



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 05:59 am GMT

PDB ID : 4FGU
Title : Crystal structure of prolegumain
Authors : Dall, E.; Brandstetter, H.
Deposited on : 2012-06-05
Resolution : 3.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

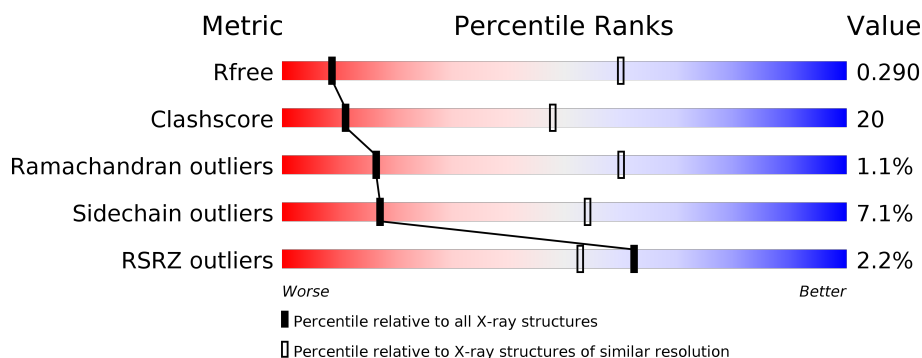
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.90 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	1007 (4.20-3.60)
Clashscore	112137	1103 (4.20-3.60)
Ramachandran outliers	110173	1062 (4.20-3.60)
Sidechain outliers	110143	1053 (4.20-3.60)
RSRZ outliers	101464	1020 (4.20-3.60)

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	429	<div> <div>2%</div> <div> <div></div> <div>64%</div> <div>27%</div> <div>• 5%</div> </div> </div>
1	B	429	<div> <div>3%</div> <div> <div></div> <div>66%</div> <div>25%</div> <div>• 6%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	A	501	-	-	-	X
3	NAG	A	503	-	-	X	X
3	NAG	B	503	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6688 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 Legumain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	407	Total	C	N	O	S	0	0	0
			3287	2082	569	614	22			
1	B	404	Total	C	N	O	S	0	0	0
			3261	2065	564	610	22			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	13	GLY	-	EXPRESSION TAG	UNP Q99538
A	14	ALA	-	EXPRESSION TAG	UNP Q99538
A	15	SER	-	EXPRESSION TAG	UNP Q99538
A	16	LEU	-	EXPRESSION TAG	UNP Q99538
A	17	GLU	-	EXPRESSION TAG	UNP Q99538
A	18	ILE	VAL	VARIANT	UNP Q99538
A	272	GLN	ASN	ENGINEERED MUTATION	UNP Q99538
A	434	GLY	-	EXPRESSION TAG	UNP Q99538
A	435	THR	-	EXPRESSION TAG	UNP Q99538
A	436	HIS	-	EXPRESSION TAG	UNP Q99538
A	437	HIS	-	EXPRESSION TAG	UNP Q99538
A	438	HIS	-	EXPRESSION TAG	UNP Q99538
A	439	HIS	-	EXPRESSION TAG	UNP Q99538
A	440	HIS	-	EXPRESSION TAG	UNP Q99538
A	441	HIS	-	EXPRESSION TAG	UNP Q99538
B	13	GLY	-	EXPRESSION TAG	UNP Q99538
B	14	ALA	-	EXPRESSION TAG	UNP Q99538
B	15	SER	-	EXPRESSION TAG	UNP Q99538
B	16	LEU	-	EXPRESSION TAG	UNP Q99538
B	17	GLU	-	EXPRESSION TAG	UNP Q99538
B	18	ILE	VAL	VARIANT	UNP Q99538
B	272	GLN	ASN	ENGINEERED MUTATION	UNP Q99538
B	434	GLY	-	EXPRESSION TAG	UNP Q99538
B	435	THR	-	EXPRESSION TAG	UNP Q99538
B	436	HIS	-	EXPRESSION TAG	UNP Q99538

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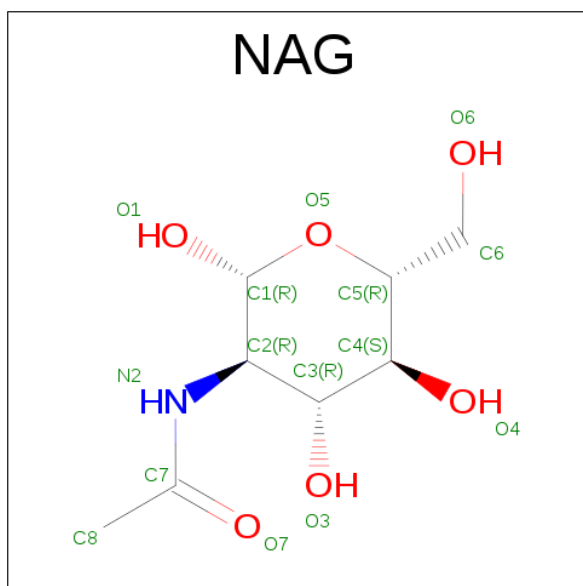
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Chain	Residue	Modelled	Actual	Comment	Reference
B	437	HIS	-	EXPRESSION TAG	UNP Q99538
B	438	HIS	-	EXPRESSION TAG	UNP Q99538
B	439	HIS	-	EXPRESSION TAG	UNP Q99538
B	440	HIS	-	EXPRESSION TAG	UNP Q99538
B	441	HIS	-	EXPRESSION TAG	UNP Q99538

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

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

- 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	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

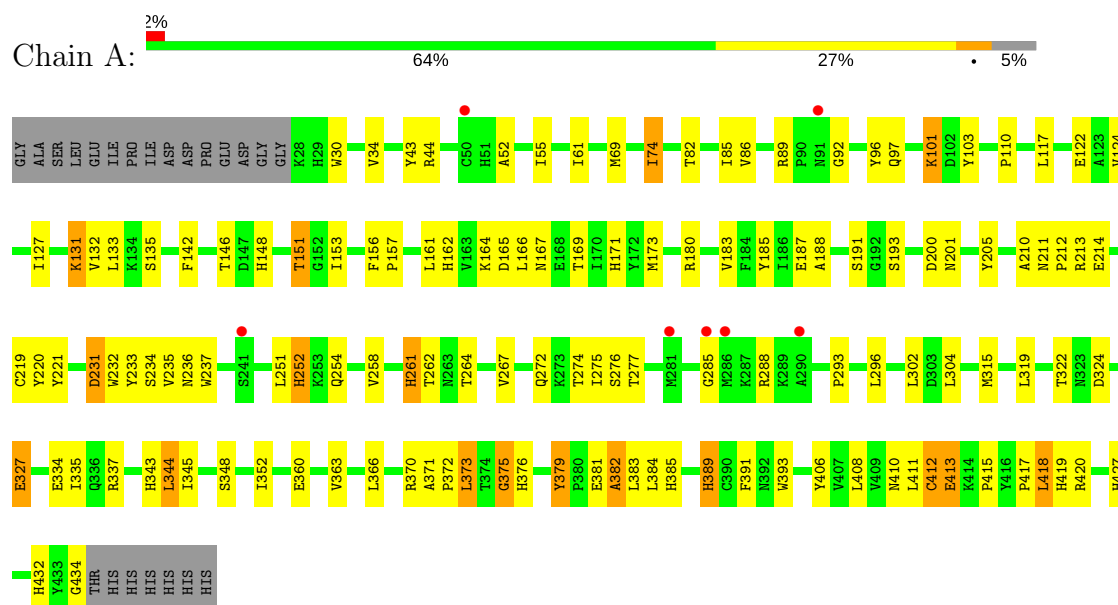
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	32	Total 32	O 32	0	0
4	B	24	Total 24	O 24	0	0

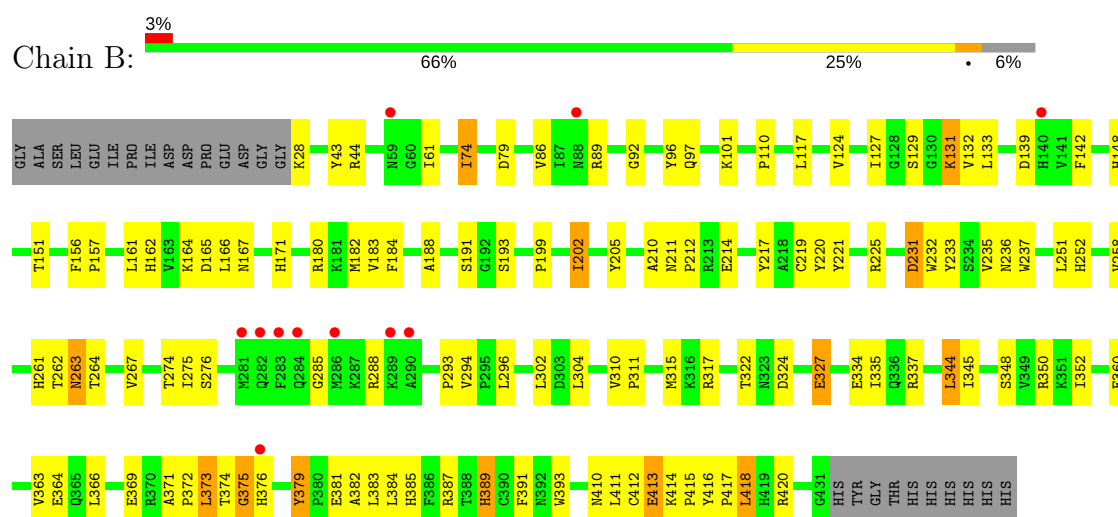
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Legumain



• Molecule 1: Legumain



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	185.10Å 185.10Å 173.07Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.63 – 3.90 49.63 – 3.90	Depositor EDS
% Data completeness (in resolution range)	99.2 (49.63-3.90) 99.3 (49.63-3.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 3.88Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.286 , 0.300 0.287 , 0.290	Depositor DCC
R_{free} test set	503 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	39.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 55.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.77	EDS
Total number of atoms	6688	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.08% 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 [i](#)

5.1 Standard geometry [i](#)

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	A	0.46	0/3377	0.56	1/4589 (0.0%)
1	B	0.47	0/3349	0.55	1/4551 (0.0%)
All	All	0.46	0/6726	0.56	2/9140 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	375	GLY	N-CA-C	5.60	127.10	113.10
1	B	375	GLY	N-CA-C	5.55	126.97	113.10

There are no chirality outliers.

There are no planarity outliers.

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	3287	0	3179	140	0
1	B	3261	0	3159	119	0
2	A	28	0	25	2	0
2	B	28	0	25	1	0
3	A	14	0	13	8	0
3	B	14	0	13	5	0
4	A	32	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	24	0	0	0	0
All	All	6688	0	6414	259	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 20.

All (259) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:171:HIS:HB2	3:A:503:NAG:C8	1.28	1.60
1:B:171:HIS:HB2	3:B:503:NAG:C8	1.55	1.36
1:B:371:ALA:HB1	1:B:372:PRO:CD	1.60	1.30
1:A:371:ALA:HB1	1:A:372:PRO:CD	1.63	1.29
1:A:201:ASN:HA	1:A:272:GLN:NE2	1.52	1.24
1:A:371:ALA:HB1	1:A:372:PRO:HD2	1.26	1.17
1:A:171:HIS:CB	3:A:503:NAG:H82	1.75	1.16
1:A:171:HIS:CB	3:A:503:NAG:C8	2.23	1.16
1:A:151:THR:HG22	1:A:373:LEU:HD23	1.22	1.15
1:A:201:ASN:CA	1:A:272:GLN:HE21	1.58	1.15
1:A:151:THR:CG2	1:A:373:LEU:HD23	1.81	1.11
1:B:151:THR:CG2	1:B:373:LEU:HD23	1.79	1.10
1:B:151:THR:HG22	1:B:373:LEU:HD23	1.30	1.10
1:B:371:ALA:HB1	1:B:372:PRO:HD2	1.30	1.09
1:A:171:HIS:HB2	3:A:503:NAG:H82	1.07	1.04
1:A:151:THR:CG2	1:A:373:LEU:CD2	2.36	1.03
1:A:151:THR:HG22	1:A:373:LEU:CD2	1.89	1.02
1:B:171:HIS:CB	3:B:503:NAG:H82	1.88	1.02
1:B:151:THR:CG2	1:B:373:LEU:CD2	2.38	1.02
1:B:171:HIS:HB2	3:B:503:NAG:H82	1.01	1.00
1:B:371:ALA:HB1	1:B:372:PRO:HD3	1.36	1.00
1:B:220:TYR:HE2	1:B:235:VAL:HG11	1.27	1.00
1:A:371:ALA:HB1	1:A:372:PRO:HD3	1.45	0.99
1:B:371:ALA:CB	1:B:372:PRO:CD	2.37	0.98
1:A:372:PRO:HG2	1:A:413:GLU:OE1	1.61	0.98
1:A:171:HIS:HB2	3:A:503:NAG:H83	1.40	0.97
1:A:220:TYR:HE2	1:A:235:VAL:HG11	1.26	0.97
1:A:371:ALA:CB	1:A:372:PRO:HD2	1.95	0.96
1:B:89:ARG:HH12	1:B:296:LEU:HD12	1.32	0.95
1:B:371:ALA:CB	1:B:372:PRO:HD2	1.95	0.94
1:A:201:ASN:HA	1:A:272:GLN:HE21	1.09	0.93
1:B:151:THR:HG22	1:B:373:LEU:CD2	1.99	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:171:HIS:HB2	3:A:503:NAG:H81	1.49	0.91
1:B:220:TYR:CE2	1:B:235:VAL:HG11	2.06	0.90
1:A:372:PRO:O	1:A:373:LEU:HD12	1.73	0.87
1:B:89:ARG:NH1	1:B:296:LEU:HD12	1.90	0.86
1:A:103:TYR:CE1	1:A:124:VAL:HG21	2.10	0.85
1:B:372:PRO:O	1:B:373:LEU:HD12	1.76	0.85
1:A:220:TYR:CE2	1:A:235:VAL:HG11	2.13	0.83
1:A:372:PRO:CG	1:A:413:GLU:OE1	2.26	0.82
1:B:372:PRO:HG2	1:B:413:GLU:OE1	1.80	0.81
1:A:262:THR:HG21	1:A:267:VAL:HG23	1.64	0.79
1:B:262:THR:HG21	1:B:267:VAL:HG23	1.64	0.78
1:A:89:ARG:NH1	1:A:296:LEU:HD12	1.98	0.78
1:A:86:VAL:HG22	1:A:96:TYR:HB2	1.66	0.78
1:B:171:HIS:HB2	3:B:503:NAG:H83	1.64	0.78
1:A:201:ASN:CA	1:A:272:GLN:NE2	2.28	0.77
1:A:344:LEU:C	1:A:344:LEU:HD23	2.06	0.76
1:B:372:PRO:CG	1:B:413:GLU:OE1	2.33	0.75
1:B:86:VAL:HG22	1:B:96:TYR:HB2	1.68	0.74
1:B:151:THR:HG23	1:B:373:LEU:CD2	2.16	0.74
1:A:372:PRO:HG3	1:A:413:GLU:HG2	1.69	0.73
1:B:164:LYS:NZ	1:B:376:HIS:O	2.21	0.73
1:B:372:PRO:HG2	1:B:413:GLU:CD	2.11	0.72
1:B:372:PRO:HG3	1:B:413:GLU:HG2	1.73	0.71
1:A:236:ASN:ND2	1:A:261:HIS:HB3	2.06	0.70
1:A:201:ASN:C	1:A:272:GLN:HE21	1.94	0.70
1:B:236:ASN:ND2	1:B:261:HIS:HB3	2.07	0.70
1:B:28:LYS:O	1:B:139:ASP:HB3	1.91	0.69
1:B:344:LEU:HD23	1:B:344:LEU:C	2.11	0.69
1:B:360:GLU:H	1:B:363:VAL:HG12	1.57	0.68
1:B:156:PHE:HB3	1:B:157:PRO:HD2	1.76	0.68
1:B:236:ASN:CG	1:B:261:HIS:HB3	2.14	0.68
1:A:151:THR:HG23	1:A:373:LEU:CD2	2.22	0.67
1:A:381:GLU:O	1:A:383:LEU:N	2.28	0.67
1:B:232:TRP:HE1	1:B:262:THR:HA	1.59	0.67
1:A:117:LEU:HA	1:A:133:LEU:HD12	1.77	0.67
1:A:264:THR:HG22	2:A:501:NAG:H82	1.76	0.67
1:A:193:SER:HB3	1:A:212:PRO:HB3	1.77	0.67
1:B:217:TYR:CD1	1:B:264:THR:HB	2.30	0.66
1:A:61:ILE:HD11	1:A:285:GLY:HA3	1.76	0.66
1:B:381:GLU:O	1:B:383:LEU:N	2.29	0.66
1:A:372:PRO:O	1:A:373:LEU:CD1	2.42	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:TYR:HE1	1:A:124:VAL:HG21	1.59	0.65
1:A:372:PRO:CG	1:A:413:GLU:HG2	2.26	0.65
1:A:236:ASN:CG	1:A:261:HIS:HB3	2.16	0.65
1:B:236:ASN:OD1	1:B:261:HIS:HB3	1.96	0.64
1:A:151:THR:CG2	1:A:373:LEU:HD21	2.28	0.64
1:B:334:GLU:HG3	1:B:337:ARG:HH21	1.62	0.64
1:B:264:THR:HG23	1:B:304:LEU:HD23	1.80	0.63
1:A:236:ASN:OD1	1:A:261:HIS:HB3	1.99	0.63
1:B:372:PRO:O	1:B:373:LEU:CD1	2.46	0.62
1:A:372:PRO:O	1:A:373:LEU:CG	2.47	0.62
1:A:151:THR:HG23	1:A:373:LEU:HD21	1.80	0.62
1:B:151:THR:HG23	1:B:373:LEU:HD21	1.80	0.62
1:B:236:ASN:OD1	1:B:261:HIS:CB	2.48	0.61
1:B:89:ARG:HG2	1:B:92:GLY:HA3	1.83	0.61
1:A:124:VAL:O	1:A:124:VAL:HG22	2.01	0.60
1:A:372:PRO:C	1:A:373:LEU:HG	2.20	0.60
1:B:171:HIS:CB	3:B:503:NAG:C8	2.52	0.60
1:A:201:ASN:C	1:A:272:GLN:NE2	2.54	0.60
1:A:232:TRP:HE1	1:A:262:THR:HA	1.66	0.60
1:A:89:ARG:HH12	1:A:296:LEU:HD12	1.64	0.60
1:B:251:LEU:HB2	1:B:276:SER:HA	1.84	0.60
1:B:372:PRO:C	1:B:373:LEU:HG	2.21	0.59
1:B:262:THR:O	1:B:262:THR:HG23	2.03	0.59
1:B:188:ALA:HB3	1:B:191:SER:HB3	1.84	0.58
1:B:344:LEU:HD23	1:B:345:ILE:N	2.19	0.58
1:A:142:PHE:CE1	1:A:185:TYR:HD2	2.22	0.58
1:B:232:TRP:O	1:B:233:TYR:HB3	2.03	0.58
1:A:171:HIS:HB3	3:A:503:NAG:H82	1.77	0.57
1:B:372:PRO:O	1:B:373:LEU:CG	2.52	0.57
1:B:263:ASN:N	1:B:263:ASN:OD1	2.37	0.57
1:A:142:PHE:HE1	1:A:185:TYR:HD2	1.51	0.57
1:B:117:LEU:HA	1:B:133:LEU:HD12	1.87	0.57
1:A:344:LEU:HD21	1:A:391:PHE:CZ	2.39	0.57
1:A:262:THR:O	1:A:262:THR:HG23	2.06	0.56
1:A:142:PHE:HD1	1:A:183:VAL:HG13	1.71	0.55
1:B:142:PHE:HD1	1:B:183:VAL:HG13	1.70	0.55
1:B:372:PRO:CG	1:B:413:GLU:HG2	2.36	0.55
1:B:96:TYR:OH	1:B:101:LYS:HE3	2.07	0.55
1:A:427:HIS:NE2	1:A:434:GLY:HA2	2.21	0.55
1:B:61:ILE:HD11	1:B:285:GLY:HA3	1.89	0.55
1:B:142:PHE:CD1	1:B:183:VAL:HG13	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:417:PRO:HG2	1:B:420:ARG:HE	1.72	0.54
1:B:302:LEU:HG	1:B:304:LEU:HD13	1.89	0.54
1:B:199:PRO:HG2	1:B:202:ILE:HG12	1.89	0.54
1:B:302:LEU:CD2	2:B:501:NAG:H5	2.37	0.54
1:B:182:MET:SD	1:B:184:PHE:HE1	2.30	0.54
1:A:302:LEU:HG	1:A:304:LEU:HD13	1.89	0.54
1:A:156:PHE:HB3	1:A:157:PRO:HD2	1.89	0.53
1:A:344:LEU:HD23	1:A:345:ILE:N	2.23	0.53
1:B:371:ALA:HB3	1:B:410:ASN:HB3	1.90	0.53
1:B:131:LYS:HD2	1:B:132:VAL:O	2.08	0.53
1:A:96:TYR:OH	1:A:101:LYS:CE	2.57	0.53
1:A:96:TYR:OH	1:A:101:LYS:HE3	2.07	0.53
1:A:131:LYS:HD3	1:A:132:VAL:N	2.23	0.53
1:A:334:GLU:HG3	1:A:337:ARG:HH21	1.74	0.53
1:A:236:ASN:HD21	1:A:261:HIS:HB3	1.74	0.53
1:A:344:LEU:CD2	1:A:344:LEU:C	2.78	0.52
1:B:366:LEU:HD11	1:B:414:LYS:HG3	1.90	0.52
1:A:372:PRO:HG2	1:A:413:GLU:CD	2.28	0.52
1:A:375:GLY:O	1:A:376:HIS:HB2	2.09	0.51
1:A:360:GLU:HA	1:A:363:VAL:HG12	1.92	0.51
1:B:375:GLY:O	1:B:376:HIS:HB2	2.09	0.51
1:B:315:MET:HB3	1:B:335:ILE:CG2	2.41	0.50
1:A:324:ASP:HB3	1:A:327:GLU:HB2	1.93	0.50
1:B:315:MET:HB3	1:B:335:ILE:HG22	1.92	0.50
1:A:274:THR:O	1:A:277:THR:HG23	2.12	0.50
1:B:372:PRO:CG	1:B:413:GLU:CD	2.77	0.50
1:A:131:LYS:HD2	1:A:132:VAL:O	2.10	0.50
1:B:156:PHE:HB3	1:B:157:PRO:CD	2.41	0.50
1:A:379:TYR:HA	1:A:383:LEU:HB2	1.93	0.50
1:B:379:TYR:HA	1:B:383:LEU:HB2	1.93	0.50
1:A:372:PRO:O	1:A:373:LEU:CB	2.60	0.50
1:B:124:VAL:HA	1:B:127:ILE:HD12	1.94	0.50
1:A:52:ALA:O	1:A:55:ILE:HG22	2.12	0.49
1:A:153:ILE:HD13	1:A:384:LEU:HD21	1.95	0.49
1:A:219:CYS:SG	1:A:231:ASP:C	2.91	0.49
1:A:344:LEU:CD2	1:A:391:PHE:CZ	2.96	0.49
1:B:366:LEU:HD23	1:B:411:LEU:HD12	1.95	0.49
1:B:205:TYR:HB2	1:B:275:ILE:HB	1.96	0.48
1:A:348:SER:O	1:A:352:ILE:HG12	2.14	0.48
1:A:372:PRO:C	1:A:373:LEU:CG	2.80	0.48
1:B:372:PRO:C	1:B:373:LEU:CG	2.82	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:TRP:O	1:A:233:TYR:HB3	2.12	0.48
1:A:210:ALA:HB1	1:A:214:GLU:HG3	1.94	0.48
1:A:366:LEU:HD23	1:A:411:LEU:HD12	1.95	0.48
1:A:103:TYR:HE1	1:A:124:VAL:CG2	2.24	0.48
1:B:219:CYS:SG	1:B:231:ASP:C	2.92	0.48
1:A:103:TYR:CZ	1:A:124:VAL:HG21	2.47	0.48
1:A:213:ARG:HA	1:A:370:ARG:NH1	2.29	0.48
1:B:96:TYR:OH	1:B:101:LYS:CE	2.62	0.48
1:A:110:PRO:HD3	1:A:156:PHE:CD2	2.49	0.48
1:A:161:LEU:HD21	1:A:166:LEU:HD13	1.96	0.48
1:A:322:THR:HG22	1:A:324:ASP:H	1.78	0.48
1:A:200:ASP:C	1:A:272:GLN:HG2	2.34	0.47
1:B:389:HIS:ND1	1:B:389:HIS:N	2.62	0.47
1:A:103:TYR:OH	1:A:124:VAL:HG11	2.15	0.47
1:B:384:LEU:O	1:B:387:ARG:HB3	2.14	0.47
1:A:142:PHE:CD1	1:A:183:VAL:HG13	2.49	0.47
1:A:254:GLN:O	1:A:258:VAL:HG23	2.15	0.47
1:B:110:PRO:HD3	1:B:156:PHE:CD2	2.50	0.47
1:A:162:HIS:CE1	1:A:384:LEU:HD22	2.50	0.46
1:A:406:TYR:O	1:A:410:ASN:ND2	2.49	0.46
1:B:372:PRO:O	1:B:373:LEU:CB	2.62	0.46
1:B:193:SER:HB3	1:B:212:PRO:HB3	1.98	0.46
1:B:161:LEU:HD21	1:B:166:LEU:HD13	1.97	0.46
1:B:205:TYR:CZ	1:B:276:SER:HB3	2.51	0.46
1:B:210:ALA:HB1	1:B:214:GLU:HG3	1.97	0.45
1:B:236:ASN:HD21	1:B:261:HIS:HB3	1.78	0.45
1:A:188:ALA:HB3	1:A:191:SER:HB3	1.99	0.45
1:A:389:HIS:N	1:A:389:HIS:ND1	2.64	0.45
1:B:360:GLU:N	1:B:363:VAL:HG12	2.29	0.45
1:B:221:TYR:HE2	1:B:317:ARG:HG2	1.81	0.45
1:B:383:LEU:HD21	1:B:418:LEU:O	2.17	0.45
1:B:371:ALA:CB	1:B:410:ASN:HB3	2.47	0.45
1:B:89:ARG:CG	1:B:92:GLY:HA3	2.46	0.45
1:A:89:ARG:HG2	1:A:92:GLY:HA3	1.98	0.45
1:A:44:ARG:HH22	1:A:148:HIS:CE1	2.35	0.45
1:B:44:ARG:HH22	1:B:148:HIS:CE1	2.35	0.45
1:B:322:THR:HG21	1:B:327:GLU:HB3	1.99	0.45
1:B:372:PRO:HG3	1:B:413:GLU:CG	2.46	0.45
1:A:30:TRP:CD1	1:A:135:SER:HB3	2.52	0.44
1:A:417:PRO:HG2	1:A:420:ARG:HE	1.82	0.44
1:A:419:HIS:CE1	1:A:420:ARG:HG3	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:414:LYS:O	1:B:416:TYR:HD2	2.01	0.44
1:B:156:PHE:HE2	1:B:161:LEU:HB2	1.83	0.44
1:B:236:ASN:OD1	1:B:261:HIS:HB2	2.16	0.44
1:B:372:PRO:HG3	1:B:413:GLU:OE1	2.15	0.44
1:A:82:THR:HB	1:A:85:ILE:HB	1.99	0.44
1:B:344:LEU:CD2	1:B:344:LEU:C	2.83	0.44
1:B:360:GLU:HA	1:B:363:VAL:HG12	1.99	0.44
1:B:374:THR:HB	1:B:413:GLU:OE1	2.17	0.44
1:A:124:VAL:HA	1:A:127:ILE:HD12	2.00	0.44
1:A:89:ARG:CG	1:A:92:GLY:HA3	2.47	0.44
1:B:372:PRO:CG	1:B:413:GLU:CG	2.95	0.44
1:A:343:HIS:ND1	1:A:343:HIS:O	2.51	0.43
1:A:418:LEU:H	1:A:418:LEU:HD23	1.82	0.43
1:B:348:SER:O	1:B:352:ILE:HG12	2.17	0.43
1:A:131:LYS:HD3	1:A:132:VAL:H	1.82	0.43
1:A:142:PHE:HE1	1:A:185:TYR:CD2	2.34	0.43
1:A:372:PRO:CG	1:A:413:GLU:CG	2.95	0.43
1:A:252:HIS:ND1	1:A:276:SER:OG	2.47	0.43
1:A:344:LEU:CD2	1:A:345:ILE:N	2.81	0.43
1:A:187:GLU:HG2	1:A:234:SER:HB2	1.99	0.43
1:A:110:PRO:HD3	1:A:156:PHE:HD2	1.83	0.43
1:A:44:ARG:NH2	1:A:148:HIS:CE1	2.87	0.43
1:A:302:LEU:CD2	2:A:501:NAG:H5	2.49	0.42
1:A:344:LEU:HD21	1:A:391:PHE:CE1	2.54	0.42
1:A:171:HIS:CG	3:A:503:NAG:H83	2.54	0.42
1:B:264:THR:CG2	1:B:304:LEU:HD23	2.47	0.42
1:A:251:LEU:HB2	1:A:276:SER:HA	2.01	0.42
1:A:272:GLN:HB2	1:A:275:ILE:HD12	2.00	0.42
1:B:237:TRP:HB3	1:B:258:VAL:HG21	2.02	0.42
1:A:381:GLU:O	1:A:382:ALA:C	2.59	0.42
1:B:344:LEU:HD21	1:B:391:PHE:CZ	2.55	0.41
1:B:344:LEU:CD2	1:B:345:ILE:N	2.83	0.41
1:B:43:TYR:CD1	1:B:74:ILE:HD11	2.55	0.41
1:A:131:LYS:CD	1:A:132:VAL:N	2.83	0.41
1:B:315:MET:CB	1:B:335:ILE:HG22	2.50	0.41
1:A:371:ALA:CB	1:A:372:PRO:HD3	2.31	0.41
1:A:43:TYR:CD1	1:A:74:ILE:HD11	2.56	0.41
1:B:310:VAL:N	1:B:311:PRO:HD2	2.34	0.41
1:B:350:ARG:NH1	1:B:364:GLU:OE2	2.47	0.41
1:A:360:GLU:H	1:A:363:VAL:HG12	1.85	0.41
1:A:334:GLU:HG3	1:A:337:ARG:NH2	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:374:THR:CB	1:B:413:GLU:OE1	2.68	0.41
1:A:183:VAL:HG23	1:A:205:TYR:HD2	1.86	0.41
1:A:237:TRP:HB3	1:A:258:VAL:HG21	2.03	0.41
1:A:315:MET:HB3	1:A:335:ILE:CG2	2.51	0.41
1:A:164:LYS:NZ	1:A:376:HIS:O	2.48	0.41
1:B:324:ASP:HB3	1:B:327:GLU:HB2	2.03	0.41
1:A:236:ASN:OD1	1:A:261:HIS:CB	2.68	0.41
1:A:34:VAL:HG22	1:A:69:MET:HB2	2.03	0.41
1:B:379:TYR:CZ	1:B:384:LEU:HD12	2.56	0.41
1:A:379:TYR:CZ	1:A:384:LEU:HD12	2.56	0.40
1:B:131:LYS:HD3	1:B:132:VAL:N	2.36	0.40
1:B:162:HIS:CE1	1:B:381:GLU:H	2.38	0.40
1:B:131:LYS:CD	1:B:132:VAL:N	2.83	0.40
1:B:416:TYR:HA	1:B:417:PRO:HD3	1.94	0.40
1:A:183:VAL:HA	1:A:205:TYR:O	2.21	0.40
1:A:61:ILE:HD11	1:A:285:GLY:CA	2.47	0.40
1:A:169:THR:O	1:A:173:MET:HG3	2.22	0.40
1:B:369:GLU:O	1:B:410:ASN:ND2	2.54	0.40
1:A:193:SER:N	1:A:212:PRO:HG3	2.37	0.40
1:A:371:ALA:HB3	1:A:410:ASN:HB3	2.02	0.40
1:A:408:LEU:O	1:A:412:CYS:N	2.55	0.40
1:A:383:LEU:HD12	1:A:412:CYS:SG	2.62	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	405/429 (94%)	363 (90%)	37 (9%)	5 (1%)	15	58
1	B	402/429 (94%)	364 (90%)	34 (8%)	4 (1%)	18	61
All	All	807/858 (94%)	727 (90%)	71 (9%)	9 (1%)	17	60

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	293	PRO
1	B	293	PRO
1	A	151	THR
1	A	382	ALA
1	B	382	ALA
1	A	101	LYS
1	A	415	PRO
1	B	415	PRO
1	B	202	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	A	365/383 (95%)	339 (93%)	26 (7%)	17	54
1	B	363/383 (95%)	337 (93%)	26 (7%)	17	53
All	All	728/766 (95%)	676 (93%)	52 (7%)	17	54

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

Mol	Chain	Res	Type
1	A	74	ILE
1	A	97	GLN
1	A	122	GLU
1	A	131	LYS
1	A	146	THR
1	A	165	ASP
1	A	167	ASN
1	A	180	ARG
1	A	211	ASN
1	A	221	TYR
1	A	231	ASP
1	A	252	HIS
1	A	261	HIS
1	A	288	ARG

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Mol	Chain	Res	Type
1	A	319	LEU
1	A	327	GLU
1	A	344	LEU
1	A	373	LEU
1	A	379	TYR
1	A	385	HIS
1	A	389	HIS
1	A	393	TRP
1	A	412	CYS
1	A	413	GLU
1	A	418	LEU
1	A	432	HIS
1	B	74	ILE
1	B	79	ASP
1	B	97	GLN
1	B	129	SER
1	B	131	LYS
1	B	165	ASP
1	B	167	ASN
1	B	180	ARG
1	B	211	ASN
1	B	225	ARG
1	B	231	ASP
1	B	252	HIS
1	B	263	ASN
1	B	274	THR
1	B	288	ARG
1	B	294	VAL
1	B	327	GLU
1	B	344	LEU
1	B	373	LEU
1	B	379	TYR
1	B	385	HIS
1	B	389	HIS
1	B	393	TRP
1	B	412	CYS
1	B	413	GLU
1	B	418	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	HIS
1	A	148	HIS
1	A	162	HIS
1	A	272	GLN
1	A	404	HIS
1	A	410	ASN
1	B	111	GLN
1	B	140	HIS
1	B	148	HIS
1	B	162	HIS
1	B	385	HIS
1	B	410	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

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	$\# Z > 2$
2	NAG	A	501	1,2	14,14,15	0.54	0	15,19,21	1.43	2 (13%)
2	NAG	A	502	2	14,14,15	0.43	0	15,19,21	1.02	1 (6%)
2	NAG	B	501	1,2	14,14,15	0.64	0	15,19,21	1.79	3 (20%)
2	NAG	B	502	2	14,14,15	0.41	0	15,19,21	1.01	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
2	NAG	A	501	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	502	2	-	0/6/23/26	0/1/1/1
2	NAG	B	501	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	502	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	NAG	O5-C1-C2	-3.00	107.30	111.47
2	A	502	NAG	O5-C1-C2	-2.14	108.50	111.47
2	A	501	NAG	C4-C3-C2	2.74	115.03	111.02
2	B	501	NAG	C4-C3-C2	2.87	115.22	111.02
2	A	501	NAG	C1-O5-C5	4.19	117.94	112.17
2	B	501	NAG	C1-O5-C5	5.26	119.41	112.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	NAG	2	0
2	B	501	NAG	1	0

5.6 Ligand geometry

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
3	NAG	A	503	1	14,14,15	0.43	0	15,19,21	0.95	1 (6%)
3	NAG	B	503	1	14,14,15	0.43	0	15,19,21	0.88	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
3	NAG	A	503	1	-	0/6/23/26	0/1/1/1
3	NAG	B	503	1	-	0/6/23/26	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(°)
3	B	503	NAG	C1-O5-C5	2.29	115.32	112.17
3	A	503	NAG	C1-O5-C5	2.74	115.94	112.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	503	NAG	8	0
3	B	503	NAG	5	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 [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	407/429 (94%)	0.20	7 (1%) 70 61	22, 53, 93, 125	0
1	B	404/429 (94%)	0.23	11 (2%) 55 45	26, 55, 96, 141	0
All	All	811/858 (94%)	0.21	18 (2%) 62 53	22, 54, 95, 141	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	290	ALA	4.2
1	A	290	ALA	3.4
1	A	286	MET	3.3
1	A	285	GLY	2.8
1	A	281	MET	2.7
1	B	286	MET	2.6
1	B	376	HIS	2.5
1	B	289	LYS	2.5
1	A	50	CYS	2.4
1	B	283	PHE	2.4
1	B	282	GLN	2.3
1	A	241	SER	2.3
1	A	91	ASN	2.2
1	B	88	ASN	2.2
1	B	140	HIS	2.1
1	B	59	ASN	2.1
1	B	281	MET	2.1
1	B	284	GLN	2.0

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

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

6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	NAG	A	501	14/15	0.79	0.36	2.72	17,21,22,22	0
2	NAG	A	502	14/15	0.61	0.54	-	5,6,6,7	0
2	NAG	B	502	14/15	0.79	0.44	-	5,6,6,6	0
2	NAG	B	501	14/15	0.84	0.32	-	16,20,21,21	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(Å ²)	Q<0.9
3	NAG	A	503	14/15	0.50	0.50	4.45	29,29,30,30	0
3	NAG	B	503	14/15	0.47	0.55	1.37	30,30,31,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.