B	1111	14/15	0.89	0.29	-	85,93,98,103	0
4	NAG	A	1106	14/15	0.80	0.32	-	100,121,130,139	0
4	NAG	B	1108	14/15	0.82	0.30	-	131,142,151,153	0
4	NAG	B	1112	14/15	0.89	0.20	-	84,97,111,116	0

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
5	NAG	B	1104	14/15	0.90	0.16	-0.66	84,96,104,113	0
5	NAG	B	1105	14/15	0.79	0.28	-	105,117,131,131	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
3	LYS	A	1102	9/10	0.88	0.48	11.62	41,63,88,93	4
3	LYS	B	1102	9/10	0.88	0.53	10.92	74,77,114,123	5
4	NAG	B	1106	14/15	0.86	0.28	4.22	96,104,111,118	0
4	NAG	B	1109	14/15	0.72	0.42	3.71	118,128,135,135	0
4	NAG	B	1103	14/15	0.73	0.36	1.29	102,119,123,127	0
2	ZN	A	1101	1/1	0.99	0.23	1.08	52,52,52,52	0
4	NAG	A	1105	14/15	0.68	0.26	0.92	110,128,154,156	0
4	NAG	A	1104	14/15	0.91	0.17	0.84	101,111,114,116	0
4	NAG	A	1103	14/15	0.83	0.29	0.64	101,116,119,122	0
2	ZN	B	1101	1/1	0.99	0.20	-0.96	46,46,46,46	0
4	NAG	B	1110	14/15	0.82	0.21	-	97,108,119,123	0
4	NAG	B	1112	14/15	0.89	0.20	-	84,97,111,116	0
4	NAG	A	1108	14/15	0.82	0.31	-	116,130,143,146	0
4	NAG	B	1111	14/15	0.89	0.29	-	85,93,98,103	0
4	NAG	A	1109	14/15	0.87	0.26	-	99,106,118,118	0
4	NAG	A	1107	14/15	0.90	0.10	-	86,93,100,106	0
4	NAG	A	1106	14/15	0.80	0.32	-	100,121,130,139	0
4	NAG	B	1108	14/15	0.82	0.30	-	131,142,151,153	0
4	NAG	B	1107	14/15	0.85	0.23	-	96,101,107,107	0

6.5 Other polymers [i](#)

There are no such residues in this entry.