



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 10, 2019 – 08:47 AM EST

PDB ID : 6N4X
Title : Metabotropic Glutamate Receptor 5 Apo Form Ligand Binding Domain
Authors : Koehl, A.; Hu, H.; Feng, D.; Sun, B.; Weis, W.I.; Skiniotis, G.S.; Mathiesen, J.M.; Kobilka, B.K.
Deposited on : 2018-11-20
Resolution : 4.00 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : rb-20031633
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20031633

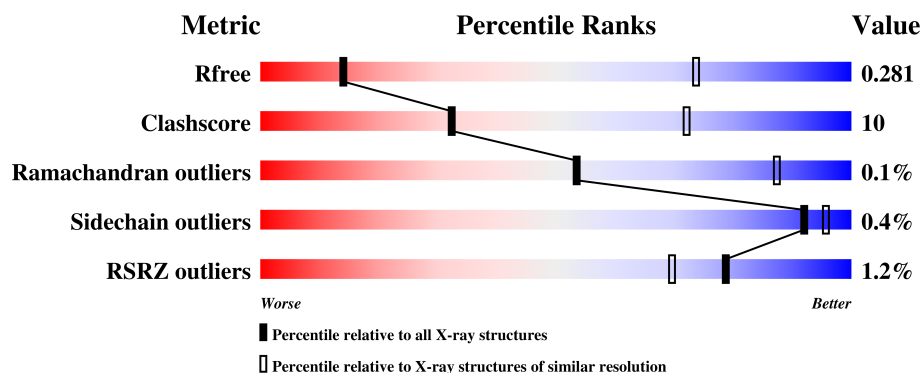
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.00 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}	111664	1008 (4.38-3.62)
Clashscore	122126	1012 (4.34-3.66)
Ramachandran outliers	120053	1000 (4.36-3.64)
Sidechain outliers	120020	1023 (4.38-3.62)
RSRZ outliers	108989	1107 (4.40-3.60)

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	877	<div> <div></div> <div>45%13%42%</div> </div>
1	B	877	<div> <div></div> <div>43%13%44%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7736 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 Metabotropic glutamate receptor 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	506	Total	C	N	O	S	0	0	0
			3897	2456	666	739	36			
1	B	487	Total	C	N	O	S	0	0	0
			3754	2362	646	712	34			

There are 62 discrepancies between the modelled and reference sequences:

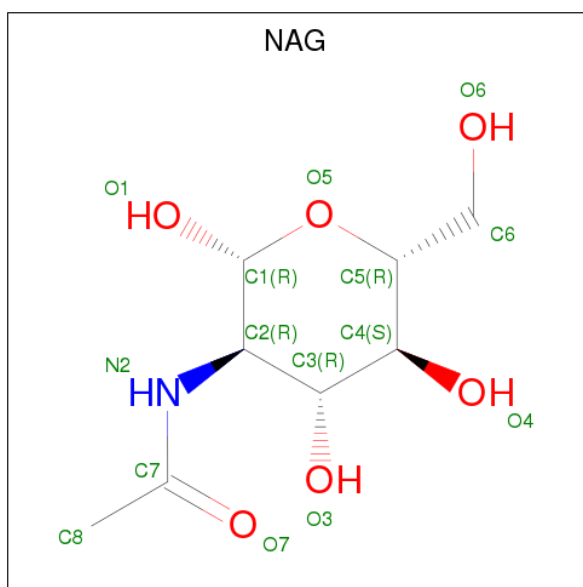
Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	MET	-	initiating methionine	UNP P41594
A	-4	LYS	-	expression tag	UNP P41594
A	-3	THR	-	expression tag	UNP P41594
A	-2	ILE	-	expression tag	UNP P41594
A	-1	ILE	-	expression tag	UNP P41594
A	0	ALA	-	expression tag	UNP P41594
A	1	LEU	-	expression tag	UNP P41594
A	2	SER	-	expression tag	UNP P41594
A	3	TYR	-	expression tag	UNP P41594
A	4	ILE	-	expression tag	UNP P41594
A	5	PHE	-	expression tag	UNP P41594
A	6	CYS	-	expression tag	UNP P41594
A	7	LEU	-	expression tag	UNP P41594
A	8	VAL	-	expression tag	UNP P41594
A	9	PHE	-	expression tag	UNP P41594
A	10	ALA	-	expression tag	UNP P41594
A	11	ASP	-	expression tag	UNP P41594
A	12	TYR	-	expression tag	UNP P41594
A	13	LYS	-	expression tag	UNP P41594
A	14	ASP	-	expression tag	UNP P41594
A	15	ASP	-	expression tag	UNP P41594
A	16	ASP	-	expression tag	UNP P41594
A	17	ASP	-	expression tag	UNP P41594
A	18	ALA	-	expression tag	UNP P41594
A	19	ALA	-	expression tag	UNP P41594

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Chain	Residue	Modelled	Actual	Comment	Reference
A	866	HIS	-	expression tag	UNP P41594
A	867	HIS	-	expression tag	UNP P41594
A	868	HIS	-	expression tag	UNP P41594
A	869	HIS	-	expression tag	UNP P41594
A	870	HIS	-	expression tag	UNP P41594
A	871	HIS	-	expression tag	UNP P41594
B	-5	MET	-	initiating methionine	UNP P41594
B	-4	LYS	-	expression tag	UNP P41594
B	-3	THR	-	expression tag	UNP P41594
B	-2	ILE	-	expression tag	UNP P41594
B	-1	ILE	-	expression tag	UNP P41594
B	0	ALA	-	expression tag	UNP P41594
B	1	LEU	-	expression tag	UNP P41594
B	2	SER	-	expression tag	UNP P41594
B	3	TYR	-	expression tag	UNP P41594
B	4	ILE	-	expression tag	UNP P41594
B	5	PHE	-	expression tag	UNP P41594
B	6	CYS	-	expression tag	UNP P41594
B	7	LEU	-	expression tag	UNP P41594
B	8	VAL	-	expression tag	UNP P41594
B	9	PHE	-	expression tag	UNP P41594
B	10	ALA	-	expression tag	UNP P41594
B	11	ASP	-	expression tag	UNP P41594
B	12	TYR	-	expression tag	UNP P41594
B	13	LYS	-	expression tag	UNP P41594
B	14	ASP	-	expression tag	UNP P41594
B	15	ASP	-	expression tag	UNP P41594
B	16	ASP	-	expression tag	UNP P41594
B	17	ASP	-	expression tag	UNP P41594
B	18	ALA	-	expression tag	UNP P41594
B	19	ALA	-	expression tag	UNP P41594
B	866	HIS	-	expression tag	UNP P41594
B	867	HIS	-	expression tag	UNP P41594
B	868	HIS	-	expression tag	UNP P41594
B	869	HIS	-	expression tag	UNP P41594
B	870	HIS	-	expression tag	UNP P41594
B	871	HIS	-	expression tag	UNP P41594

- Molecule 2 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

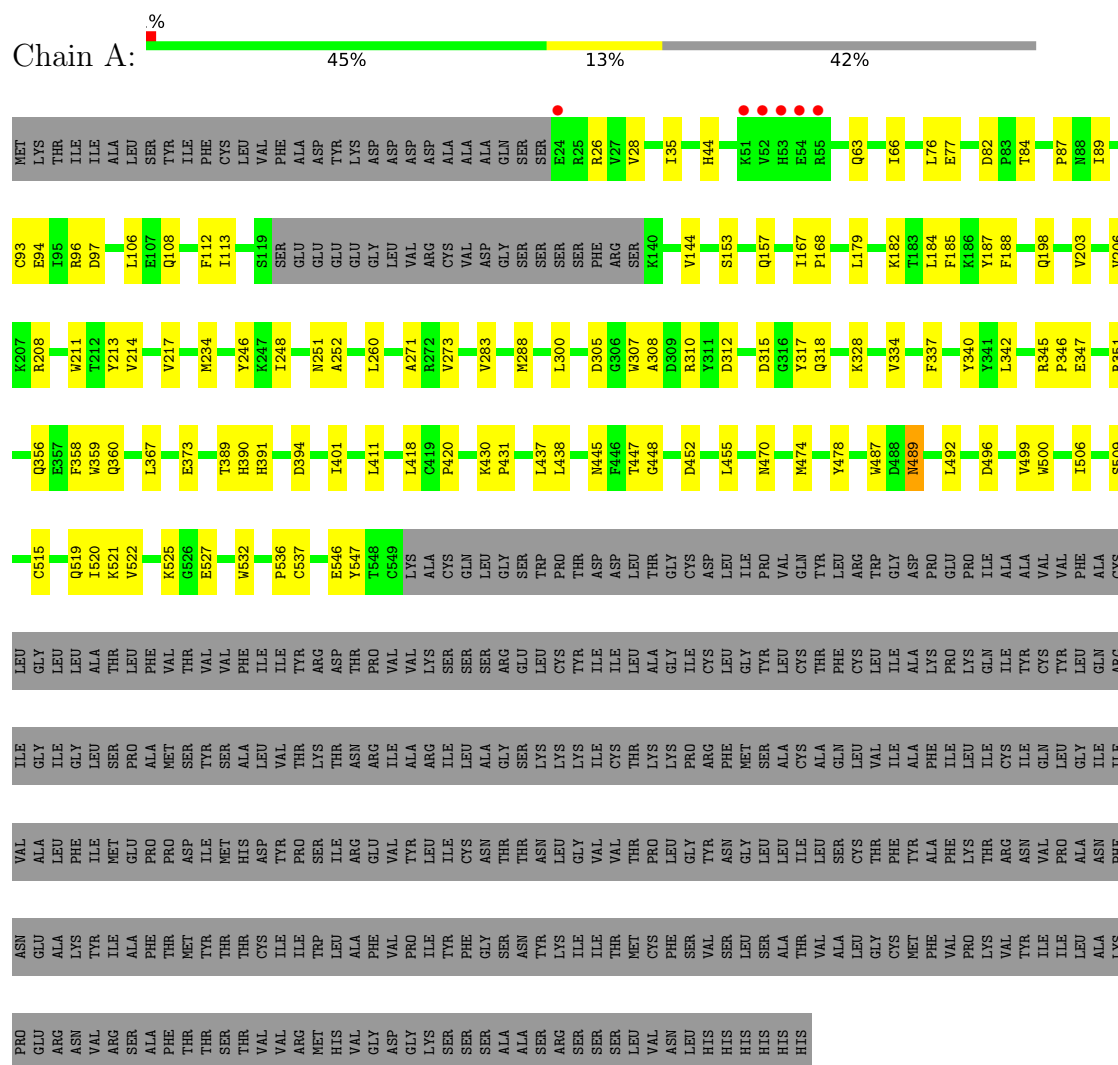
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		

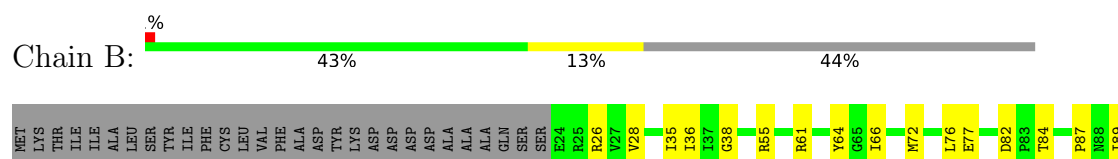
3 Residue-property plots

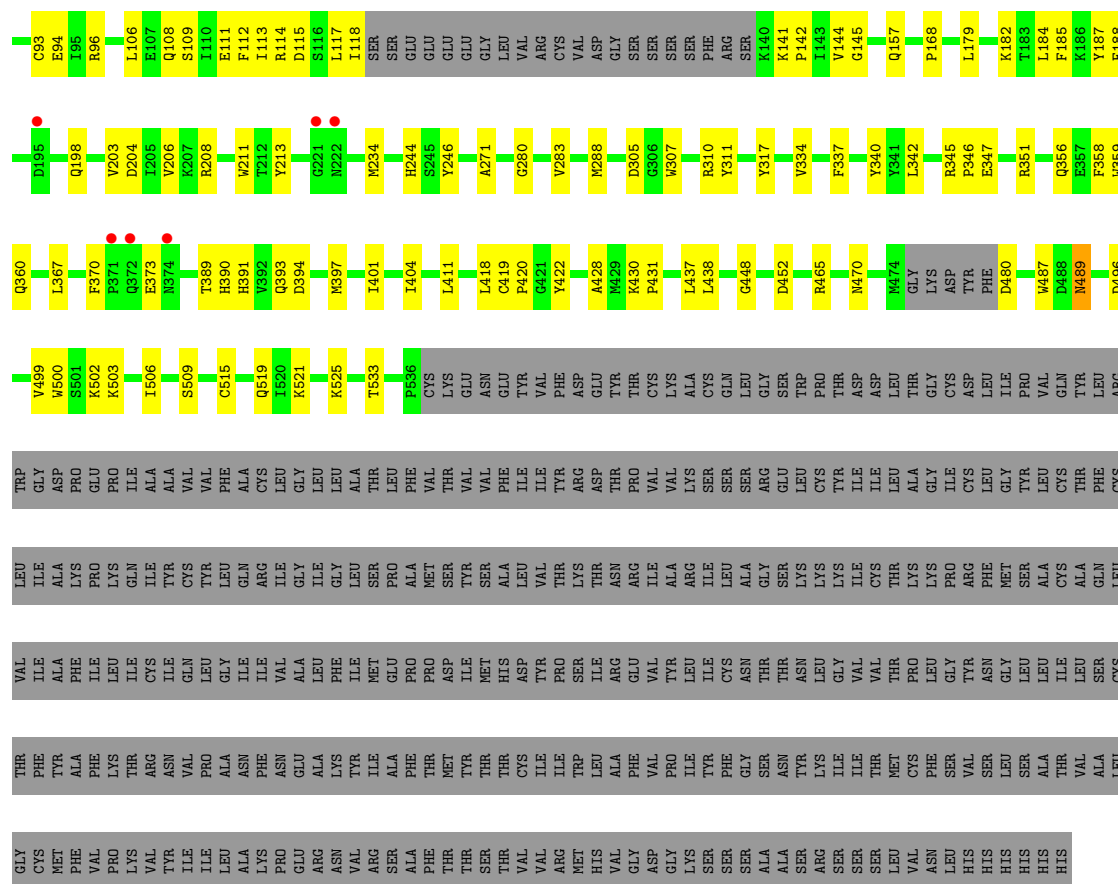
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Metabotropic glutamate receptor 5



• Molecule 1: Metabotropic glutamate receptor 5





4 Data and refinement statistics

Property	Value	Source
Space group	P 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	119.08Å 174.10Å 180.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.57 – 4.00 49.57 – 4.00	Depositor EDS
% Data completeness (in resolution range)	93.2 (49.57-4.00) 83.5 (49.57-4.00)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 4.00Å)	Xtriage
Refinement program	PHENIX 1.14_3211	Depositor
R, R_{free}	0.270 , 0.282 0.268 , 0.281	Depositor DCC
R_{free} test set	1540 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	172.6	Xtriage
Anisotropy	0.246	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 160.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.064 for -h,l,k	Xtriage
F_o, F_c correlation	0.82	EDS
Total number of atoms	7736	wwPDB-VP
Average B, all atoms (Å ²)	184.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.29% 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: MG, 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	0.25	0/3985	0.45	0/5407
1	B	0.24	0/3836	0.45	0/5200
All	All	0.25	0/7821	0.45	0/10607

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	3897	0	3649	73	0
1	B	3754	0	3546	71	0
2	A	56	0	52	0	0
2	B	28	0	26	0	0
3	B	1	0	0	0	0
All	All	7736	0	7273	143	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 (143) 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:520:ILE:HD13	1:A:546:GLU:C	1.90	0.90
1:A:77:GLU:OE2	1:A:351:ARG:NH1	2.18	0.76
1:B:77:GLU:OE2	1:B:351:ARG:NH1	2.19	0.74
1:B:360:GLN:HE22	1:B:373:GLU:HA	1.52	0.74
1:A:360:GLN:HE22	1:A:373:GLU:HA	1.53	0.73
1:A:522:VAL:HG22	1:A:547:TYR:HA	1.72	0.70
1:B:419:CYS:HB3	1:B:422:TYR:HB2	1.74	0.69
1:B:198:GLN:HE22	1:B:305:ASP:H	1.40	0.69
1:A:182:LYS:NZ	1:A:188:PHE:O	2.27	0.67
1:A:198:GLN:HE22	1:A:305:ASP:H	1.42	0.66
1:B:182:LYS:NZ	1:B:188:PHE:O	2.29	0.65
1:A:82:ASP:OD2	1:A:84:THR:OG1	2.11	0.64
1:A:94:GLU:OE2	1:A:96:ARG:NH2	2.32	0.63
1:A:447:THR:HG21	1:B:428:ALA:HB2	1.80	0.63
1:B:213:TYR:O	1:B:509:SER:OG	2.16	0.63
1:B:106:LEU:HD21	1:B:157:GLN:HB3	1.81	0.61
1:A:198:GLN:NE2	1:A:305:ASP:H	1.97	0.61
1:B:525:LYS:HG2	1:B:533:THR:HG23	1.82	0.61
1:A:345:ARG:HD3	1:A:346:PRO:HD2	1.83	0.60
1:A:106:LEU:HD21	1:A:157:GLN:HB3	1.83	0.60
1:B:311:TYR:OH	1:B:480:ASP:OD1	2.16	0.60
1:B:109:SER:HA	1:B:112:PHE:HD2	1.65	0.60
1:A:520:ILE:HD13	1:A:546:GLU:O	2.02	0.60
1:B:345:ARG:HD3	1:B:346:PRO:HD2	1.84	0.59
1:A:35:ILE:HG23	1:A:144:VAL:HG21	1.83	0.59
1:A:310:ARG:HG2	1:A:312:ASP:OD1	2.03	0.59
1:B:208:ARG:NH1	1:B:496:ASP:OD1	2.31	0.58
1:B:342:LEU:HD21	1:B:391:HIS:CG	2.39	0.57
1:B:35:ILE:HG23	1:B:144:VAL:HG21	1.86	0.57
1:A:487:TRP:HD1	1:A:492:LEU:HD13	1.71	0.56
1:A:213:TYR:O	1:A:509:SER:OG	2.22	0.56
1:A:342:LEU:HD21	1:A:391:HIS:CG	2.41	0.55
1:A:525:LYS:HG3	1:A:527:GLU:H	1.71	0.55
1:A:26:ARG:HD2	1:A:108:GLN:NE2	2.21	0.55
1:A:168:PRO:HG2	1:A:437:LEU:HD23	1.88	0.55
1:A:203:VAL:HG21	1:A:234:MET:HB2	1.89	0.54
1:A:208:ARG:NH1	1:A:496:ASP:OD1	2.27	0.54
1:A:340:TYR:OH	1:A:351:ARG:NH2	2.33	0.54
1:A:28:VAL:HG11	1:A:94:GLU:HG3	1.88	0.54
1:A:519:GLN:HB3	1:A:536:PRO:HA	1.89	0.54
1:B:198:GLN:NE2	1:B:305:ASP:H	2.04	0.54
1:B:206:VAL:HG13	1:B:211:TRP:HB2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:94:GLU:OE2	1:B:96:ARG:NH2	2.41	0.53
1:A:520:ILE:HG13	1:A:521:LYS:H	1.74	0.53
1:B:203:VAL:HG21	1:B:234:MET:HB2	1.90	0.53
1:B:82:ASP:OD2	1:B:84:THR:OG1	2.13	0.53
1:A:345:ARG:HG3	1:A:347:GLU:H	1.74	0.53
1:B:109:SER:O	1:B:113:ILE:HG12	2.09	0.53
1:A:182:LYS:HZ3	1:A:188:PHE:HB3	1.73	0.52
1:B:288:MET:HG2	1:B:317:TYR:CZ	2.45	0.52
1:B:345:ARG:HG3	1:B:347:GLU:H	1.74	0.52
1:A:44:HIS:ND1	1:A:97:ASP:OD2	2.42	0.51
1:A:500:TRP:CD1	1:A:506:ILE:HG12	2.46	0.51
1:A:273:VAL:HG22	1:A:300:LEU:HD23	1.93	0.51
1:A:519:GLN:HB3	1:A:537:CYS:H	1.76	0.51
1:A:334:VAL:HG21	1:A:401:ILE:HD12	1.94	0.50
1:B:179:LEU:HD13	1:B:188:PHE:CZ	2.47	0.50
1:B:418:LEU:O	1:B:420:PRO:HD3	2.11	0.50
1:B:26:ARG:HD2	1:B:108:GLN:NE2	2.27	0.49
1:B:499:VAL:HG13	1:B:500:TRP:CD1	2.47	0.49
1:B:61:ARG:HG2	1:B:64:TYR:HB2	1.93	0.49
1:A:520:ILE:CD1	1:A:546:GLU:C	2.75	0.49
1:A:499:VAL:HG13	1:A:500:TRP:CD1	2.47	0.49
1:B:36:ILE:HD12	1:B:141:LYS:HB3	1.94	0.49
1:A:448:GLY:N	1:A:452:ASP:O	2.31	0.49
1:B:28:VAL:HG11	1:B:94:GLU:HG3	1.93	0.49
1:B:76:LEU:CD1	1:B:93:CYS:HB3	2.43	0.48
1:B:334:VAL:HG21	1:B:401:ILE:HD12	1.94	0.48
1:A:515:CYS:SG	1:A:532:TRP:NE1	2.85	0.48
1:A:248:ILE:HB	1:A:260:LEU:HD13	1.94	0.48
1:A:474:MET:O	1:A:478:TYR:HB3	2.12	0.48
1:A:418:LEU:O	1:A:420:PRO:HD3	2.13	0.48
1:A:179:LEU:HD13	1:A:188:PHE:CZ	2.49	0.47
1:A:288:MET:HG2	1:A:317:TYR:CZ	2.49	0.47
1:A:445:ASN:HB2	1:A:455:LEU:HD12	1.95	0.47
1:A:206:VAL:HG13	1:A:211:TRP:HB2	1.96	0.47
1:A:144:VAL:HG11	1:A:411:LEU:HD21	1.96	0.47
1:B:168:PRO:HG2	1:B:437:LEU:HD23	1.97	0.47
1:B:76:LEU:HD11	1:B:93:CYS:HB3	1.96	0.47
1:A:76:LEU:CD1	1:A:93:CYS:HB3	2.45	0.47
1:B:28:VAL:HG22	1:B:96:ARG:HG2	1.97	0.47
1:B:489:ASN:C	1:B:489:ASN:HD22	2.18	0.47
1:B:470:ASN:ND2	1:B:499:VAL:HG23	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:187:TYR:CD1	1:B:438:LEU:HD22	2.50	0.46
1:A:489:ASN:C	1:A:489:ASN:HD22	2.19	0.46
1:B:117:LEU:HD13	1:B:117:LEU:O	2.15	0.45
1:B:87:PRO:O	1:B:89:ILE:HD12	2.16	0.45
1:A:521:LYS:HD3	1:A:537:CYS:HA	1.99	0.45
1:A:187:TYR:CD1	1:A:438:LEU:HD22	2.51	0.45
1:B:502:LYS:HG2	1:B:503:LYS:N	2.32	0.45
1:A:96:ARG:HD3	1:A:108:GLN:HB3	1.98	0.45
1:A:28:VAL:HG22	1:A:96:ARG:HG2	1.97	0.45
1:A:470:ASN:ND2	1:A:499:VAL:HG23	2.31	0.45
1:B:283:VAL:HG11	1:B:307:TRP:CE3	2.52	0.45
1:B:367:LEU:HB3	1:B:370:PHE:CD2	2.51	0.44
1:B:244:HIS:CE1	1:B:246:TYR:CZ	3.05	0.44
1:B:448:GLY:N	1:B:452:ASP:O	2.37	0.44
1:A:389:THR:HG22	1:A:390:HIS:ND1	2.33	0.44
1:B:114:ARG:O	1:B:117:LEU:HB2	2.18	0.44
1:B:184:LEU:HD22	1:B:185:PHE:CZ	2.53	0.44
1:B:356:GLN:O	1:B:360:GLN:HG3	2.18	0.44
1:A:251:ASN:OD1	1:A:252:ALA:N	2.51	0.44
1:B:389:THR:HG22	1:B:390:HIS:ND1	2.32	0.44
1:A:87:PRO:O	1:A:89:ILE:HD12	2.16	0.44
1:B:394:ASP:HB3	1:B:397:MET:HB2	2.00	0.44
1:B:500:TRP:CD1	1:B:506:ILE:HG12	2.52	0.44
1:A:315:ASP:HA	1:A:318:GLN:HE21	1.82	0.44
1:B:144:VAL:HG11	1:B:411:LEU:HD21	2.00	0.44
1:A:94:GLU:HG2	1:A:112:PHE:HE1	1.83	0.44
1:A:356:GLN:O	1:A:360:GLN:HG3	2.18	0.43
1:B:204:ASP:OD2	1:B:487:TRP:HH2	2.01	0.43
1:B:346:PRO:HD3	1:B:359:TRP:CE3	2.53	0.43
1:A:213:TYR:CD2	1:A:271:ALA:HB2	2.54	0.43
1:A:63:GLN:HG2	1:A:394:ASP:HA	2.00	0.43
1:B:430:LYS:HA	1:B:431:PRO:HA	1.86	0.43
1:B:26:ARG:NH2	1:B:55:ARG:O	2.52	0.43
1:B:337:PHE:HD2	1:B:393:GLN:HE21	1.66	0.42
1:B:38:GLY:N	1:B:145:GLY:O	2.52	0.42
1:B:213:TYR:CD2	1:B:271:ALA:HB2	2.55	0.42
1:A:334:VAL:HG12	1:A:337:PHE:H	1.84	0.42
1:A:346:PRO:HD3	1:A:359:TRP:CE3	2.54	0.42
1:B:66:ILE:HG13	1:B:358:PHE:CG	2.54	0.42
1:B:334:VAL:HG12	1:B:337:PHE:H	1.85	0.42
1:A:184:LEU:HD22	1:A:185:PHE:CZ	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:ILE:HG13	1:A:358:PHE:CG	2.55	0.42
1:A:283:VAL:HG11	1:A:307:TRP:CE3	2.55	0.41
1:A:308:ALA:HB1	1:A:328:LYS:HA	2.01	0.41
1:B:182:LYS:HZ3	1:B:188:PHE:HB3	1.85	0.41
1:B:280:GLY:HA3	1:B:310:ARG:NH1	2.34	0.41
1:B:515:CYS:HB3	1:B:519:GLN:O	2.20	0.41
1:B:346:PRO:HG3	1:B:359:TRP:CG	2.55	0.41
1:A:214:VAL:HG12	1:A:273:VAL:HB	2.02	0.41
1:B:359:TRP:CD1	1:B:367:LEU:HD21	2.56	0.41
1:B:168:PRO:HA	1:B:187:TYR:HB3	2.03	0.41
1:B:118:ILE:HD12	1:B:142:PRO:HD3	2.03	0.41
1:A:113:ILE:HG22	1:A:167:ILE:HD12	2.03	0.41
1:B:340:TYR:OH	1:B:351:ARG:NH2	2.36	0.41
1:B:72:MET:HA	1:B:404:ILE:HD13	2.03	0.41
1:A:217:VAL:HG22	1:A:246:TYR:HB2	2.02	0.41
1:A:153:SER:O	1:A:157:GLN:NE2	2.54	0.41
1:A:359:TRP:CD1	1:A:367:LEU:HD21	2.55	0.41
1:A:430:LYS:HA	1:A:431:PRO:HA	1.84	0.40
1:B:111:GLU:O	1:B:115:ASP:HB2	2.22	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	502/877 (57%)	471 (94%)	31 (6%)	0	100	100
1	B	481/877 (55%)	453 (94%)	27 (6%)	1 (0%)	49	84
All	All	983/1754 (56%)	924 (94%)	58 (6%)	1 (0%)	53	87

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	521	LYS

5.3.2 Protein sidechains [i](#)

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	412/763 (54%)	411 (100%)	1 (0%)	94	97
1	B	399/763 (52%)	397 (100%)	2 (0%)	90	95
All	All	811/1526 (53%)	808 (100%)	3 (0%)	92	96

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	489	ASN
1	B	465	ARG
1	B	489	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	HIS
1	A	157	GLN
1	A	166	ASN
1	A	198	GLN
1	A	364	GLN
1	B	45	HIS
1	B	74	HIS
1	B	157	GLN
1	B	166	ASN
1	B	198	GLN
1	B	218	HIS
1	B	244	HIS
1	B	364	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	NAG	A	901	1	14,14,15	0.35	0	17,19,21	0.50	0
2	NAG	A	902	1	14,14,15	0.27	0	17,19,21	0.62	0
2	NAG	A	903	1	14,14,15	0.36	0	17,19,21	0.52	0
2	NAG	A	904	1	14,14,15	0.46	0	17,19,21	0.69	1 (5%)
2	NAG	B	902	1	14,14,15	0.36	0	17,19,21	0.65	1 (5%)
2	NAG	B	903	1	14,14,15	0.45	0	17,19,21	0.58	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	901	1	-	0/6/23/26	0/1/1/1
2	NAG	A	902	1	-	0/6/23/26	0/1/1/1
2	NAG	A	903	1	-	0/6/23/26	0/1/1/1
2	NAG	A	904	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	902	1	-	0/6/23/26	0/1/1/1
2	NAG	B	903	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(°)
2	A	904	NAG	C1-O5-C5	2.16	115.15	112.19
2	B	902	NAG	C1-O5-C5	2.19	115.20	112.19

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 [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	506/877 (57%)	-0.19	6 (1%) 79 69	127, 180, 254, 270	0
1	B	487/877 (55%)	-0.20	6 (1%) 79 69	114, 183, 252, 292	0
All	All	993/1754 (56%)	-0.19	12 (1%) 79 69	114, 181, 254, 292	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	54	GLU	4.6
1	B	222	ASN	3.8
1	A	51	LYS	3.5
1	B	372	GLN	3.3
1	A	52	VAL	3.2
1	B	371	PRO	2.9
1	B	221	GLY	2.7
1	A	55	ARG	2.7
1	A	24	GLU	2.5
1	A	53	HIS	2.5
1	B	374	ASN	2.4
1	B	195	ASP	2.3

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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	NAG	A	903	14/15	0.59	0.26	214,233,238,244	0
2	NAG	A	901	14/15	0.62	0.28	200,209,217,221	0
2	NAG	B	902	14/15	0.64	0.27	184,210,226,226	0
2	NAG	A	902	14/15	0.72	0.24	200,219,226,230	0
2	NAG	B	903	14/15	0.84	0.24	222,240,245,248	0
2	NAG	A	904	14/15	0.84	0.16	193,206,212,217	0
3	MG	B	901	1/1	0.92	0.10	165,165,165,165	0

6.5 Other polymers [i](#)

There are no such residues in this entry.