



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 8, 2017 – 07:40 PM EST

PDB ID : 5OJM
Title : Structure of a chimaeric beta3-alpha5 GABAA receptor in complex with nanobody Nb25
Authors : Miller, P.S.; Scott, S.; Masiulis, S.; De Colibus, L.; Pardon, E.; Steyaert, J.; Aricescu, A.R.
Deposited on : unknown
Resolution : 3.30 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	:	rb-20030345
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)	:	rb-20030345

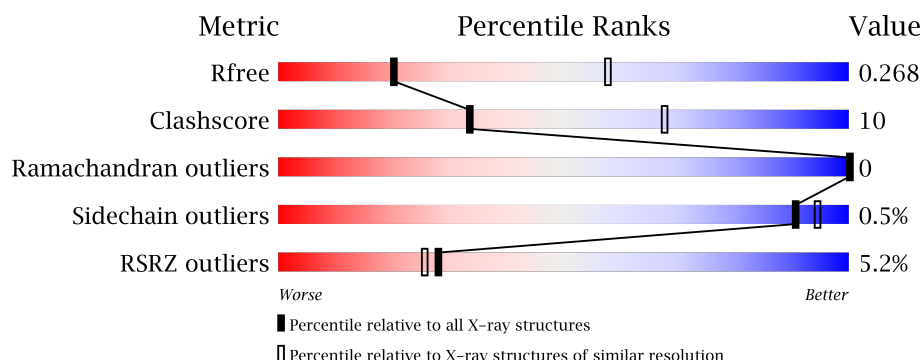
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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	1034 (3.36-3.24)
Clashscore	112137	1100 (3.36-3.24)
Ramachandran outliers	110173	1081 (3.36-3.24)
Sidechain outliers	110143	1080 (3.36-3.24)
RSRZ outliers	101464	1039 (3.36-3.24)

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	395	<div> <div>2%</div> <div>70%14%16%</div> </div>
1	B	395	<div> <div>3%</div> <div>68%16%16%</div> </div>
1	C	395	<div> <div>2%</div> <div>67%16%16%</div> </div>
1	D	395	<div> <div>0%</div> <div>66%18%16%</div> </div>
1	E	395	<div> <div>3%</div> <div>68%16%16%</div> </div>

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Mol	Chain	Length	Quality of chain
2	K	125	
2	L	125	
2	M	125	
2	N	125	
2	O	125	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 18538 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 Human GABAA receptor chimera beta3-alpha5, Gamma-aminobutyric acid receptor subunit beta-3, Gamma-aminobutyric acid receptor subunit alpha-5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	0	0
			2670	1736	442	476	16			
1	B	333	Total	C	N	O	S	0	0	0
			2692	1748	446	481	17			
1	C	330	Total	C	N	O	S	0	0	0
			2670	1736	442	476	16			
1	D	332	Total	C	N	O	S	0	0	0
			2684	1744	444	479	17			
1	E	330	Total	C	N	O	S	0	0	0
			2670	1736	442	476	16			

There are 100 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	316	SER	-	linker	UNP P31644
A	317	GLN	-	linker	UNP P31644
A	318	PRO	-	linker	UNP P31644
A	319	ALA	-	linker	UNP P31644
A	320	ARG	-	linker	UNP P31644
A	321	ALA	-	linker	UNP P31644
A	322	ALA	-	linker	UNP P31644
A	404	ILE	VAL	conflict	UNP P31644
A	428	THR	ALA	conflict	UNP P31644
A	432	GLY	-	expression tag	UNP P31644
A	433	THR	-	expression tag	UNP P31644
A	434	THR	-	expression tag	UNP P31644
A	435	GLU	-	expression tag	UNP P31644
A	436	THR	-	expression tag	UNP P31644
A	437	SER	-	expression tag	UNP P31644
A	438	GLN	-	expression tag	UNP P31644
A	439	VAL	-	expression tag	UNP P31644
A	440	ALA	-	expression tag	UNP P31644

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Chain	Residue	Modelled	Actual	Comment	Reference
A	441	PRO	-	expression tag	UNP P31644
A	442	ALA	-	expression tag	UNP P31644
B	316	SER	-	linker	UNP P31644
B	317	GLN	-	linker	UNP P31644
B	318	PRO	-	linker	UNP P31644
B	319	ALA	-	linker	UNP P31644
B	320	ARG	-	linker	UNP P31644
B	321	ALA	-	linker	UNP P31644
B	322	ALA	-	linker	UNP P31644
B	404	ILE	VAL	conflict	UNP P31644
B	428	THR	ALA	conflict	UNP P31644
B	432	GLY	-	expression tag	UNP P31644
B	433	THR	-	expression tag	UNP P31644
B	434	THR	-	expression tag	UNP P31644
B	435	GLU	-	expression tag	UNP P31644
B	436	THR	-	expression tag	UNP P31644
B	437	SER	-	expression tag	UNP P31644
B	438	GLN	-	expression tag	UNP P31644
B	439	VAL	-	expression tag	UNP P31644
B	440	ALA	-	expression tag	UNP P31644
B	441	PRO	-	expression tag	UNP P31644
B	442	ALA	-	expression tag	UNP P31644
C	316	SER	-	linker	UNP P31644
C	317	GLN	-	linker	UNP P31644
C	318	PRO	-	linker	UNP P31644
C	319	ALA	-	linker	UNP P31644
C	320	ARG	-	linker	UNP P31644
C	321	ALA	-	linker	UNP P31644
C	322	ALA	-	linker	UNP P31644
C	404	ILE	VAL	conflict	UNP P31644
C	428	THR	ALA	conflict	UNP P31644
C	432	GLY	-	expression tag	UNP P31644
C	433	THR	-	expression tag	UNP P31644
C	434	THR	-	expression tag	UNP P31644
C	435	GLU	-	expression tag	UNP P31644
C	436	THR	-	expression tag	UNP P31644
C	437	SER	-	expression tag	UNP P31644
C	438	GLN	-	expression tag	UNP P31644
C	439	VAL	-	expression tag	UNP P31644
C	440	ALA	-	expression tag	UNP P31644
C	441	PRO	-	expression tag	UNP P31644
C	442	ALA	-	expression tag	UNP P31644

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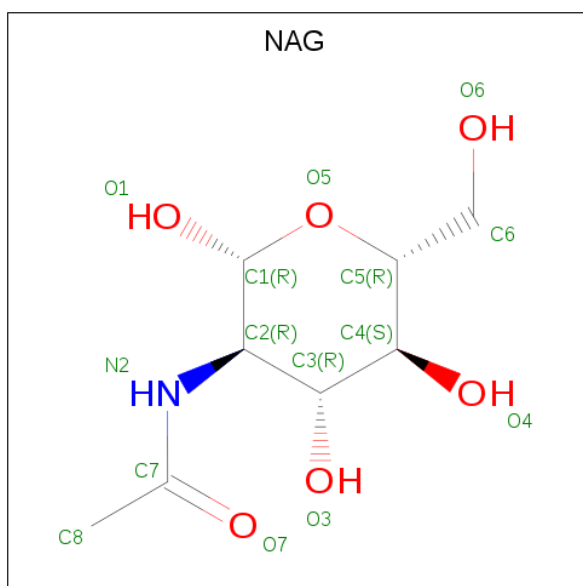
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Chain	Residue	Modelled	Actual	Comment	Reference
D	316	SER	-	linker	UNP P31644
D	317	GLN	-	linker	UNP P31644
D	318	PRO	-	linker	UNP P31644
D	319	ALA	-	linker	UNP P31644
D	320	ARG	-	linker	UNP P31644
D	321	ALA	-	linker	UNP P31644
D	322	ALA	-	linker	UNP P31644
D	404	ILE	VAL	conflict	UNP P31644
D	428	THR	ALA	conflict	UNP P31644
D	432	GLY	-	expression tag	UNP P31644
D	433	THR	-	expression tag	UNP P31644
D	434	THR	-	expression tag	UNP P31644
D	435	GLU	-	expression tag	UNP P31644
D	436	THR	-	expression tag	UNP P31644
D	437	SER	-	expression tag	UNP P31644
D	438	GLN	-	expression tag	UNP P31644
D	439	VAL	-	expression tag	UNP P31644
D	440	ALA	-	expression tag	UNP P31644
D	441	PRO	-	expression tag	UNP P31644
D	442	ALA	-	expression tag	UNP P31644
E	316	SER	-	linker	UNP P31644
E	317	GLN	-	linker	UNP P31644
E	318	PRO	-	linker	UNP P31644
E	319	ALA	-	linker	UNP P31644
E	320	ARG	-	linker	UNP P31644
E	321	ALA	-	linker	UNP P31644
E	322	ALA	-	linker	UNP P31644
E	404	ILE	VAL	conflict	UNP P31644
E	428	THR	ALA	conflict	UNP P31644
E	432	GLY	-	expression tag	UNP P31644
E	433	THR	-	expression tag	UNP P31644
E	434	THR	-	expression tag	UNP P31644
E	435	GLU	-	expression tag	UNP P31644
E	436	THR	-	expression tag	UNP P31644
E	437	SER	-	expression tag	UNP P31644
E	438	GLN	-	expression tag	UNP P31644
E	439	VAL	-	expression tag	UNP P31644
E	440	ALA	-	expression tag	UNP P31644
E	441	PRO	-	expression tag	UNP P31644
E	442	ALA	-	expression tag	UNP P31644

- Molecule 2 is a protein called Nanobody Nb25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	K	123	Total	C	N	O	S	0	0	0
			951	598	164	185	4			
2	L	123	Total	C	N	O	S	0	0	0
			951	598	164	185	4			
2	M	123	Total	C	N	O	S	0	0	0
			951	598	164	185	4			
2	N	123	Total	C	N	O	S	0	0	0
			951	598	164	185	4			
2	O	123	Total	C	N	O	S	0	0	0
			951	598	164	185	4			

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



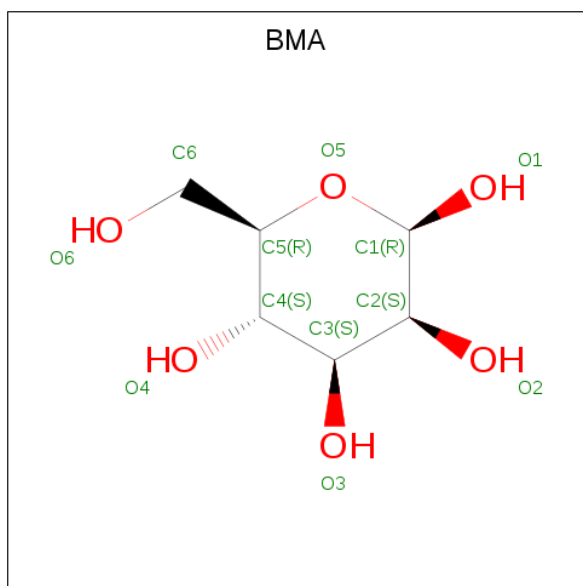
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		

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

- Molecule 4 is BETA-D-MANNOSE (three-letter code: BMA) (formula: $C_6H_{12}O_6$).



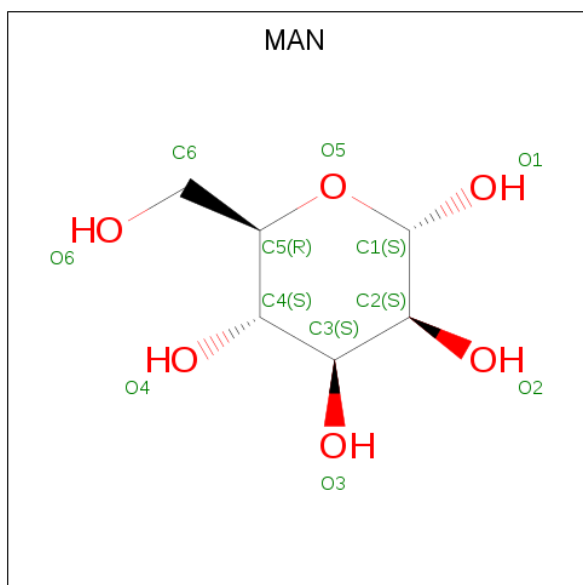
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			11	6	5		
4	B	1	Total	C	O	0	0
			11	6	5		
4	C	1	Total	C	O	0	0
			11	6	5		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	C	O	0	0
			11	6	5		
4	E	1	Total	C	O	0	0
			11	6	5		

- Molecule 5 is ALPHA-D-MANNOSE (three-letter code: MAN) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			11	6	5		
5	A	1	Total	C	O	0	0
			11	6	5		
5	B	1	Total	C	O	0	0
			11	6	5		
5	B	1	Total	C	O	0	0
			11	6	5		
5	C	1	Total	C	O	0	0
			11	6	5		
5	C	1	Total	C	O	0	0
			11	6	5		
5	D	1	Total	C	O	0	0
			11	6	5		
5	D	1	Total	C	O	0	0
			11	6	5		
5	D	1	Total	C	O	0	0
			11	6	5		

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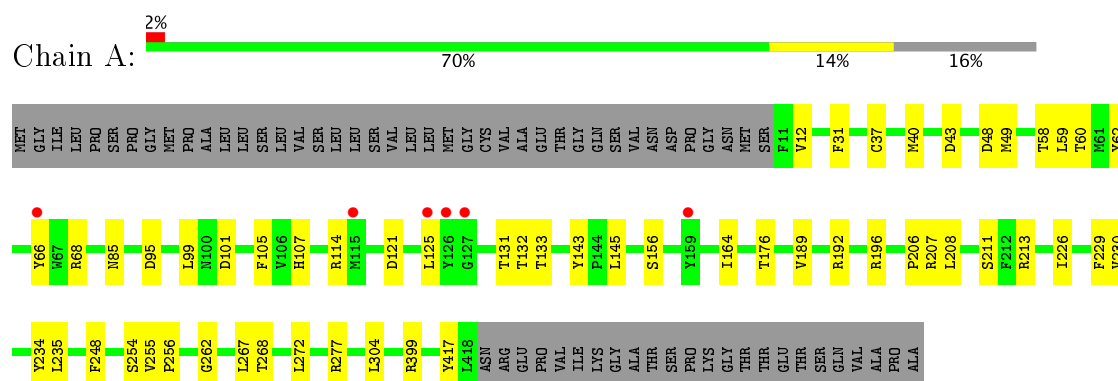
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	E	1	Total	C	O	0	0
			11	6	5		
5	E	1	Total	C	O	0	0
			11	6	5		
5	E	1	Total	C	O	0	0
			11	6	5		

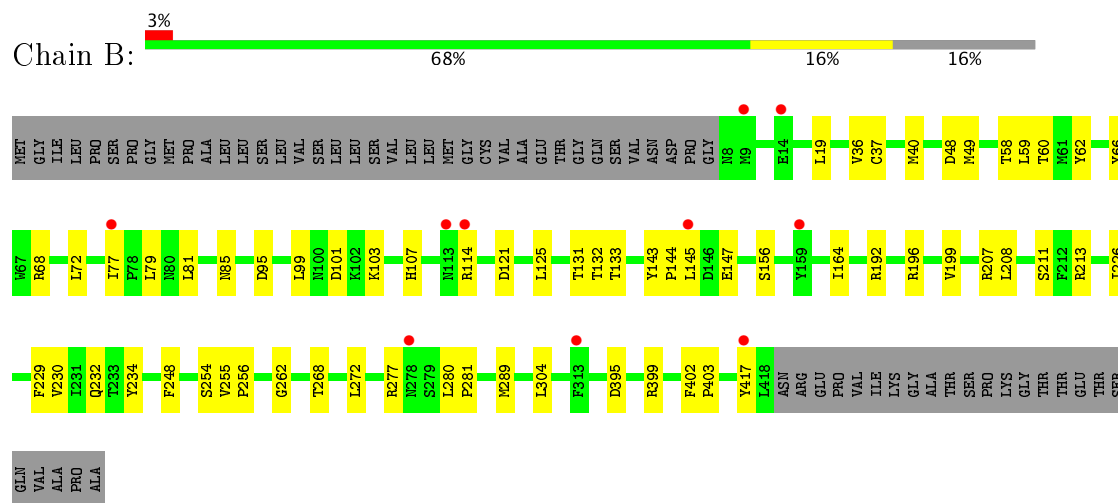
3 Residue-property plots [i](#)

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.

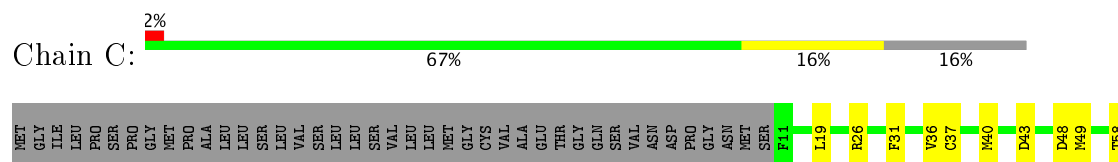
- Molecule 1: Human GABAA receptor chimera beta3-alpha5, Gamma-aminobutyric acid receptor subunit beta-3, Gamma-aminobutyric acid receptor subunit alpha-5

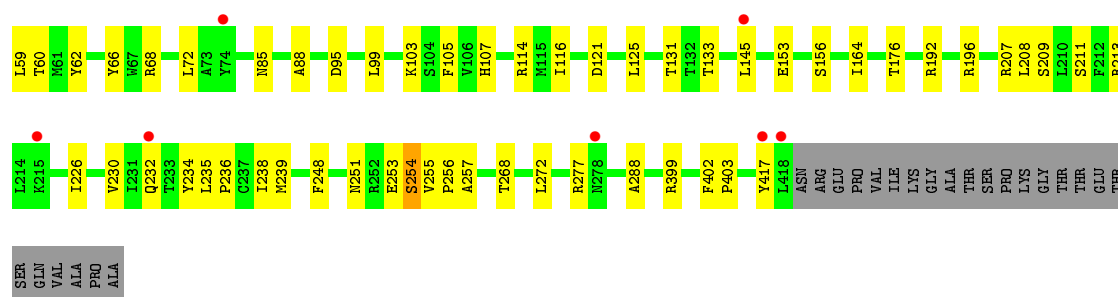


- Molecule 1: Human GABAA receptor chimera beta3-alpha5, Gamma-aminobutyric acid receptor subunit beta-3, Gamma-aminobutyric acid receptor subunit alpha-5

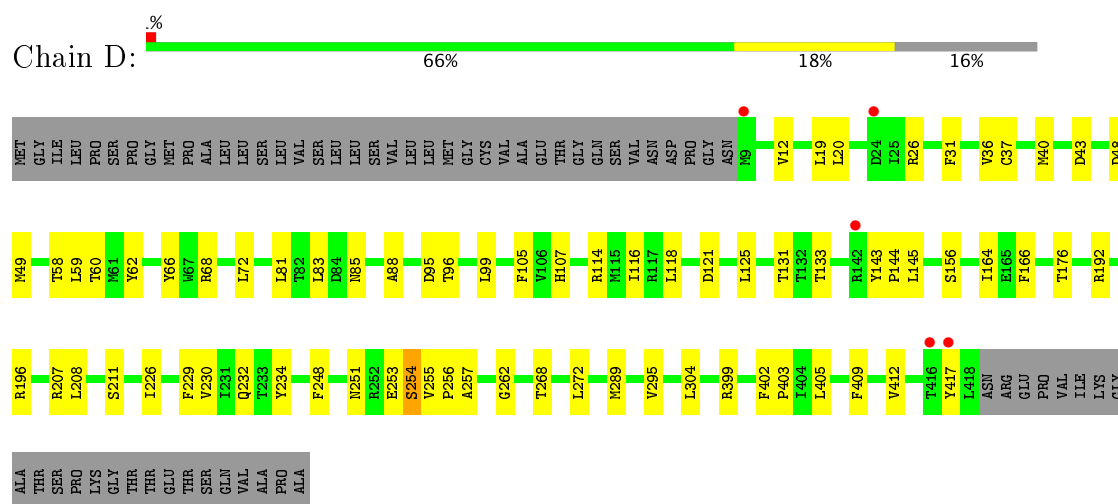


- Molecule 1: Human GABAA receptor chimera beta3-alpha5, Gamma-aminobutyric acid receptor subunit beta-3, Gamma-aminobutyric acid receptor subunit alpha-5

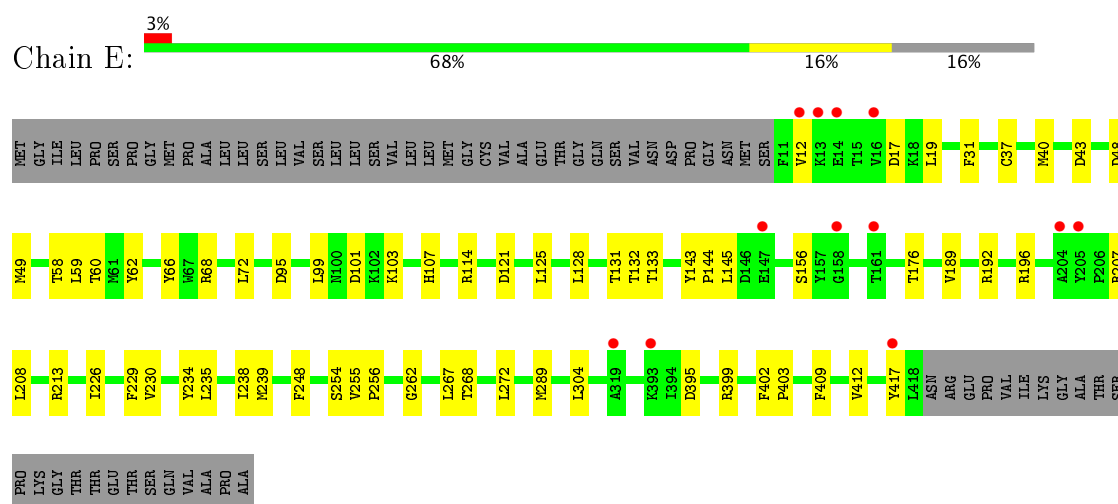




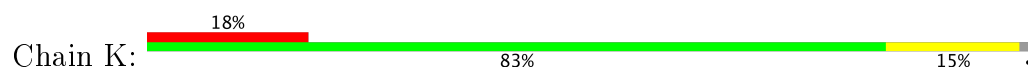
- Molecule 1: Human GABAA receptor chimera beta3-alpha5, Gamma-aminobutyric acid receptor subunit beta-3, Gamma-aminobutyric acid receptor subunit alpha-5

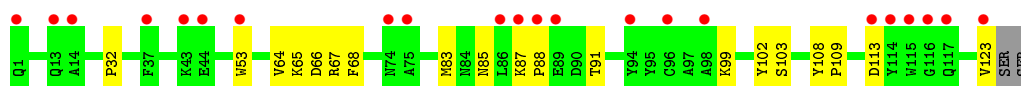


- Molecule 1: Human GABAA receptor chimera beta3-alpha5, Gamma-aminobutyric acid receptor subunit beta-3, Gamma-aminobutyric acid receptor subunit alpha-5

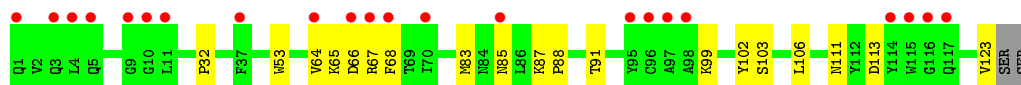
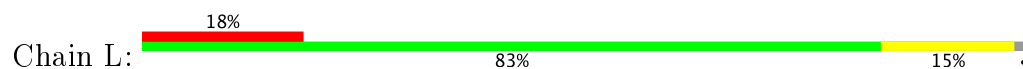


- Molecule 2: Nanobody Nb25

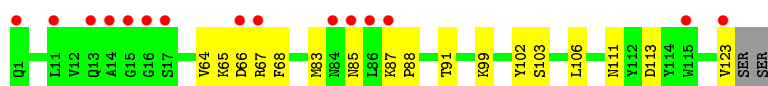
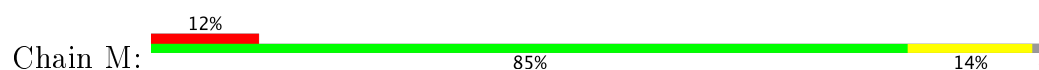




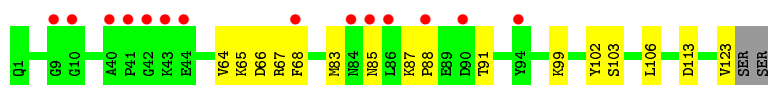
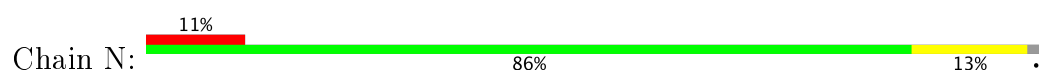
- Molecule 2: Nanobody Nb25



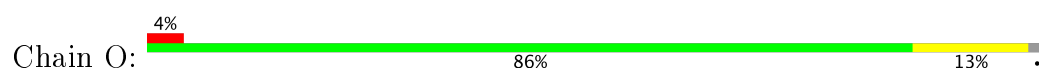
- Molecule 2: Nanobody Nb25



- Molecule 2: Nanobody Nb25



- Molecule 2: Nanobody Nb25



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	177.35Å 139.94Å 191.50Å 90.00° 102.25° 90.00°	Depositor
Resolution (Å)	48.88 – 3.30 48.88 – 3.28	Depositor EDS
% Data completeness (in resolution range)	73.6 (48.88-3.30) 99.1 (48.88-3.28)	Depositor EDS
R_{merge}	0.35	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.17 (at 3.25Å)	Xtriage
Refinement program	REFMAC 5.8.0158, PHENIX (1.12rc)-2787-000)	Depositor
R, R_{free}	0.234 , 0.250 0.251 , 0.268	Depositor DCC
R_{free} test set	3382 reflections (5.16%)	DCC
Wilson B-factor (Å ²)	85.8	Xtriage
Anisotropy	0.424	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 9.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	18538	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.26% 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: BMA, NAG, MAN

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	0.50	0/2737	0.44	0/3724
1	B	0.50	0/2759	0.43	0/3753
1	C	0.50	0/2737	0.44	0/3724
1	D	0.52	0/2751	0.44	0/3742
1	E	0.50	0/2737	0.44	0/3724
2	K	0.41	0/975	0.38	0/1321
2	L	0.38	0/975	0.37	0/1321
2	M	0.40	0/975	0.37	0/1321
2	N	0.41	0/975	0.37	0/1321
2	O	0.39	0/975	0.37	0/1321
All	All	0.48	0/18596	0.42	0/25272

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	2670	0	2678	60	0
1	B	2692	0	2698	70	0
1	C	2670	0	2678	66	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2684	0	2692	67	0
1	E	2670	0	2678	63	0
2	K	951	0	895	15	0
2	L	951	0	895	16	0
2	M	951	0	895	15	0
2	N	951	0	895	14	0
2	O	951	0	895	14	0
3	A	42	0	37	0	0
3	B	42	0	37	0	0
3	C	42	0	37	1	0
3	D	42	0	37	0	0
3	E	42	0	37	1	0
4	A	11	0	8	0	0
4	B	11	0	8	0	0
4	C	11	0	8	0	0
4	D	11	0	8	0	0
4	E	11	0	8	1	0
5	A	22	0	20	0	0
5	B	22	0	20	0	0
5	C	22	0	20	0	0
5	D	33	0	29	0	0
5	E	33	0	29	1	0
All	All	18538	0	18242	355	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 (355) 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:62:TYR:OH	1:E:99:LEU:CD1	2.22	0.87
1:D:207:ARG:NH1	2:N:102:TYR:O	2.07	0.87
1:A:99:LEU:CD1	1:B:62:TYR:OH	2.24	0.86
1:B:143:TYR:OH	1:B:229:PHE:HE2	1.57	0.85
1:C:196:ARG:NH1	2:M:113:ASP:OD1	2.10	0.83
1:B:196:ARG:NH1	2:L:113:ASP:OD1	2.10	0.83
1:A:99:LEU:HD12	1:B:62:TYR:OH	1.79	0.82
1:D:99:LEU:CD1	1:E:62:TYR:OH	2.28	0.81
1:B:99:LEU:CD1	1:C:62:TYR:OH	2.28	0.81
1:C:99:LEU:CD1	1:D:62:TYR:OH	2.31	0.79
1:A:207:ARG:NH1	2:K:102:TYR:O	2.15	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:64:VAL:HG13	2:O:68:PHE:CD1	2.18	0.79
1:B:207:ARG:NH1	2:L:102:TYR:O	2.15	0.78
2:L:64:VAL:HG13	2:L:68:PHE:CD1	2.18	0.78
2:M:64:VAL:HG13	2:M:68:PHE:CD1	2.18	0.78
1:E:196:ARG:NH1	2:O:113:ASP:OD1	2.15	0.78
2:N:64:VAL:HG13	2:N:68:PHE:CD1	2.18	0.78
1:E:99:LEU:HD22	1:E:207:ARG:NH2	1.99	0.78
2:K:64:VAL:HG13	2:K:68:PHE:CD1	2.18	0.77
1:B:99:LEU:HD22	1:B:207:ARG:NH2	1.99	0.77
1:C:99:LEU:HD22	1:C:207:ARG:NH2	2.00	0.77
1:C:207:ARG:NH1	2:M:102:TYR:O	2.19	0.76
1:A:99:LEU:HD22	1:A:207:ARG:NH2	2.00	0.76
1:E:207:ARG:NH1	2:O:102:TYR:O	2.18	0.76
1:D:49:MET:HB3	1:D:58:THR:HG22	1.68	0.76
1:B:49:MET:HB3	1:B:58:THR:HG22	1.68	0.76
1:A:62:TYR:OH	1:E:99:LEU:HD12	1.87	0.75
1:D:99:LEU:HD22	1:D:207:ARG:NH2	2.02	0.75
1:D:99:LEU:HD12	1:E:62:TYR:OH	1.88	0.73
1:C:49:MET:HB3	1:C:58:THR:HG22	1.71	0.72
1:E:49:MET:HB3	1:E:58:THR:HG22	1.71	0.72
1:E:192:ARG:NH1	3:E:502:NAG:O7	2.25	0.70
1:D:196:ARG:NH1	2:N:113:ASP:OD1	2.23	0.70
1:A:49:MET:HB3	1:A:58:THR:HG22	1.71	0.70
1:B:99:LEU:HD12	1:C:62:TYR:OH	1.92	0.70
1:B:145:LEU:CD2	1:B:226:ILE:CD1	2.72	0.67
1:E:60:THR:OG1	1:E:131:THR:HG22	1.95	0.67
1:A:145:LEU:CD2	1:A:226:ILE:CD1	2.72	0.67
1:D:145:LEU:CD2	1:D:226:ILE:CD1	2.73	0.66
1:A:60:THR:OG1	1:A:131:THR:HG22	1.95	0.66
1:A:248:PHE:O	1:A:399:ARG:NE	2.28	0.66
1:C:145:LEU:CD2	1:C:226:ILE:CD1	2.73	0.66
1:C:99:LEU:HD12	1:D:62:TYR:OH	1.94	0.66
1:C:60:THR:OG1	1:C:131:THR:HG22	1.96	0.66
1:E:248:PHE:O	1:E:399:ARG:NE	2.29	0.66
1:C:248:PHE:O	1:C:399:ARG:NE	2.29	0.65
1:D:19:LEU:HD22	1:D:72:LEU:HA	1.79	0.65
1:B:60:THR:OG1	1:B:131:THR:HG22	1.97	0.65
1:C:153:GLU:OE1	1:C:209:SER:OG	2.14	0.65
1:D:248:PHE:O	1:D:399:ARG:NE	2.30	0.65
1:D:60:THR:OG1	1:D:131:THR:HG22	1.96	0.65
1:D:58:THR:HB	1:D:133:THR:HG22	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:145:LEU:CD2	1:E:226:ILE:CD1	2.74	0.64
2:O:91:THR:HG22	2:O:123:VAL:H	1.63	0.64
1:C:58:THR:CB	1:C:133:THR:HG22	2.29	0.63
1:D:40:MET:HG3	1:D:208:LEU:HD12	1.80	0.63
2:K:91:THR:HG22	2:K:123:VAL:H	1.63	0.63
1:C:58:THR:HB	1:C:133:THR:HG22	1.80	0.63
2:L:91:THR:HG22	2:L:123:VAL:H	1.63	0.63
2:M:64:VAL:HG13	2:M:68:PHE:CG	2.33	0.63
1:E:40:MET:HG3	1:E:208:LEU:HD12	1.81	0.63
1:D:58:THR:CB	1:D:133:THR:HG22	2.29	0.62
1:A:58:THR:CB	1:A:133:THR:HG22	2.30	0.62
1:A:58:THR:HB	1:A:133:THR:HG22	1.80	0.62
1:B:143:TYR:OH	1:B:229:PHE:CE2	2.41	0.62
2:M:91:THR:HG22	2:M:123:VAL:H	1.65	0.62
2:O:64:VAL:HG13	2:O:68:PHE:CG	2.34	0.62
2:K:64:VAL:HG13	2:K:68:PHE:CG	2.34	0.61
1:D:99:LEU:HD12	1:E:62:TYR:HH	1.65	0.61
1:B:58:THR:HB	1:B:133:THR:HG22	1.83	0.61
2:L:64:VAL:HG13	2:L:68:PHE:CG	2.36	0.61
1:B:49:MET:HB3	1:B:58:THR:CG2	2.31	0.60
2:N:91:THR:HG22	2:N:123:VAL:H	1.65	0.60
2:N:64:VAL:HG13	2:N:68:PHE:CG	2.35	0.60
1:E:58:THR:CB	1:E:133:THR:HG22	2.32	0.60
1:A:164:ILE:HG23	1:A:206:PRO:HG3	1.81	0.60
1:D:68:ARG:NH1	1:D:121:ASP:OD1	2.34	0.60
1:E:68:ARG:NH1	1:E:121:ASP:OD1	2.35	0.60
1:C:68:ARG:NH1	1:C:121:ASP:OD1	2.34	0.60
1:B:68:ARG:NH1	1:B:121:ASP:OD1	2.34	0.59
1:A:68:ARG:NH1	1:A:121:ASP:OD1	2.35	0.59
1:C:145:LEU:CD2	1:C:226:ILE:HD13	2.31	0.59
1:A:62:TYR:OH	1:E:99:LEU:HD11	2.03	0.59
1:B:40:MET:HG3	1:B:208:LEU:HD12	1.85	0.59
1:C:232:GLN:O	1:C:236:PRO:HG2	2.03	0.59
1:D:49:MET:HB3	1:D:58:THR:CG2	2.33	0.58
1:B:58:THR:CB	1:B:133:THR:HG22	2.32	0.58
1:B:19:LEU:HD22	1:B:72:LEU:HA	1.84	0.58
1:C:40:MET:HG3	1:C:208:LEU:HD12	1.84	0.58
1:E:58:THR:HB	1:E:133:THR:HG22	1.84	0.58
1:E:99:LEU:HD22	1:E:207:ARG:CZ	2.33	0.58
1:E:145:LEU:CD2	1:E:226:ILE:HD13	2.33	0.58
1:E:49:MET:HB3	1:E:58:THR:CG2	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:99:LEU:HD22	1:C:207:ARG:CZ	2.33	0.58
1:A:196:ARG:NH1	2:K:113:ASP:OD1	2.37	0.57
2:O:67:ARG:NH2	2:O:85:ASN:O	2.37	0.57
1:C:255:VAL:N	1:C:256:PRO:HD2	2.19	0.57
1:B:255:VAL:N	1:B:256:PRO:HD2	2.19	0.57
1:E:255:VAL:N	1:E:256:PRO:HD2	2.19	0.57
1:C:19:LEU:HD22	1:C:72:LEU:HA	1.86	0.57
1:B:145:LEU:HD23	1:B:226:ILE:CD1	2.35	0.57
1:D:255:VAL:N	1:D:256:PRO:HD2	2.19	0.57
1:E:19:LEU:HD22	1:E:72:LEU:HA	1.86	0.57
1:A:255:VAL:N	1:A:256:PRO:HD2	2.19	0.57
2:L:67:ARG:NH2	2:L:85:ASN:O	2.38	0.57
1:A:99:LEU:HD22	1:A:207:ARG:CZ	2.34	0.57
1:D:88:ALA:HB2	1:D:116:ILE:HD11	1.87	0.57
1:D:99:LEU:HD22	1:D:207:ARG:CZ	2.35	0.57
1:D:145:LEU:CD2	1:D:226:ILE:HD13	2.35	0.57
1:C:288:ALA:HB2	1:C:417:TYR:CE2	2.40	0.56
1:A:40:MET:HG3	1:A:208:LEU:HD12	1.86	0.56
1:E:235:LEU:HD23	1:E:238:ILE:HD11	1.86	0.56
2:M:67:ARG:NH2	2:M:85:ASN:O	2.38	0.56
2:N:67:ARG:NH2	2:N:85:ASN:O	2.38	0.56
2:K:67:ARG:NH2	2:K:85:ASN:O	2.38	0.56
1:E:226:ILE:HG22	1:E:226:ILE:O	2.06	0.55
1:C:226:ILE:HG22	1:C:226:ILE:O	2.06	0.55
1:B:277:ARG:HB3	1:C:232:GLN:HE21	1.71	0.55
1:B:145:LEU:CD2	1:B:226:ILE:HD11	2.37	0.54
1:B:145:LEU:CD2	1:B:226:ILE:HD13	2.37	0.54
1:D:68:ARG:NH1	1:D:121:ASP:O	2.37	0.54
1:A:145:LEU:HD23	1:A:226:ILE:CD1	2.37	0.54
1:A:226:ILE:HG22	1:A:226:ILE:O	2.07	0.54
2:N:65:LYS:O	2:N:66:ASP:HB2	2.08	0.54
1:A:248:PHE:O	1:A:399:ARG:NH2	2.40	0.54
1:C:145:LEU:HD23	1:C:226:ILE:CD1	2.38	0.54
1:A:49:MET:HB3	1:A:58:THR:CG2	2.36	0.54
1:D:143:TYR:HH	1:D:229:PHE:HE1	1.56	0.54
1:C:248:PHE:O	1:C:399:ARG:NH2	2.41	0.54
2:M:65:LYS:O	2:M:66:ASP:HB2	2.07	0.54
1:A:145:LEU:CD2	1:A:226:ILE:HD13	2.38	0.53
1:B:226:ILE:HG22	1:B:226:ILE:O	2.07	0.53
1:C:277:ARG:HB3	1:D:232:GLN:HE21	1.73	0.53
1:B:99:LEU:HD22	1:B:207:ARG:CZ	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:68:PHE:HE2	2:N:83:MET:SD	2.31	0.53
1:B:103:LYS:HE3	1:C:105:PHE:CE2	2.44	0.53
2:K:68:PHE:CE2	2:K:83:MET:SD	3.02	0.53
1:C:49:MET:HB3	1:C:58:THR:CG2	2.36	0.53
1:D:226:ILE:HG22	1:D:226:ILE:O	2.07	0.53
1:D:248:PHE:O	1:D:399:ARG:NH2	2.41	0.53
2:O:68:PHE:HE2	2:O:83:MET:SD	2.32	0.53
2:L:65:LYS:O	2:L:66:ASP:HB2	2.09	0.53
2:L:68:PHE:CE2	2:L:83:MET:SD	3.03	0.53
2:N:68:PHE:CE2	2:N:83:MET:SD	3.03	0.53
2:K:68:PHE:HE2	2:K:83:MET:SD	2.31	0.52
2:L:68:PHE:HE2	2:L:83:MET:SD	2.31	0.52
1:C:230:VAL:HA	1:C:234:TYR:HB2	1.91	0.52
2:K:65:LYS:O	2:K:66:ASP:HB2	2.07	0.52
2:M:68:PHE:HE2	2:M:83:MET:SD	2.32	0.52
2:O:68:PHE:CE2	2:O:83:MET:SD	3.03	0.52
1:A:145:LEU:CD2	1:A:226:ILE:HD11	2.37	0.52
1:B:143:TYR:HH	1:B:229:PHE:HE2	0.74	0.52
2:M:68:PHE:CE2	2:M:83:MET:SD	3.03	0.52
2:O:65:LYS:O	2:O:66:ASP:HB2	2.08	0.52
1:E:48:ASP:HB2	1:E:58:THR:HG23	1.92	0.52
1:B:277:ARG:HB3	1:C:232:GLN:NE2	2.24	0.52
1:C:235:LEU:HD23	1:C:238:ILE:HD11	1.91	0.51
1:D:145:LEU:HD23	1:D:226:ILE:CD1	2.40	0.51
1:D:48:ASP:HB2	1:D:58:THR:CG2	2.40	0.51
1:E:248:PHE:O	1:E:399:ARG:NH2	2.42	0.51
2:O:91:THR:HG22	2:O:123:VAL:N	2.25	0.51
1:C:48:ASP:HB2	1:C:58:THR:CG2	2.41	0.51
1:E:68:ARG:NH1	1:E:121:ASP:O	2.38	0.51
2:L:91:THR:HG22	2:L:123:VAL:N	2.25	0.51
1:B:68:ARG:NH1	1:B:121:ASP:O	2.38	0.51
1:B:48:ASP:HB2	1:B:58:THR:HG23	1.93	0.51
2:M:91:THR:HG22	2:M:123:VAL:N	2.26	0.51
2:K:91:THR:HG22	2:K:123:VAL:N	2.26	0.51
1:B:48:ASP:HB2	1:B:58:THR:CG2	2.41	0.51
1:E:114:ARG:HB2	1:E:128:LEU:HD23	1.93	0.51
1:D:48:ASP:HB2	1:D:58:THR:HG23	1.92	0.50
1:D:230:VAL:HA	1:D:234:TYR:HB2	1.94	0.50
1:A:230:VAL:HA	1:A:234:TYR:HB2	1.92	0.50
1:E:145:LEU:HD23	1:E:226:ILE:CD1	2.40	0.50
1:E:143:TYR:HH	1:E:229:PHE:HE1	1.60	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:ASP:HB2	1:A:58:THR:CG2	2.42	0.50
1:D:66:TYR:CZ	1:D:125:LEU:HD13	2.47	0.50
1:C:48:ASP:HB2	1:C:58:THR:HG23	1.93	0.50
1:A:277:ARG:HB3	1:B:232:GLN:HE21	1.76	0.49
1:D:96:THR:HG22	1:D:166:PHE:CZ	2.47	0.49
1:A:277:ARG:HB3	1:B:232:GLN:NE2	2.28	0.49
1:C:59:LEU:O	1:C:131:THR:HA	2.13	0.49
2:N:91:THR:HG22	2:N:123:VAL:N	2.26	0.49
1:D:226:ILE:HG22	1:D:230:VAL:HG23	1.94	0.49
1:B:230:VAL:HA	1:B:234:TYR:HB2	1.95	0.49
1:B:77:ILE:HG21	1:B:81:LEU:HD21	1.94	0.49
1:A:48:ASP:HB2	1:A:58:THR:HG23	1.93	0.49
1:E:48:ASP:HB2	1:E:58:THR:CG2	2.41	0.49
2:N:87:LYS:HB3	2:N:88:PRO:HD2	1.94	0.49
1:A:189:VAL:HG12	1:A:213:ARG:HB3	1.94	0.48
1:C:235:LEU:HA	1:C:238:ILE:HG12	1.94	0.48
1:C:66:TYR:CZ	1:C:125:LEU:HD13	2.48	0.48
1:D:59:LEU:O	1:D:131:THR:HA	2.13	0.48
2:K:87:LYS:HB3	2:K:88:PRO:HD2	1.94	0.48
1:C:251:ASN:OD1	1:C:253:GLU:HB2	2.13	0.48
1:B:59:LEU:O	1:B:131:THR:HA	2.14	0.48
1:E:230:VAL:HA	1:E:234:TYR:HB2	1.94	0.48
1:E:59:LEU:O	1:E:131:THR:HA	2.14	0.48
1:C:85:ASN:HB2	1:C:114:ARG:HG3	1.95	0.48
1:C:103:LYS:HE3	1:D:105:PHE:CE2	2.48	0.48
1:B:268:THR:O	1:B:272:LEU:HG	2.14	0.48
1:C:68:ARG:NH1	1:C:121:ASP:O	2.39	0.48
1:C:31:PHE:CE2	1:D:12:VAL:HG11	2.49	0.48
1:D:99:LEU:HD11	1:E:62:TYR:OH	2.14	0.48
1:A:62:TYR:HH	1:E:99:LEU:HD12	1.77	0.47
2:O:87:LYS:HB3	2:O:88:PRO:HD2	1.95	0.47
1:D:145:LEU:CD1	1:D:417:TYR:HB3	2.44	0.47
2:M:87:LYS:HB3	2:M:88:PRO:HD2	1.95	0.47
1:D:145:LEU:CD2	1:D:226:ILE:HD11	2.43	0.47
2:L:87:LYS:HB3	2:L:88:PRO:HD2	1.96	0.47
1:A:59:LEU:O	1:A:131:THR:HA	2.14	0.47
1:C:145:LEU:CD2	1:C:226:ILE:HD11	2.45	0.47
1:A:12:VAL:HG11	1:E:31:PHE:CE2	2.50	0.47
1:C:268:THR:O	1:C:272:LEU:HG	2.15	0.46
1:D:145:LEU:HD23	1:D:226:ILE:HG12	1.97	0.46
1:D:251:ASN:OD1	1:D:253:GLU:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:288:ALA:CB	1:C:417:TYR:CE2	2.99	0.46
1:E:235:LEU:HA	1:E:238:ILE:HG12	1.96	0.46
1:B:36:VAL:HB	1:B:164:ILE:HD13	1.97	0.46
1:D:107:HIS:NE2	1:D:131:THR:HG23	2.30	0.46
1:B:145:LEU:CD1	1:B:417:TYR:HB3	2.46	0.46
1:B:143:TYR:CZ	1:B:229:PHE:CE2	3.03	0.46
1:B:79:LEU:HD12	1:B:81:LEU:HD23	1.97	0.46
1:A:66:TYR:CZ	1:A:125:LEU:HD13	2.51	0.46
1:A:143:TYR:HH	1:A:229:PHE:HE1	1.63	0.46
1:A:31:PHE:CE2	1:B:81:LEU:HD22	2.51	0.46
1:A:68:ARG:NH1	1:A:121:ASP:O	2.38	0.46
1:A:255:VAL:HB	1:B:254:SER:OG	2.15	0.46
1:B:107:HIS:NE2	1:B:131:THR:HG23	2.31	0.46
1:A:145:LEU:CD1	1:A:417:TYR:HB3	2.46	0.45
1:E:66:TYR:CZ	1:E:125:LEU:HD13	2.51	0.45
2:K:99:LYS:HG2	2:K:103:SER:HB2	1.99	0.45
2:O:68:PHE:CE2	2:O:83:MET:HB3	2.52	0.45
1:A:85:ASN:HB2	1:A:114:ARG:CG	2.47	0.45
1:C:26:ARG:NH1	1:D:20:LEU:HD12	2.31	0.45
1:E:268:THR:O	1:E:272:LEU:HG	2.16	0.45
1:E:145:LEU:CD1	1:E:417:TYR:HB3	2.46	0.45
1:B:66:TYR:CZ	1:B:125:LEU:HD13	2.51	0.45
4:E:504:BMA:H62	5:E:505:MAN:H2	1.77	0.45
2:K:68:PHE:CE2	2:K:83:MET:HB3	2.51	0.45
1:D:85:ASN:HB2	1:D:114:ARG:CG	2.46	0.45
2:N:68:PHE:CE2	2:N:83:MET:HB3	2.51	0.45
1:E:145:LEU:HD23	1:E:226:ILE:HG12	1.98	0.45
1:C:95:ASP:OD1	1:C:156:SER:HB2	2.16	0.45
1:D:85:ASN:HB2	1:D:114:ARG:HG3	1.98	0.45
2:M:68:PHE:CE2	2:M:83:MET:HB3	2.52	0.45
1:A:107:HIS:NE2	1:A:131:THR:HG23	2.31	0.44
1:D:192:ARG:HB2	1:D:211:SER:OG	2.17	0.44
1:E:226:ILE:HG22	1:E:230:VAL:HG23	2.00	0.44
1:B:144:PRO:HD3	1:B:289:MET:HB2	1.99	0.44
1:E:262:GLY:HA3	1:E:304:LEU:HD13	2.00	0.44
1:E:95:ASP:OD1	1:E:156:SER:HB2	2.17	0.44
1:B:66:TYR:CE2	1:B:125:LEU:HD13	2.53	0.44
1:B:95:ASP:OD1	1:B:156:SER:HB2	2.17	0.44
1:C:226:ILE:HG22	1:C:230:VAL:HG23	1.99	0.44
1:C:248:PHE:O	1:C:399:ARG:CZ	2.65	0.44
1:A:268:THR:O	1:A:272:LEU:HG	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:36:VAL:HB	1:D:164:ILE:HD13	1.98	0.44
2:M:99:LYS:HG2	2:M:103:SER:HB2	1.99	0.44
1:A:248:PHE:O	1:A:399:ARG:CZ	2.66	0.44
1:B:226:ILE:HG22	1:B:230:VAL:HG23	2.00	0.44
1:D:95:ASP:OD1	1:D:156:SER:HB2	2.17	0.44
1:D:268:THR:O	1:D:272:LEU:HG	2.18	0.44
1:D:66:TYR:CE2	1:D:125:LEU:HD13	2.52	0.44
1:E:107:HIS:NE2	1:E:131:THR:HG23	2.33	0.44
2:O:99:LYS:HG2	2:O:103:SER:HB2	1.99	0.44
1:B:395:ASP:O	1:B:399:ARG:HG3	2.18	0.44
2:L:68:PHE:CE2	2:L:83:MET:HB3	2.52	0.44
1:A:145:LEU:HD23	1:A:226:ILE:HG12	1.99	0.44
1:A:58:THR:O	1:A:58:THR:HG23	2.18	0.44
1:D:31:PHE:CE2	1:E:12:VAL:HG11	2.53	0.44
2:L:99:LYS:HG2	2:L:103:SER:HB2	2.00	0.44
1:A:235:LEU:HA	1:A:235:LEU:HD23	1.89	0.43
1:A:95:ASP:OD1	1:A:156:SER:HB2	2.17	0.43
1:B:145:LEU:HD23	1:B:226:ILE:HD11	2.00	0.43
1:B:147:GLU:OE1	1:B:213:ARG:NH1	2.50	0.43
1:E:145:LEU:CD2	1:E:226:ILE:HD11	2.46	0.43
1:C:145:LEU:CD1	1:C:417:TYR:HB3	2.48	0.43
1:A:226:ILE:HG22	1:A:230:VAL:HG23	2.00	0.43
1:A:99:LEU:HD12	1:B:62:TYR:HH	1.82	0.43
1:C:402:PHE:HB2	1:C:403:PRO:HD3	2.00	0.43
1:E:402:PHE:HB2	1:E:403:PRO:HD3	2.00	0.43
1:C:58:THR:O	1:C:58:THR:HG23	2.18	0.43
2:M:99:LYS:HE2	2:M:106:LEU:HD23	2.00	0.43
1:B:192:ARG:HB2	1:B:211:SER:OG	2.17	0.43
2:N:99:LYS:HG2	2:N:103:SER:HB2	2.01	0.43
1:B:101:ASP:HB2	1:B:132:THR:HG23	2.01	0.43
1:B:79:LEU:CD1	1:B:81:LEU:HD23	2.49	0.43
1:D:248:PHE:O	1:D:399:ARG:CZ	2.67	0.43
2:L:99:LYS:HE2	2:L:106:LEU:HD23	2.01	0.43
2:O:99:LYS:HE2	2:O:106:LEU:HD23	2.01	0.43
1:B:85:ASN:HB2	1:B:114:ARG:HG3	2.00	0.43
1:A:99:LEU:HD11	1:B:62:TYR:OH	2.15	0.43
1:A:262:GLY:HA3	1:A:304:LEU:HD13	2.01	0.43
1:B:199:VAL:HB	2:L:111:ASN:HD21	1.84	0.43
1:C:107:HIS:NE2	1:C:131:THR:HG23	2.33	0.43
1:E:66:TYR:CE2	1:E:125:LEU:HD13	2.54	0.43
1:A:192:ARG:HB2	1:A:211:SER:OG	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:144:PRO:HD3	1:D:289:MET:HB2	2.00	0.42
1:B:248:PHE:HB3	1:B:399:ARG:HG2	1.99	0.42
1:E:238:ILE:HG13	1:E:239:MET:N	2.33	0.42
1:A:101:ASP:HB2	1:A:132:THR:HG23	2.01	0.42
1:C:66:TYR:CE2	1:C:125:LEU:HD13	2.55	0.42
1:D:83:LEU:HD11	1:D:118:LEU:HD12	2.01	0.42
1:B:145:LEU:HD23	1:B:226:ILE:HG12	2.02	0.42
1:C:31:PHE:CE2	1:D:81:LEU:HD22	2.54	0.42
1:D:262:GLY:HA3	1:D:304:LEU:HD13	2.02	0.42
1:D:58:THR:O	1:D:58:THR:HG23	2.19	0.42
1:B:402:PHE:HB2	1:B:403:PRO:HD3	2.01	0.42
1:E:43:ASP:HA	1:E:176:THR:OG1	2.20	0.42
1:E:248:PHE:O	1:E:399:ARG:CZ	2.67	0.41
1:B:280:LEU:HB3	1:B:281:PRO:HD2	2.02	0.41
1:C:88:ALA:HB2	1:C:116:ILE:HD11	2.02	0.41
1:C:255:VAL:N	1:C:256:PRO:CD	2.83	0.41
1:D:145:LEU:HD22	1:D:226:ILE:HD11	2.02	0.41
1:A:66:TYR:CE2	1:A:125:LEU:HD13	2.54	0.41
1:C:192:ARG:HB2	1:C:211:SER:OG	2.20	0.41
1:A:255:VAL:N	1:A:256:PRO:CD	2.83	0.41
1:C:254:SER:HB3	1:C:257:ALA:HB3	2.02	0.41
1:E:144:PRO:HD3	1:E:289:MET:HB2	2.03	0.41
1:B:85:ASN:HB2	1:B:114:ARG:CG	2.49	0.41
1:E:101:ASP:HB2	1:E:132:THR:HG23	2.02	0.41
1:E:189:VAL:CG1	1:E:213:ARG:HB3	2.50	0.41
1:C:36:VAL:HB	1:C:164:ILE:HD13	2.02	0.41
1:D:405:LEU:HA	1:D:405:LEU:HD23	1.90	0.41
1:A:267:LEU:HD22	1:B:268:THR:HG21	2.01	0.41
1:C:145:LEU:HD23	1:C:226:ILE:HG12	2.01	0.41
1:D:26:ARG:NH2	1:E:17:ASP:OD1	2.48	0.41
2:K:108:TYR:HA	2:K:109:PRO:HD3	1.86	0.41
1:C:85:ASN:HB2	1:C:114:ARG:CG	2.50	0.41
1:E:58:THR:HG23	1:E:58:THR:O	2.20	0.41
1:B:58:THR:HG23	1:B:58:THR:O	2.19	0.41
1:A:43:ASP:HA	1:A:176:THR:OG1	2.21	0.41
1:B:255:VAL:N	1:B:256:PRO:CD	2.84	0.41
1:D:402:PHE:HB2	1:D:403:PRO:HD3	2.02	0.41
1:D:145:LEU:HD12	1:D:417:TYR:HB3	2.03	0.41
1:A:268:THR:HG21	1:E:267:LEU:HD22	2.03	0.41
1:E:395:ASP:O	1:E:399:ARG:HG3	2.21	0.41
2:L:32:PRO:HB2	2:L:53:TRP:HB2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:262:GLY:HA3	1:B:304:LEU:HD13	2.03	0.41
1:D:254:SER:HB3	1:D:257:ALA:HB3	2.02	0.41
1:D:409:PHE:O	1:D:412:VAL:HG22	2.20	0.41
1:E:235:LEU:HD23	1:E:235:LEU:HA	1.91	0.41
2:N:99:LYS:HE2	2:N:106:LEU:HD23	2.03	0.41
1:B:145:LEU:HD13	1:B:417:TYR:HB3	2.03	0.40
1:B:99:LEU:HD11	1:C:62:TYR:OH	2.15	0.40
1:D:295:VAL:HG11	1:D:409:PHE:CZ	2.57	0.40
1:D:43:ASP:HA	1:D:176:THR:OG1	2.21	0.40
1:A:105:PHE:CE2	1:E:103:LYS:HE3	2.56	0.40
1:C:238:ILE:HG13	1:C:239:MET:N	2.36	0.40
1:C:43:ASP:HA	1:C:176:THR:OG1	2.21	0.40
2:M:103:SER:O	2:M:111:ASN:HB3	2.22	0.40
1:C:213:ARG:HB2	3:C:502:NAG:H81	2.02	0.40
1:E:409:PHE:O	1:E:412:VAL:HG22	2.21	0.40
2:K:32:PRO:HB2	2:K:53:TRP:HB2	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	328/395 (83%)	319 (97%)	9 (3%)	0	100	100
1	B	331/395 (84%)	322 (97%)	9 (3%)	0	100	100
1	C	328/395 (83%)	320 (98%)	8 (2%)	0	100	100
1	D	330/395 (84%)	320 (97%)	10 (3%)	0	100	100
1	E	328/395 (83%)	318 (97%)	10 (3%)	0	100	100
2	K	121/125 (97%)	115 (95%)	6 (5%)	0	100	100
2	L	121/125 (97%)	115 (95%)	6 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	M	121/125 (97%)	116 (96%)	5 (4%)	0	100	100
2	N	121/125 (97%)	116 (96%)	5 (4%)	0	100	100
2	O	121/125 (97%)	116 (96%)	5 (4%)	0	100	100
All	All	2250/2600 (86%)	2177 (97%)	73 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	293/346 (85%)	291 (99%)	2 (1%)	87	92
1	B	296/346 (86%)	295 (100%)	1 (0%)	94	97
1	C	293/346 (85%)	291 (99%)	2 (1%)	87	92
1	D	295/346 (85%)	293 (99%)	2 (1%)	87	92
1	E	293/346 (85%)	291 (99%)	2 (1%)	87	92
2	K	97/99 (98%)	97 (100%)	0	100	100
2	L	97/99 (98%)	97 (100%)	0	100	100
2	M	97/99 (98%)	97 (100%)	0	100	100
2	N	97/99 (98%)	97 (100%)	0	100	100
2	O	97/99 (98%)	97 (100%)	0	100	100
All	All	1955/2225 (88%)	1946 (100%)	9 (0%)	91	94

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	CYS
1	A	254	SER
1	B	37	CYS
1	C	37	CYS
1	C	254	SER

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Mol	Chain	Res	Type
1	D	37	CYS
1	D	254	SER
1	E	37	CYS
1	E	254	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	85	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

32 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	501	1	14,14,15	0.34	0	15,19,21	0.42	0
3	NAG	A	502	1,3	14,14,15	0.25	0	15,19,21	0.53	0
3	NAG	A	503	3,4	14,14,15	0.21	0	15,19,21	0.52	0
4	BMA	A	504	3,5	11,11,12	0.19	0	13,15,17	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	MAN	A	505	4	11,11,12	0.24	0	13,15,17	0.39	0
5	MAN	A	506	4	11,11,12	0.28	0	13,15,17	0.26	0
3	NAG	B	501	1	14,14,15	0.25	0	15,19,21	0.54	0
3	NAG	B	502	1,3	14,14,15	0.28	0	15,19,21	0.66	0
3	NAG	B	503	3,4	14,14,15	0.19	0	15,19,21	0.55	0
4	BMA	B	504	3,5	11,11,12	0.39	0	13,15,17	0.77	0
5	MAN	B	505	4	11,11,12	0.36	0	13,15,17	0.28	0
5	MAN	B	506	4	11,11,12	0.28	0	13,15,17	0.31	0
3	NAG	C	501	1	14,14,15	0.25	0	15,19,21	0.52	0
3	NAG	C	502	1,3	14,14,15	0.21	0	15,19,21	0.60	0
3	NAG	C	503	3,4	14,14,15	0.22	0	15,19,21	0.51	0
4	BMA	C	504	3,5	11,11,12	0.37	0	13,15,17	0.57	0
5	MAN	C	505	4	11,11,12	0.24	0	13,15,17	0.59	0
5	MAN	C	506	4	11,11,12	0.17	0	13,15,17	0.22	0
3	NAG	D	501	1	14,14,15	0.34	0	15,19,21	0.53	0
3	NAG	D	502	1,3	14,14,15	0.25	0	15,19,21	0.59	0
3	NAG	D	503	3,4	14,14,15	0.29	0	15,19,21	0.46	0
4	BMA	D	504	3,5	11,11,12	0.26	0	13,15,17	0.33	0
5	MAN	D	505	5,4	11,11,12	0.43	0	13,15,17	0.33	0
5	MAN	D	506	5	11,11,12	0.51	0	13,15,17	0.56	0
5	MAN	D	507	4	11,11,12	0.24	0	13,15,17	0.22	0
3	NAG	E	501	1	14,14,15	0.21	0	15,19,21	0.69	0
3	NAG	E	502	1,3	14,14,15	0.23	0	15,19,21	0.60	0
3	NAG	E	503	3,4	14,14,15	0.23	0	15,19,21	0.44	0
4	BMA	E	504	3,5	11,11,12	0.33	0	13,15,17	0.35	0
5	MAN	E	505	5,4	11,11,12	0.32	0	13,15,17	0.50	0
5	MAN	E	506	5	11,11,12	0.30	0	13,15,17	0.31	0
5	MAN	E	507	4	11,11,12	0.45	0	13,15,17	0.46	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	501	1	-	0/6/23/26	0/1/1/1
3	NAG	A	502	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	503	3,4	-	0/6/23/26	0/1/1/1
4	BMA	A	504	3,5	-	0/2/19/22	0/1/1/1
5	MAN	A	505	4	-	0/2/19/22	0/1/1/1
5	MAN	A	506	4	-	0/2/19/22	0/1/1/1
3	NAG	B	501	1	-	0/6/23/26	0/1/1/1
3	NAG	B	502	1,3	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	B	503	3,4	-	0/6/23/26	0/1/1/1
4	BMA	B	504	3,5	-	0/2/19/22	0/1/1/1
5	MAN	B	505	4	-	0/2/19/22	0/1/1/1
5	MAN	B	506	4	-	0/2/19/22	0/1/1/1
3	NAG	C	501	1	-	0/6/23/26	0/1/1/1
3	NAG	C	502	1,3	-	0/6/23/26	0/1/1/1
3	NAG	C	503	3,4	-	0/6/23/26	0/1/1/1
4	BMA	C	504	3,5	-	0/2/19/22	0/1/1/1
5	MAN	C	505	4	-	0/2/19/22	0/1/1/1
5	MAN	C	506	4	-	0/2/19/22	0/1/1/1
3	NAG	D	501	1	-	0/6/23/26	0/1/1/1
3	NAG	D	502	1,3	-	0/6/23/26	0/1/1/1
3	NAG	D	503	3,4	-	0/6/23/26	0/1/1/1
4	BMA	D	504	3,5	-	0/2/19/22	0/1/1/1
5	MAN	D	505	5,4	-	0/2/19/22	0/1/1/1
5	MAN	D	506	5	-	0/2/19/22	0/1/1/1
5	MAN	D	507	4	-	0/2/19/22	0/1/1/1
3	NAG	E	501	1	-	0/6/23/26	0/1/1/1
3	NAG	E	502	1,3	-	0/6/23/26	0/1/1/1
3	NAG	E	503	3,4	-	0/6/23/26	0/1/1/1
4	BMA	E	504	3,5	-	0/2/19/22	0/1/1/1
5	MAN	E	505	5,4	-	0/2/19/22	0/1/1/1
5	MAN	E	506	5	-	0/2/19/22	0/1/1/1
5	MAN	E	507	4	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	502	NAG	1	0
3	E	502	NAG	1	0
4	E	504	BMA	1	0
5	E	505	MAN	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	330/395 (83%)	-0.20	6 (1%) 69 66	30, 58, 101, 133	0
1	B	333/395 (84%)	0.04	10 (3%) 51 49	31, 62, 94, 138	0
1	C	330/395 (83%)	-0.02	7 (2%) 64 61	31, 61, 104, 137	0
1	D	332/395 (84%)	-0.14	5 (1%) 74 70	28, 51, 85, 155	0
1	E	330/395 (83%)	0.04	12 (3%) 43 40	26, 54, 89, 141	0
2	K	123/125 (98%)	0.96	22 (17%) 2 1	54, 101, 132, 151	0
2	L	123/125 (98%)	0.81	22 (17%) 2 1	78, 125, 168, 182	0
2	M	123/125 (98%)	0.45	15 (12%) 5 4	58, 97, 129, 150	0
2	N	123/125 (98%)	0.39	14 (11%) 6 5	52, 102, 141, 153	0
2	O	123/125 (98%)	-0.21	5 (4%) 38 35	60, 91, 117, 123	0
All	All	2270/2600 (87%)	0.09	118 (5%) 28 26	26, 65, 129, 182	0

All (118) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	68	PHE	6.1
2	L	115	TRP	6.1
1	A	127	GLY	5.7
2	L	116	GLY	5.6
2	L	117	GLN	5.4
2	K	88	PRO	5.2
2	N	43	LYS	5.0
2	K	1	GLN	4.8
1	B	417	TYR	4.8
1	E	204	ALA	4.5
2	K	115	TRP	4.5
2	N	44	GLU	4.2
2	M	1	GLN	4.1

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Mol	Chain	Res	Type	RSRZ
2	L	85	ASN	4.1
2	K	114	TYR	4.1
2	M	86	LEU	4.0
1	D	417	TYR	3.9
2	L	1	GLN	3.9
2	K	96	CYS	3.9
2	L	4	LEU	3.8
2	L	98	ALA	3.8
1	B	278	ASN	3.7
2	M	123	VAL	3.6
2	L	114	TYR	3.6
1	C	278	ASN	3.5
2	N	88	PRO	3.5
2	L	96	CYS	3.5
2	N	68	PHE	3.5
2	K	123	VAL	3.4
1	C	232	GLN	3.4
1	C	215	LYS	3.4
2	K	75	ALA	3.3
2	N	41	PRO	3.3
2	K	14	ALA	3.3
2	N	10	GLY	3.3
2	M	11	LEU	3.3
2	K	74	ASN	3.1
1	B	113	ASN	3.1
2	L	10	GLY	3.0
2	M	13	GLN	3.0
1	E	14	GLU	3.0
2	M	15	GLY	3.0
1	A	66	TYR	3.0
2	L	70	ILE	3.0
2	M	85	ASN	3.0
1	E	13	LYS	2.9
1	B	159	TYR	2.8
1	E	417	TYR	2.8
2	M	115	TRP	2.7
2	K	98	ALA	2.7
2	K	43	LYS	2.7
2	L	11	LEU	2.7
2	K	44	GLU	2.7
2	K	89	GLU	2.6
2	M	66	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
2	K	116	GLY	2.6
2	L	9	GLY	2.6
1	B	77	ILE	2.6
2	K	53	TRP	2.6
2	L	67	ARG	2.6
2	L	37	PHE	2.6
2	O	14	ALA	2.6
2	N	40	ALA	2.6
1	C	417	TYR	2.6
2	N	42	GLY	2.6
2	M	14	ALA	2.5
2	K	113	ASP	2.5
2	L	95	TYR	2.5
2	M	67	ARG	2.5
2	M	16	GLY	2.5
1	A	126	TYR	2.5
1	A	125	LEU	2.5
2	K	13	GLN	2.5
2	N	84	ASN	2.5
1	D	9	MET	2.5
1	D	416	THR	2.5
2	N	86	LEU	2.5
1	B	145	LEU	2.4
1	E	12	VAL	2.4
1	E	158	GLY	2.4
2	O	75	ALA	2.4
2	L	3	GLN	2.4
2	L	97	ALA	2.4
1	D	142	ARG	2.4
2	L	64	VAL	2.4
1	A	115	MET	2.4
1	C	74	TYR	2.3
1	B	14	GLU	2.3
1	D	24	ASP	2.3
2	K	86	LEU	2.3
2	L	66	ASP	2.3
2	N	90	ASP	2.2
1	A	159	TYR	2.2
2	K	94	TYR	2.2
2	N	94	TYR	2.2
2	L	5	GLN	2.2
2	M	84	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	313	PHE	2.2
2	M	17	SER	2.1
2	O	85	ASN	2.1
2	K	37	PHE	2.1
1	B	9	MET	2.1
2	M	87	LYS	2.1
1	E	147	GLU	2.1
1	C	145	LEU	2.1
1	E	319	ALA	2.1
2	K	87	LYS	2.1
1	B	114	ARG	2.1
2	O	86	LEU	2.1
1	C	418	LEU	2.1
2	N	9	GLY	2.1
2	K	117	GLN	2.0
1	E	161	THR	2.0
1	E	393	LYS	2.0
2	O	44	GLU	2.0
2	N	85	ASN	2.0
1	E	16	VAL	2.0
1	E	205	TYR	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 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. 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	D	501	14/15	0.92	0.20	0.88	93,102,115,117	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	NAG	E	503	14/15	0.95	0.15	-0.37	56,70,77,80	0
3	NAG	C	502	14/15	0.96	0.16	-0.71	68,75,81,85	0
3	NAG	B	502	14/15	0.93	0.16	-0.80	71,77,81,84	0
3	NAG	B	501	14/15	0.94	0.17	-0.88	96,100,112,119	0
3	NAG	E	502	14/15	0.94	0.13	-1.04	66,68,74,77	0
3	NAG	D	502	14/15	0.96	0.16	-1.09	46,49,53,55	0
3	NAG	D	503	14/15	0.92	0.15	-1.21	53,64,74,87	0
3	NAG	A	503	14/15	0.93	0.18	-1.31	65,69,73,79	0
3	NAG	A	502	14/15	0.95	0.14	-1.40	54,59,63,71	0
3	NAG	B	503	14/15	0.95	0.14	-1.65	84,88,98,101	0
3	NAG	C	503	14/15	0.95	0.10	-3.25	65,78,87,99	0
5	MAN	C	505	11/12	0.62	0.26	-	113,123,130,133	0
5	MAN	A	505	11/12	0.72	0.23	-	102,109,116,117	0
5	MAN	E	506	11/12	0.90	0.17	-	93,115,122,123	0
4	BMA	A	504	11/12	0.83	0.14	-	86,96,110,122	0
3	NAG	A	501	14/15	0.86	0.27	-	122,141,148,149	0
3	NAG	C	501	14/15	0.92	0.14	-	140,153,164,166	0
5	MAN	C	506	11/12	0.83	0.40	-	105,120,130,132	0
3	NAG	E	501	14/15	0.92	0.15	-	101,108,114,117	0
5	MAN	E	505	11/12	0.77	0.35	-	112,120,129,139	0
4	BMA	E	504	11/12	0.90	0.16	-	78,83,90,105	0
5	MAN	B	506	11/12	0.87	0.15	-	98,108,114,118	0
5	MAN	D	506	11/12	0.77	0.19	-	121,129,137,141	0
5	MAN	A	506	11/12	0.93	0.12	-	93,99,112,112	0
4	BMA	C	504	11/12	0.81	0.12	-	104,113,127,138	0
5	MAN	B	505	11/12	0.61	0.31	-	98,114,118,120	0
5	MAN	E	507	11/12	0.91	0.16	-	80,86,92,94	0
5	MAN	D	505	11/12	0.76	0.24	-	104,123,129,138	0
5	MAN	D	507	11/12	0.90	0.20	-	108,123,134,135	0
4	BMA	D	504	11/12	0.91	0.11	-	93,99,114,118	0
4	BMA	B	504	11/12	0.85	0.13	-	101,108,113,121	0

6.5 Other polymers [i](#)

There are no such residues in this entry.