



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:26 AM UTC

PDB ID : 2XGM / pdb_00002xgm
Title : Substrate and product analogues as human O-GlcNAc transferase inhibitors.
Authors : Dorfmueller, H.C.; Borodkin, V.S.; Blair, D.E.; Pathak, S.; Navratilova, I.;
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Deposited on : 2010-06-07
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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

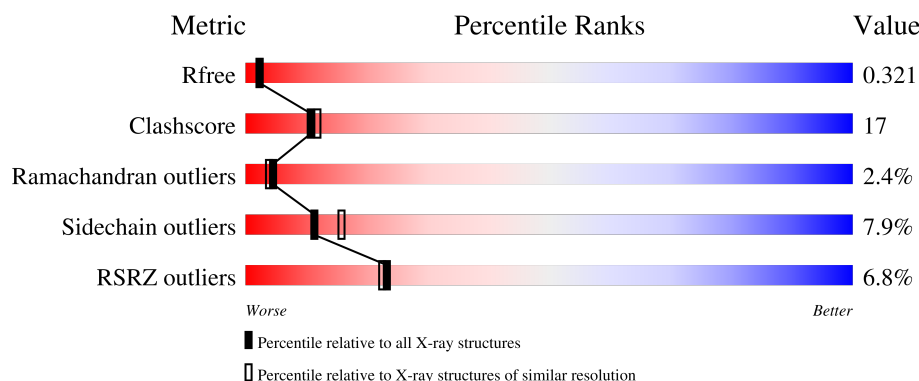
1 Overall quality at a glance

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}	180053	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	568	6% 65% 21% • 10%
1	B	568	7% 67% 24% • • 5%

2 Entry composition [i](#)

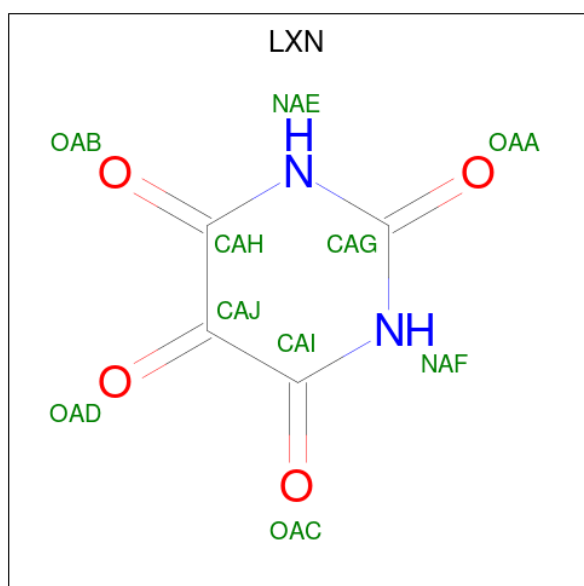
There are 3 unique types of molecules in this entry. The entry contains 8113 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 XCOGT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	512	Total	C	N	O	S	0	2	0
			3923	2480	730	700	13			
1	B	542	Total	C	N	O	S	0	0	0
			4112	2595	761	740	16			

- Molecule 2 is ALLOXAN (CCD ID: LXN) (formula: $C_4H_2N_2O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			10	4	2	4		
2	B	1	Total	C	N	O	0	0
			10	4	2	4		

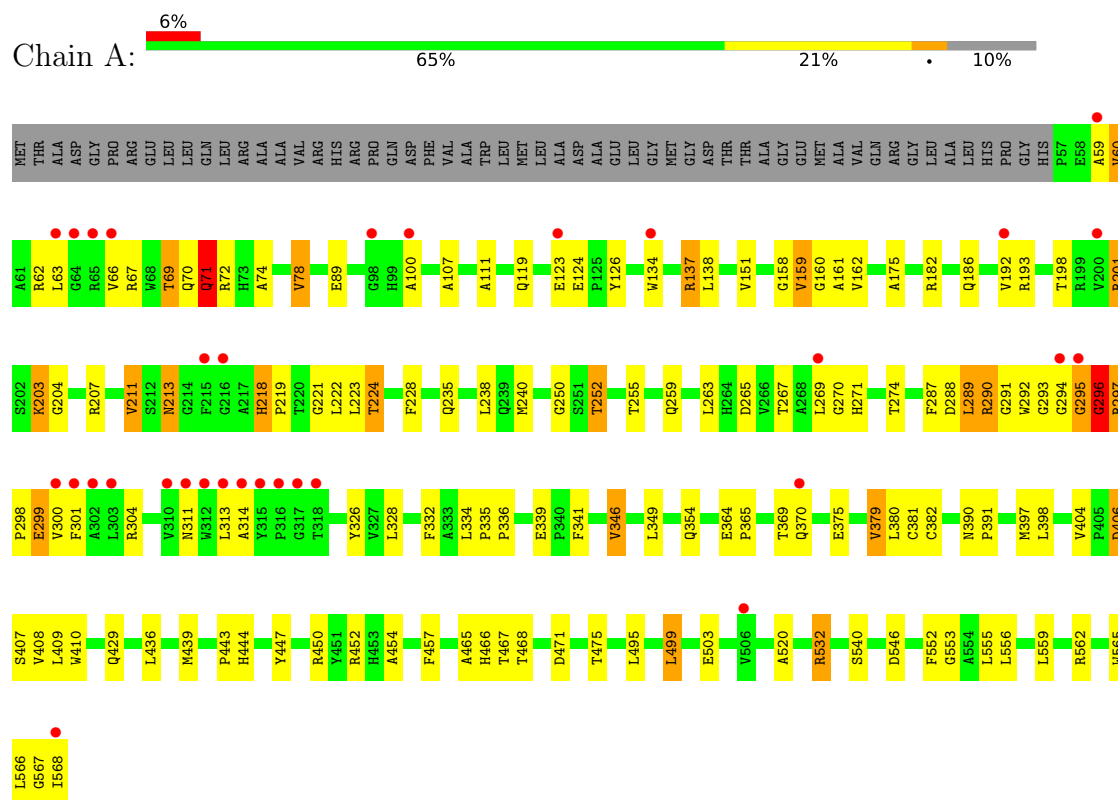
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	21	Total 21	O 21	0	0
3	B	37	Total 37	O 37	0	0

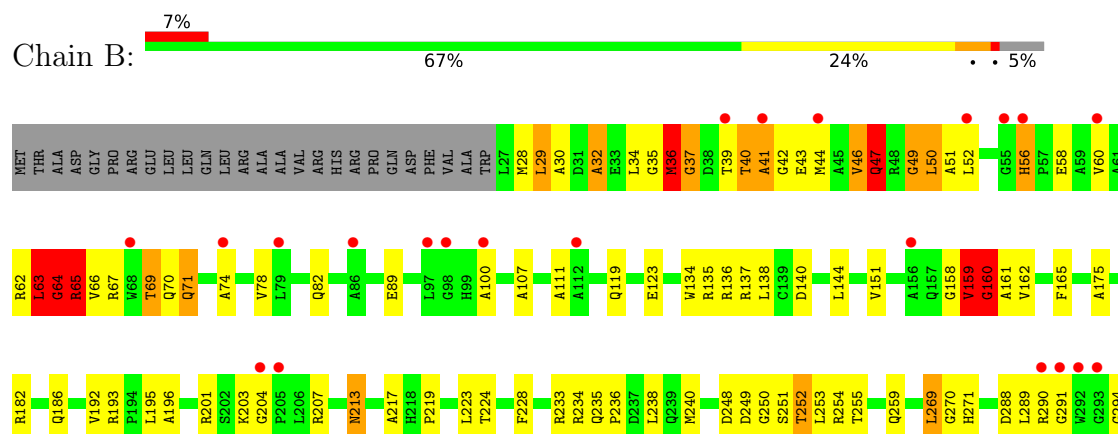
3 Residue-property plots [i](#)

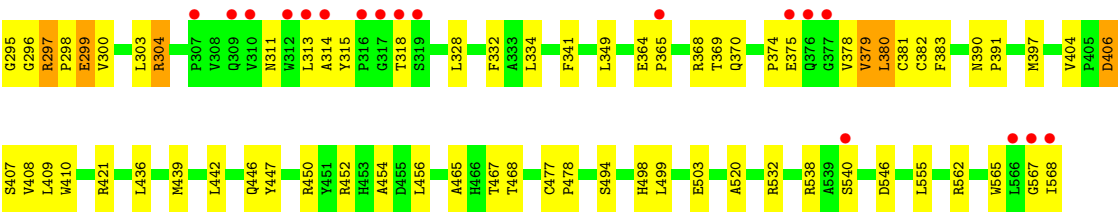
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: XCOGT



• Molecule 1: XCOGT





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.53Å 99.98Å 157.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.55 20.00 – 2.55	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.55) 97.5 (20.00-2.55)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.94 (at 2.56Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.267 , 0.309 0.278 , 0.321	Depositor DCC
R_{free} test set	1334 reflections (3.13%)	wwPDB-VP
Wilson B-factor (Å ²)	74.6	Xtriage
Anisotropy	0.227	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 10.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	8113	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.64% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.95	0/4031	1.12	14/5501 (0.3%)
1	B	0.81	0/4217	1.11	20/5753 (0.3%)
All	All	0.88	0/8248	1.11	34/11254 (0.3%)

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	3
1	B	0	5
All	All	0	8

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	56	HIS	CA-C-N	9.35	131.53	119.84
1	B	56	HIS	C-N-CA	9.35	131.53	119.84
1	A	160	GLY	CA-C-N	8.75	133.61	120.31
1	A	160	GLY	C-N-CA	8.75	133.61	120.31
1	B	63	LEU	N-CA-C	7.89	119.88	111.28
1	B	35	GLY	N-CA-C	7.13	122.92	113.37
1	B	203	LYS	CA-C-N	7.07	134.43	121.70
1	B	203	LYS	C-N-CA	7.07	134.43	121.70
1	B	269	LEU	CA-C-N	6.96	134.24	121.70
1	B	269	LEU	C-N-CA	6.96	134.24	121.70
1	A	218	HIS	CA-C-N	-6.81	111.33	119.84
1	A	218	HIS	C-N-CA	-6.81	111.33	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	46	VAL	N-CA-C	6.79	123.47	109.34
1	A	289	LEU	CA-C-N	-6.40	112.50	122.42
1	A	289	LEU	C-N-CA	-6.40	112.50	122.42
1	B	160	GLY	CA-C-N	6.35	133.13	121.70
1	B	160	GLY	C-N-CA	6.35	133.13	121.70
1	B	46	VAL	CA-C-N	6.17	132.81	121.70
1	B	46	VAL	C-N-CA	6.17	132.81	121.70
1	B	64	GLY	CA-C-N	6.11	133.20	121.54
1	B	64	GLY	C-N-CA	6.11	133.20	121.54
1	B	562	ARG	NE-CZ-NH1	-6.05	115.45	121.50
1	A	224	THR	N-CA-C	5.87	120.44	113.16
1	B	562	ARG	NE-CZ-NH2	5.82	124.44	119.20
1	B	36	MET	N-CA-C	5.72	122.98	110.80
1	B	406	ASP	N-CA-C	5.67	119.01	111.75
1	A	292	TRP	N-CA-C	-5.61	106.20	113.16
1	A	269	LEU	CA-C-N	5.53	128.60	121.13
1	A	269	LEU	C-N-CA	5.53	128.60	121.13
1	A	221	GLY	N-CA-C	-5.42	106.11	113.37
1	A	346	VAL	N-CA-C	5.37	116.47	108.73
1	B	65	ARG	N-CA-C	5.03	121.52	110.80
1	A	124	GLU	CA-C-N	5.00	124.60	119.05
1	A	124	GLU	C-N-CA	5.00	124.60	119.05

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	203	LYS	Peptide
1	A	296	GLY	Peptide
1	A	406	ASP	Peptide
1	B	160	GLY	Peptide
1	B	296	GLY	Peptide
1	B	406	ASP	Peptide
1	B	63	LEU	Peptide
1	B	64	GLY	Peptide

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	3923	0	3855	127	0
1	B	4112	0	4043	145	0
2	A	10	0	2	0	0
2	B	10	0	2	0	0
3	A	21	0	0	5	0
3	B	37	0	0	12	0
All	All	8113	0	7902	266	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 17.

All (266) 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:442:LEU:HB3	3:B:2031:HOH:O	1.31	1.26
1:A:67:ARG:HD2	3:A:2001:HOH:O	1.44	1.15
1:B:382:CYS:HB2	1:B:397:MET:HE1	1.34	1.05
1:A:439:MET:HE1	1:A:447:TYR:HD1	1.22	1.05
1:B:382:CYS:CB	1:B:397:MET:HE1	1.89	1.01
1:B:40:THR:HA	1:B:41:ALA:HB3	1.44	0.99
1:B:382:CYS:HB2	1:B:397:MET:CE	1.93	0.98
1:A:290:ARG:NE	1:A:297:ARG:H	1.62	0.97
1:A:290:ARG:HE	1:A:297:ARG:N	1.63	0.96
1:B:383:PHE:CE1	3:B:2027:HOH:O	2.19	0.95
1:A:290:ARG:HE	1:A:297:ARG:H	0.97	0.95
1:B:439:MET:HE1	1:B:447:TYR:HD1	1.31	0.95
1:B:446:GLN:HB2	3:B:2031:HOH:O	1.66	0.94
1:A:354:GLN:HE21	1:A:466[A]:HIS:CE1	1.87	0.93
1:B:328:LEU:HD11	1:B:349:LEU:HG	1.52	0.92
1:B:43:GLU:C	1:B:47:GLN:HE22	1.79	0.90
1:A:158:GLY:HA2	1:A:159:VAL:O	1.71	0.90
1:A:382:CYS:CB	1:A:397:MET:HE1	2.03	0.88
1:A:406:ASP:HB2	3:A:2015:HOH:O	1.75	0.86
1:A:354:GLN:NE2	1:A:466[A]:HIS:CE1	2.43	0.86
1:A:382:CYS:HB2	1:A:397:MET:CE	2.05	0.86
1:A:328:LEU:HD11	1:A:349:LEU:HG	1.59	0.84
1:B:223:LEU:HD11	1:B:467:THR:HG22	1.59	0.83
1:A:67:ARG:NH1	1:A:70:GLN:OE1	2.10	0.83
1:B:213:ASN:HD22	1:B:213:ASN:H	1.26	0.83
1:A:223:LEU:HD21	1:A:467:THR:HG22	1.60	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:GLY:O	1:A:296:GLY:O	2.00	0.80
1:A:354:GLN:NE2	1:A:466[A]:HIS:ND1	2.30	0.79
1:B:151:VAL:HG13	1:B:162:VAL:HG22	1.65	0.78
1:A:382:CYS:SG	1:A:397:MET:HE1	2.23	0.78
1:B:378:VAL:HB	3:B:2025:HOH:O	1.84	0.76
1:B:151:VAL:HG13	1:B:162:VAL:CG2	2.16	0.75
1:A:382:CYS:HB2	1:A:397:MET:HE1	1.65	0.74
1:B:288:ASP:CG	1:B:290:ARG:HH21	1.96	0.74
1:A:159:VAL:O	1:A:159:VAL:HG23	1.86	0.74
1:A:439:MET:HE1	1:A:447:TYR:CD1	2.15	0.73
1:A:213:ASN:H	1:A:213:ASN:HD22	1.37	0.73
1:B:290:ARG:HH11	1:B:297:ARG:H	1.38	0.72
1:B:381:CYS:HB2	3:B:2027:HOH:O	1.89	0.72
1:A:159:VAL:O	1:A:159:VAL:CG2	2.38	0.71
1:B:158:GLY:HA2	1:B:159:VAL:O	1.90	0.71
1:A:288:ASP:CG	1:A:290:ARG:HH21	1.97	0.71
1:A:290:ARG:HH11	1:A:298:PRO:HD3	1.56	0.70
1:A:568:ILE:HG22	1:A:568:ILE:O	1.90	0.70
1:A:252:THR:HA	1:B:252:THR:HG22	1.73	0.70
1:B:439:MET:HE1	1:B:447:TYR:CD1	2.22	0.70
1:B:368:ARG:HB3	3:B:2024:HOH:O	1.92	0.69
1:B:63:LEU:O	1:B:66:VAL:HB	1.93	0.68
1:B:40:THR:HG21	1:B:70:GLN:HE21	1.59	0.68
1:A:382:CYS:HB2	1:A:397:MET:HE2	1.74	0.68
1:B:382:CYS:SG	1:B:397:MET:HE1	2.34	0.67
1:B:40:THR:CA	1:B:41:ALA:HB3	2.22	0.67
1:B:238:LEU:HB3	1:B:240:MET:HE2	1.76	0.67
1:B:111:ALA:HB2	1:B:134:TRP:HD1	1.58	0.67
1:A:235:GLN:HE21	1:A:235:GLN:HA	1.59	0.67
1:B:46:VAL:N	1:B:47:GLN:O	2.28	0.66
1:B:290:ARG:HH11	1:B:298:PRO:HD3	1.60	0.66
1:B:56:HIS:HD2	1:B:58:GLU:H	1.44	0.65
1:A:158:GLY:HA2	1:A:159:VAL:C	2.19	0.64
1:B:297:ARG:H	1:B:298:PRO:HD3	1.63	0.64
1:B:288:ASP:OD2	1:B:290:ARG:NH2	2.31	0.63
1:B:201:ARG:NH1	1:B:207:ARG:HE	1.96	0.63
1:B:313:LEU:O	1:B:313:LEU:HG	1.97	0.63
1:A:182:ARG:HH21	1:A:186:GLN:HE22	1.47	0.63
1:A:111:ALA:HB2	1:A:134:TRP:HD1	1.63	0.62
1:B:47:GLN:O	1:B:47:GLN:CD	2.42	0.62
1:B:568:ILE:O	1:B:568:ILE:HG22	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:ARG:HA	1:B:67:ARG:HE	1.65	0.60
1:A:290:ARG:NH1	1:A:298:PRO:HD3	2.15	0.60
1:B:46:VAL:N	1:B:47:GLN:C	2.60	0.60
1:B:56:HIS:CD2	1:B:58:GLU:HB2	2.36	0.60
1:B:43:GLU:HA	1:B:47:GLN:OE1	2.01	0.60
1:A:252:THR:HG22	1:B:252:THR:HA	1.82	0.59
1:B:288:ASP:CG	1:B:290:ARG:NH2	2.61	0.59
1:A:298:PRO:HD2	1:A:299:GLU:OE1	2.03	0.59
1:B:503:GLU:OE2	1:B:532:ARG:NH1	2.33	0.58
1:B:297:ARG:H	1:B:298:PRO:CD	2.17	0.58
1:B:290:ARG:HB2	1:B:290:ARG:CZ	2.33	0.58
1:B:43:GLU:HA	1:B:47:GLN:CD	2.29	0.57
1:B:158:GLY:HA2	1:B:159:VAL:C	2.29	0.57
1:B:204:GLY:O	1:B:565:TRP:HZ2	1.86	0.57
1:A:235:GLN:HA	1:A:235:GLN:NE2	2.19	0.57
1:A:204:GLY:O	1:A:565:TRP:HZ2	1.87	0.57
1:A:290:ARG:CZ	1:A:297:ARG:H	2.17	0.57
1:B:235:GLN:HE21	1:B:235:GLN:HA	1.70	0.56
1:B:47:GLN:O	1:B:47:GLN:NE2	2.39	0.56
1:B:410:TRP:C	3:B:2027:HOH:O	2.48	0.56
1:A:288:ASP:OD2	1:A:290:ARG:NH2	2.39	0.56
1:B:43:GLU:HA	1:B:47:GLN:NE2	2.20	0.56
1:B:43:GLU:HG3	1:B:47:GLN:NE2	2.21	0.56
1:B:379:VAL:HG22	1:B:454:ALA:HA	1.88	0.56
1:A:555:LEU:HD23	1:A:555:LEU:C	2.31	0.56
1:B:43:GLU:CA	1:B:47:GLN:HE22	2.18	0.55
1:A:67:ARG:HE	1:A:67:ARG:HA	1.72	0.55
1:A:67:ARG:HH11	1:A:70:GLN:CD	2.11	0.55
1:A:204:GLY:O	1:A:565:TRP:CZ2	2.60	0.55
1:B:134:TRP:CZ3	1:B:137:ARG:HD2	2.41	0.55
1:A:406:ASP:CB	3:A:2015:HOH:O	2.45	0.54
1:A:290:ARG:CZ	1:A:290:ARG:HB2	2.37	0.54
1:A:379:VAL:HG22	1:A:454:ALA:HA	1.89	0.54
1:A:175:ALA:HB2	1:A:341:PHE:CE1	2.43	0.53
1:A:218:HIS:O	1:A:222:LEU:HB3	2.08	0.53
1:B:565:TRP:NE1	3:B:2036:HOH:O	2.40	0.53
1:A:288:ASP:HB2	1:A:301:PHE:CZ	2.43	0.53
1:A:354:GLN:HE21	1:A:466[A]:HIS:HE1	1.46	0.53
1:B:56:HIS:CD2	1:B:58:GLU:H	2.25	0.53
1:B:269:LEU:O	1:B:297:ARG:NH1	2.37	0.53
1:B:298:PRO:HD2	1:B:299:GLU:OE1	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:ALA:O	1:A:78:VAL:HG23	2.09	0.52
1:A:365:PRO:HG3	1:A:452:ARG:HB2	1.90	0.52
1:B:56:HIS:HD2	1:B:58:GLU:HB2	1.74	0.52
1:B:204:GLY:O	1:B:565:TRP:CZ2	2.61	0.52
1:A:201:ARG:NH1	1:A:207:ARG:HE	2.07	0.52
1:B:67:ARG:NH1	1:B:70:GLN:OE1	2.39	0.52
1:B:288:ASP:OD1	1:B:290:ARG:NH2	2.38	0.52
1:A:213:ASN:H	1:A:213:ASN:ND2	2.05	0.52
1:A:62:ARG:O	1:A:66:VAL:HG23	2.09	0.51
1:A:398:LEU:HD12	1:A:429:GLN:HG3	1.92	0.51
1:B:555:LEU:C	1:B:555:LEU:HD23	2.35	0.51
1:B:36:MET:H	1:B:37:GLY:CA	2.23	0.51
1:A:126:TYR:HA	1:A:161:ALA:HB1	1.91	0.51
1:A:568:ILE:O	1:A:568:ILE:CG2	2.59	0.51
1:B:40:THR:HA	1:B:41:ALA:CB	2.22	0.51
1:A:151:VAL:HG13	1:A:162:VAL:HG22	1.93	0.51
1:B:303:LEU:O	1:B:304:ARG:C	2.54	0.51
1:B:74:ALA:O	1:B:78:VAL:HG23	2.11	0.51
1:B:62:ARG:O	1:B:65:ARG:HB3	2.11	0.51
1:B:297:ARG:N	1:B:298:PRO:CD	2.73	0.50
1:B:160:GLY:HA2	1:B:161:ALA:HB3	1.93	0.50
1:B:151:VAL:HG13	1:B:162:VAL:HG21	1.92	0.50
1:B:313:LEU:O	1:B:313:LEU:CG	2.56	0.50
1:B:538:ARG:O	1:B:538:ARG:HG2	2.12	0.50
1:A:213:ASN:HD22	1:A:213:ASN:N	2.01	0.50
1:B:369:THR:OG1	1:B:375:GLU:OE2	2.28	0.49
1:A:439:MET:SD	1:A:450:ARG:HG3	2.53	0.49
1:A:252:THR:HG21	1:B:255:THR:OG1	2.12	0.49
1:B:365:PRO:HG3	1:B:452:ARG:HB2	1.94	0.49
1:A:404:VAL:HG11	1:A:520:ALA:HB1	1.94	0.49
1:A:409:LEU:HD23	1:A:436:LEU:HD22	1.94	0.49
1:B:270:GLY:O	1:B:271:HIS:C	2.55	0.49
1:A:158:GLY:CA	1:A:159:VAL:C	2.86	0.49
1:A:290:ARG:NH1	1:A:298:PRO:CD	2.76	0.49
1:B:290:ARG:NH1	1:B:298:PRO:HD3	2.25	0.49
1:A:203:LYS:HD3	3:A:2021:HOH:O	2.13	0.49
1:B:175:ALA:HB2	1:B:341:PHE:CE1	2.47	0.49
1:B:271:HIS:ND1	1:B:299:GLU:HG3	2.28	0.49
1:B:271:HIS:CE1	1:B:299:GLU:HG3	2.48	0.48
1:B:439:MET:SD	1:B:450:ARG:HG3	2.53	0.48
1:B:290:ARG:HE	1:B:297:ARG:N	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:390:ASN:HB2	1:B:391:PRO:HD2	1.94	0.48
1:A:471:ASP:O	1:A:475:THR:HG23	2.13	0.48
1:A:274:THR:OG1	1:A:297:ARG:NH2	2.43	0.48
1:A:252:THR:HG23	1:B:250:GLY:O	2.14	0.48
1:A:332:PHE:C	1:A:332:PHE:CD2	2.92	0.48
1:A:503:GLU:OE2	1:A:532:ARG:NH1	2.39	0.48
1:B:50:LEU:HB3	3:B:2001:HOH:O	2.14	0.48
1:B:62:ARG:C	1:B:64:GLY:HA3	2.39	0.48
1:A:63:LEU:O	1:A:66:VAL:HB	2.13	0.48
1:B:136:ARG:HH12	1:B:144:LEU:HD21	1.79	0.48
1:A:288:ASP:CG	1:A:290:ARG:NH2	2.69	0.48
1:B:28:MET:C	1:B:30:ALA:H	2.22	0.48
1:B:46:VAL:HB	1:B:47:GLN:HG3	1.94	0.48
1:B:235:GLN:HA	1:B:235:GLN:NE2	2.28	0.48
1:A:290:ARG:NH1	1:A:291:GLY:O	2.47	0.47
1:A:313:LEU:O	1:A:313:LEU:HG	2.14	0.47
1:B:248:ASP:OD1	1:B:254:ARG:HG2	2.14	0.47
1:A:297:ARG:O	1:A:300:VAL:HB	2.14	0.47
1:A:134:TRP:CZ3	1:A:137:ARG:HD2	2.50	0.47
1:B:304:ARG:HE	1:B:304:ARG:HB3	1.41	0.47
1:B:404:VAL:HG11	1:B:520:ALA:HB1	1.96	0.47
1:A:235:GLN:HG3	1:A:553:GLY:HA3	1.95	0.47
1:B:42:GLY:C	1:B:44:MET:H	2.23	0.47
1:A:70:GLN:O	1:A:72:ARG:N	2.48	0.47
1:B:62:ARG:O	1:B:66:VAL:HG23	2.15	0.47
1:B:234:ARG:O	1:B:236:PRO:HD3	2.15	0.47
1:B:290:ARG:NH1	1:B:291:GLY:O	2.48	0.47
1:B:410:TRP:HB3	3:B:2027:HOH:O	2.14	0.47
1:B:39:THR:HB	1:B:40:THR:HB	1.97	0.47
1:A:223:LEU:HD11	1:A:467:THR:CG2	2.45	0.46
1:B:213:ASN:HD22	1:B:213:ASN:N	2.00	0.46
1:A:465:ALA:HB1	1:A:468:THR:OG1	2.15	0.46
1:B:290:ARG:NH2	1:B:290:ARG:HB2	2.30	0.46
1:B:565:TRP:CZ3	1:B:567:GLY:HA2	2.51	0.46
1:A:250:GLY:O	1:B:252:THR:HG23	2.14	0.46
1:A:552:PHE:CE2	1:A:556:LEU:HD11	2.50	0.46
1:A:201:ARG:O	1:A:565:TRP:HD1	1.98	0.46
1:A:390:ASN:HB2	1:A:391:PRO:HD2	1.98	0.46
1:B:364:GLU:OE2	1:B:365:PRO:HD2	2.16	0.46
1:A:218:HIS:O	1:A:222:LEU:CB	2.63	0.45
1:B:182:ARG:HH21	1:B:186:GLN:HE22	1.63	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:380:LEU:HA	1:B:456:LEU:O	2.16	0.45
1:B:409:LEU:HD23	1:B:436:LEU:HD22	1.98	0.45
1:A:137:ARG:HG2	1:A:137:ARG:HH21	1.82	0.45
1:B:43:GLU:CA	1:B:47:GLN:NE2	2.79	0.45
1:B:135:ARG:HB3	1:B:140:ASP:O	2.17	0.45
1:A:69:THR:O	1:A:71:GLN:NE2	2.50	0.45
1:A:457:PHE:C	1:A:457:PHE:CD2	2.94	0.45
1:B:67:ARG:HH11	1:B:70:GLN:CD	2.23	0.45
1:A:495:LEU:O	1:A:499:LEU:HB2	2.17	0.45
1:B:165:PHE:CD1	1:B:165:PHE:C	2.95	0.45
1:B:538:ARG:HB3	1:B:538:ARG:NH2	2.31	0.45
1:A:224:THR:HB	1:A:228:PHE:CE2	2.51	0.44
1:B:67:ARG:HA	1:B:67:ARG:NE	2.32	0.44
1:A:67:ARG:HA	1:A:67:ARG:NE	2.32	0.44
1:B:40:THR:CA	1:B:41:ALA:CB	2.91	0.44
1:B:297:ARG:O	1:B:300:VAL:HB	2.18	0.44
1:A:364:GLU:OE2	1:A:364:GLU:HA	2.17	0.44
1:A:265:ASP:OD1	1:A:267:THR:HB	2.17	0.44
1:A:335:PRO:O	1:A:336:PRO:C	2.60	0.44
1:B:36:MET:H	1:B:37:GLY:HA2	1.83	0.44
1:A:369:THR:OG1	1:A:375:GLU:OE2	2.36	0.43
1:A:271:HIS:CD2	1:A:297:ARG:HG2	2.53	0.43
1:B:49:GLY:C	1:B:51:ALA:N	2.76	0.43
1:B:374:PRO:C	3:B:2024:HOH:O	2.60	0.43
1:B:36:MET:N	1:B:37:GLY:CA	2.81	0.43
1:A:290:ARG:HH11	1:A:297:ARG:N	2.16	0.43
1:A:293:GLY:C	3:A:2008:HOH:O	2.62	0.43
1:B:494:SER:O	1:B:498:HIS:ND1	2.50	0.43
1:A:151:VAL:HG13	1:A:162:VAL:CG2	2.48	0.43
1:A:364:GLU:OE2	1:A:365:PRO:HD2	2.18	0.43
1:B:43:GLU:HA	1:B:47:GLN:HE22	1.81	0.43
1:B:217:ALA:HA	1:B:253:LEU:HD12	2.00	0.43
1:A:218:HIS:O	1:A:219:PRO:C	2.60	0.43
1:A:297:ARG:HB3	1:A:300:VAL:CG2	2.48	0.43
1:B:465:ALA:HB1	1:B:468:THR:OG1	2.19	0.43
1:A:263:LEU:C	1:A:263:LEU:HD23	2.44	0.43
1:A:313:LEU:O	1:A:313:LEU:CG	2.63	0.43
1:A:381:CYS:HB2	1:A:410:TRP:HB3	2.00	0.43
1:A:443:PRO:O	1:A:444:HIS:C	2.61	0.43
1:A:107:ALA:HB1	1:A:138:LEU:HD21	2.00	0.42
1:A:339:GLU:HG2	1:A:346:VAL:HG21	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:34:LEU:HD23	1:B:69:THR:HG21	2.01	0.42
1:B:49:GLY:O	1:B:51:ALA:N	2.52	0.42
1:A:222:LEU:HD12	1:A:222:LEU:O	2.19	0.42
1:B:138:LEU:O	1:B:421:ARG:NH2	2.51	0.42
1:B:290:ARG:HB2	1:B:291:GLY:HA2	2.02	0.42
1:A:59:ALA:O	1:A:60:VAL:C	2.62	0.42
1:B:219:PRO:HB3	1:B:467:THR:HG21	2.00	0.42
1:A:238:LEU:HB3	1:A:240:MET:HE2	2.01	0.42
1:A:287:PHE:CZ	1:A:556:LEU:HD21	2.55	0.42
1:B:201:ARG:O	1:B:565:TRP:HD1	2.02	0.42
1:B:249:ASP:OD2	1:B:251:SER:OG	2.30	0.42
1:A:565:TRP:CH2	1:A:567:GLY:HA2	2.55	0.41
1:A:288:ASP:HB2	1:A:301:PHE:CE2	2.54	0.41
1:A:299:GLU:CD	1:A:299:GLU:H	2.28	0.41
1:A:271:HIS:CE1	1:A:299:GLU:HG3	2.55	0.41
1:A:290:ARG:HE	1:A:296:GLY:C	2.25	0.41
1:A:379:VAL:HG22	1:A:379:VAL:O	2.20	0.41
1:A:66:VAL:O	1:A:69:THR:N	2.53	0.41
1:A:255:THR:OG1	1:B:252:THR:HG21	2.20	0.41
1:B:477:CYS:HA	1:B:478:PRO:HD2	1.94	0.41
1:B:315:TYR:HE2	1:B:318:THR:O	2.02	0.41
1:A:326:TYR:CE2	1:A:559:LEU:HD11	2.56	0.41
1:B:213:ASN:H	1:B:213:ASN:ND2	2.05	0.41
1:A:211:VAL:HG22	1:A:290:ARG:HD2	2.01	0.41
1:B:332:PHE:CD2	1:B:332:PHE:C	2.99	0.41
1:A:270:GLY:O	1:A:271:HIS:C	2.64	0.41
1:A:271:HIS:HD2	1:A:297:ARG:HG2	1.84	0.41
1:B:158:GLY:CA	1:B:159:VAL:C	2.94	0.41
1:B:299:GLU:H	1:B:299:GLU:CD	2.29	0.41
1:A:66:VAL:O	1:A:67:ARG:C	2.64	0.41
1:A:223:LEU:CD2	1:A:467:THR:HG22	2.40	0.41
1:B:195:LEU:O	1:B:196:ALA:C	2.64	0.40
1:B:82:GLN:NE2	3:B:2005:HOH:O	2.54	0.40
1:B:30:ALA:C	1:B:32:ALA:N	2.79	0.40
1:B:224:THR:HB	1:B:228:PHE:CE2	2.56	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	512/568 (90%)	472 (92%)	32 (6%)	8 (2%)	7	9
1	B	540/568 (95%)	490 (91%)	33 (6%)	17 (3%)	3	2
All	All	1052/1136 (93%)	962 (91%)	65 (6%)	25 (2%)	4	4

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	71	GLN
1	A	159	VAL
1	A	296	GLY
1	B	49	GLY
1	B	65	ARG
1	B	297	ARG
1	A	60	VAL
1	A	314	ALA
1	B	50	LEU
1	B	71	GLN
1	A	100	ALA
1	B	32	ALA
1	B	41	ALA
1	B	314	ALA
1	A	295	GLY
1	B	29	LEU
1	B	100	ALA
1	B	107	ALA
1	B	159	VAL
1	B	294	GLY
1	B	295	GLY
1	A	294	GLY
1	B	47	GLN
1	B	37	GLY
1	B	60	VAL

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	390/429 (91%)	357 (92%)	33 (8%)	10	13
1	B	407/429 (95%)	377 (93%)	30 (7%)	13	17
All	All	797/858 (93%)	734 (92%)	63 (8%)	11	16

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	THR
1	A	71	GLN
1	A	78	VAL
1	A	89	GLU
1	A	119	GLN
1	A	123	GLU
1	A	137	ARG
1	A	192	VAL
1	A	193	ARG
1	A	198	THR
1	A	201	ARG
1	A	211	VAL
1	A	213	ASN
1	A	252	THR
1	A	259	GLN
1	A	289	LEU
1	A	290	ARG
1	A	297	ARG
1	A	299	GLU
1	A	304	ARG
1	A	311	ASN
1	A	334	LEU
1	A	370	GLN
1	A	379	VAL
1	A	380	LEU
1	A	407	SER
1	A	408	VAL

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Mol	Chain	Res	Type
1	A	499	LEU
1	A	532	ARG
1	A	540	SER
1	A	546	ASP
1	A	562	ARG
1	A	566	LEU
1	B	29	LEU
1	B	36	MET
1	B	40	THR
1	B	47	GLN
1	B	52	LEU
1	B	69	THR
1	B	71	GLN
1	B	89	GLU
1	B	119	GLN
1	B	123	GLU
1	B	159	VAL
1	B	192	VAL
1	B	193	ARG
1	B	213	ASN
1	B	233	ARG
1	B	252	THR
1	B	259	GLN
1	B	289	LEU
1	B	299	GLU
1	B	304	ARG
1	B	311	ASN
1	B	334	LEU
1	B	370	GLN
1	B	379	VAL
1	B	380	LEU
1	B	407	SER
1	B	408	VAL
1	B	499	LEU
1	B	540	SER
1	B	546	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	GLN
1	A	73	HIS

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Mol	Chain	Res	Type
1	A	99	HIS
1	A	118	HIS
1	A	119	GLN
1	A	130	GLN
1	A	177	GLN
1	A	186	GLN
1	A	213	ASN
1	A	239	GLN
1	A	259	GLN
1	A	271	HIS
1	A	281	HIS
1	A	354	GLN
1	A	370	GLN
1	A	376	GLN
1	A	385	ASN
1	A	392	GLN
1	B	47	GLN
1	B	56	HIS
1	B	82	GLN
1	B	118	HIS
1	B	133	ASN
1	B	177	GLN
1	B	186	GLN
1	B	213	ASN
1	B	235	GLN
1	B	259	GLN
1	B	271	HIS
1	B	370	GLN
1	B	392	GLN
1	B	434	GLN
1	B	446	GLN
1	B	461	HIS
1	B	557	GLN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
2	LXN	B	1569	-	10,10,10	4.93	4 (40%)	13,14,14	2.65	5 (38%)
2	LXN	A	1569	-	10,10,10	4.89	3 (30%)	13,14,14	2.37	5 (38%)

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	LXN	B	1569	-	-	-	0/1/1/1
2	LXN	A	1569	-	-	-	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1569	LXN	CAI-CAJ	-11.34	1.39	1.53
2	B	1569	LXN	CAI-CAJ	-10.83	1.39	1.53
2	B	1569	LXN	CAH-CAJ	-10.44	1.40	1.53
2	A	1569	LXN	CAH-CAJ	-10.18	1.40	1.53
2	B	1569	LXN	CAG-NAE	2.56	1.41	1.37
2	B	1569	LXN	CAG-NAF	2.04	1.40	1.37
2	A	1569	LXN	CAG-NAF	2.02	1.40	1.37

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	B	1569	LXN	CAH-NAE-CAG	-5.58	118.69	126.37
2	A	1569	LXN	CAI-NAF-CAG	-4.78	119.78	126.37
2	B	1569	LXN	CAI-NAF-CAG	-4.75	119.83	126.37
2	A	1569	LXN	CAH-NAE-CAG	-4.52	120.14	126.37
2	B	1569	LXN	CAJ-CAH-NAE	3.44	120.64	116.39
2	B	1569	LXN	NAE-CAG-NAF	3.05	120.54	115.74
2	B	1569	LXN	CAJ-CAI-NAF	2.92	120.00	116.39
2	A	1569	LXN	NAE-CAG-NAF	2.88	120.27	115.74
2	A	1569	LXN	CAJ-CAH-NAE	2.68	119.70	116.39
2	A	1569	LXN	CAJ-CAI-NAF	2.60	119.60	116.39

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [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	512/568 (90%)	0.51	32 (6%) 26 25	19, 38, 42, 74	2 (0%)
1	B	542/568 (95%)	0.64	40 (7%) 20 19	35, 38, 52, 74	0
All	All	1054/1136 (92%)	0.58	72 (6%) 23 22	19, 38, 48, 74	2 (0%)

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	39	THR	4.3
1	B	74	ALA	4.2
1	B	377	GLY	4.0
1	A	315	TYR	3.9
1	B	97	LEU	3.9
1	A	313	LEU	3.6
1	B	365	PRO	3.5
1	A	200	VAL	3.5
1	A	192	VAL	3.2
1	A	66	VAL	3.2
1	B	292	TRP	3.2
1	B	100	ALA	3.2
1	B	566	LEU	3.1
1	B	44	MET	3.1
1	B	41	ALA	3.0
1	B	312	TRP	2.9
1	B	318	THR	2.9
1	A	314	ALA	2.9
1	B	204	GLY	2.9
1	B	310	VAL	2.9
1	B	293	GLY	2.9
1	B	55	GLY	2.8
1	A	311	ASN	2.8
1	A	59	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	370	GLN	2.8
1	B	317	GLY	2.7
1	A	65	ARG	2.7
1	A	316	PRO	2.7
1	B	314	ALA	2.6
1	B	56	HIS	2.6
1	A	295	GLY	2.6
1	A	216	GLY	2.6
1	B	568	ILE	2.5
1	A	63	LEU	2.5
1	A	310	VAL	2.5
1	A	64	GLY	2.5
1	B	375	GLU	2.5
1	A	301	PHE	2.5
1	A	303	LEU	2.5
1	A	134	TRP	2.4
1	A	312	TRP	2.4
1	B	316	PRO	2.4
1	B	86	ALA	2.4
1	A	123	GLU	2.4
1	A	302	ALA	2.4
1	B	52	LEU	2.4
1	A	98	GLY	2.3
1	A	317	GLY	2.3
1	B	313	LEU	2.3
1	B	156	ALA	2.3
1	B	567	GLY	2.3
1	A	215	PHE	2.3
1	A	300	VAL	2.3
1	B	79	LEU	2.3
1	B	68	TRP	2.2
1	B	60	VAL	2.2
1	B	309	GLN	2.2
1	B	290	ARG	2.2
1	A	506	VAL	2.2
1	B	112	ALA	2.2
1	A	269	LEU	2.2
1	A	100	ALA	2.2
1	B	319	SER	2.2
1	B	540	SER	2.2
1	A	568	ILE	2.1
1	A	294	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	318	THR	2.1
1	B	376	GLN	2.1
1	B	291	GLY	2.1
1	B	205	PRO	2.1
1	B	307	PRO	2.1
1	B	98	GLY	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 oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
2	LXN	B	1569	10/10	0.78	0.10	52,54,55,55	0
2	LXN	A	1569	10/10	0.87	0.15	23,26,27,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.