



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:34 PM GMT

PDB ID : 4COF  
Title : Crystal structure of a human gamma-aminobutyric acid receptor, the GABA(A)R-beta3 homopentamer  
Authors : Miller, P.S.; Aricescu, A.R.  
Deposited on : 2014-01-28  
Resolution : 2.97 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

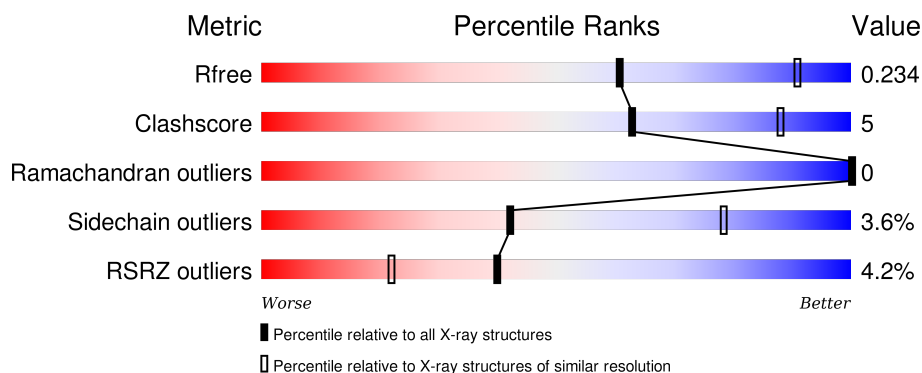
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.97 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	1992 (3.00-2.96)
Clashscore	102246	2349 (3.00-2.96)
Ramachandran outliers	100387	2274 (3.00-2.96)
Sidechain outliers	100360	2277 (3.00-2.96)
RSRZ outliers	91569	2007 (3.00-2.96)

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  $\geq 3$ , 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	355	<div> <div>2%</div> <div>82%</div> <div>13%</div> <div>• •</div> </div>
1	B	355	<div> <div>4%</div> <div>79%</div> <div>13%</div> <div>• 7%</div> </div>
1	C	355	<div> <div>5%</div> <div>79%</div> <div>13%</div> <div>• 7%</div> </div>
1	D	355	<div> <div>4%</div> <div>75%</div> <div>17%</div> <div>• 6%</div> </div>
1	E	355	<div> <div>5%</div> <div>77%</div> <div>15%</div> <div>• 7%</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	BEN	A	500	-	-	-	X
2	BEN	B	500	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13972 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 GAMMA-AMINOBUTYRIC ACID RECEPTOR SUBUNIT BETA-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	340	Total	C	N	O	S	0	0	0
			2779	1816	455	492	16			
1	B	331	Total	C	N	O	S	0	0	0
			2714	1779	443	477	15			
1	C	331	Total	C	N	O	S	0	0	0
			2714	1779	443	477	15			
1	D	332	Total	C	N	O	S	0	0	0
			2722	1783	445	479	15			
1	E	331	Total	C	N	O	S	0	0	0
			2714	1779	443	477	15			

There are 105 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLU	-	EXPRESSION TAG	UNP P28472
A	-1	THR	-	EXPRESSION TAG	UNP P28472
A	0	GLY	-	EXPRESSION TAG	UNP P28472
A	449	GLY	-	EXPRESSION TAG	UNP P28472
A	450	ALA	-	EXPRESSION TAG	UNP P28472
A	451	THR	-	EXPRESSION TAG	UNP P28472
A	452	GLU	-	EXPRESSION TAG	UNP P28472
A	453	THR	-	EXPRESSION TAG	UNP P28472
A	454	SER	-	EXPRESSION TAG	UNP P28472
A	455	GLN	-	EXPRESSION TAG	UNP P28472
A	456	VAL	-	EXPRESSION TAG	UNP P28472
A	457	ALA	-	EXPRESSION TAG	UNP P28472
A	458	PRO	-	EXPRESSION TAG	UNP P28472
A	459	ALA	-	EXPRESSION TAG	UNP P28472
A	308	SER	-	LINKER	UNP P28472
A	309	GLN	-	LINKER	UNP P28472
A	310	PRO	-	LINKER	UNP P28472
A	311	ALA	-	LINKER	UNP P28472

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Chain	Residue	Modelled	Actual	Comment	Reference
A	312	ARG	-	LINKER	UNP P28472
A	313	ALA	-	LINKER	UNP P28472
A	314	ALA	-	LINKER	UNP P28472
B	-2	GLU	-	EXPRESSION TAG	UNP P28472
B	-1	THR	-	EXPRESSION TAG	UNP P28472
B	0	GLY	-	EXPRESSION TAG	UNP P28472
B	449	GLY	-	EXPRESSION TAG	UNP P28472
B	450	ALA	-	EXPRESSION TAG	UNP P28472
B	451	THR	-	EXPRESSION TAG	UNP P28472
B	452	GLU	-	EXPRESSION TAG	UNP P28472
B	453	THR	-	EXPRESSION TAG	UNP P28472
B	454	SER	-	EXPRESSION TAG	UNP P28472
B	455	GLN	-	EXPRESSION TAG	UNP P28472
B	456	VAL	-	EXPRESSION TAG	UNP P28472
B	457	ALA	-	EXPRESSION TAG	UNP P28472
B	458	PRO	-	EXPRESSION TAG	UNP P28472
B	459	ALA	-	EXPRESSION TAG	UNP P28472
B	308	SER	-	LINKER	UNP P28472
B	309	GLN	-	LINKER	UNP P28472
B	310	PRO	-	LINKER	UNP P28472
B	311	ALA	-	LINKER	UNP P28472
B	312	ARG	-	LINKER	UNP P28472
B	313	ALA	-	LINKER	UNP P28472
B	314	ALA	-	LINKER	UNP P28472
C	-2	GLU	-	EXPRESSION TAG	UNP P28472
C	-1	THR	-	EXPRESSION TAG	UNP P28472
C	0	GLY	-	EXPRESSION TAG	UNP P28472
C	449	GLY	-	EXPRESSION TAG	UNP P28472
C	450	ALA	-	EXPRESSION TAG	UNP P28472
C	451	THR	-	EXPRESSION TAG	UNP P28472
C	452	GLU	-	EXPRESSION TAG	UNP P28472
C	453	THR	-	EXPRESSION TAG	UNP P28472
C	454	SER	-	EXPRESSION TAG	UNP P28472
C	455	GLN	-	EXPRESSION TAG	UNP P28472
C	456	VAL	-	EXPRESSION TAG	UNP P28472
C	457	ALA	-	EXPRESSION TAG	UNP P28472
C	458	PRO	-	EXPRESSION TAG	UNP P28472
C	459	ALA	-	EXPRESSION TAG	UNP P28472
C	308	SER	-	LINKER	UNP P28472
C	309	GLN	-	LINKER	UNP P28472
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C	311	ALA	-	LINKER	UNP P28472

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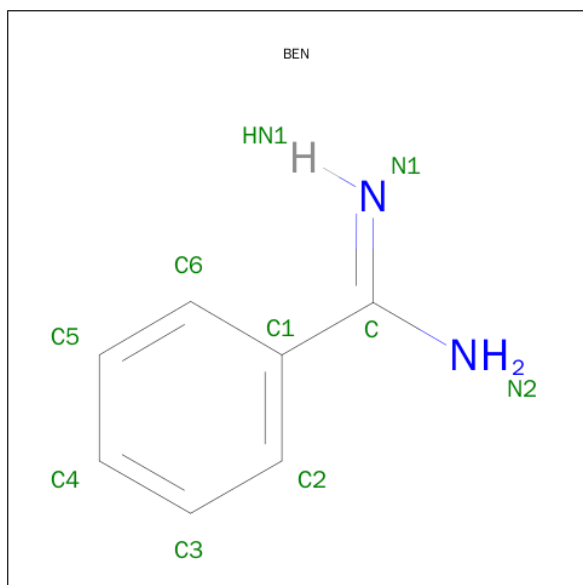
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C	312	ARG	-	LINKER	UNP P28472
C	313	ALA	-	LINKER	UNP P28472
C	314	ALA	-	LINKER	UNP P28472
D	-2	GLU	-	EXPRESSION TAG	UNP P28472
D	-1	THR	-	EXPRESSION TAG	UNP P28472
D	0	GLY	-	EXPRESSION TAG	UNP P28472
D	449	GLY	-	EXPRESSION TAG	UNP P28472
D	450	ALA	-	EXPRESSION TAG	UNP P28472
D	451	THR	-	EXPRESSION TAG	UNP P28472
D	452	GLU	-	EXPRESSION TAG	UNP P28472
D	453	THR	-	EXPRESSION TAG	UNP P28472
D	454	SER	-	EXPRESSION TAG	UNP P28472
D	455	GLN	-	EXPRESSION TAG	UNP P28472
D	456	VAL	-	EXPRESSION TAG	UNP P28472
D	457	ALA	-	EXPRESSION TAG	UNP P28472
D	458	PRO	-	EXPRESSION TAG	UNP P28472
D	459	ALA	-	EXPRESSION TAG	UNP P28472
D	308	SER	-	LINKER	UNP P28472
D	309	GLN	-	LINKER	UNP P28472
D	310	PRO	-	LINKER	UNP P28472
D	311	ALA	-	LINKER	UNP P28472
D	312	ARG	-	LINKER	UNP P28472
D	313	ALA	-	LINKER	UNP P28472
D	314	ALA	-	LINKER	UNP P28472
E	-2	GLU	-	EXPRESSION TAG	UNP P28472
E	-1	THR	-	EXPRESSION TAG	UNP P28472
E	0	GLY	-	EXPRESSION TAG	UNP P28472
E	449	GLY	-	EXPRESSION TAG	UNP P28472
E	450	ALA	-	EXPRESSION TAG	UNP P28472
E	451	THR	-	EXPRESSION TAG	UNP P28472
E	452	GLU	-	EXPRESSION TAG	UNP P28472
E	453	THR	-	EXPRESSION TAG	UNP P28472
E	454	SER	-	EXPRESSION TAG	UNP P28472
E	455	GLN	-	EXPRESSION TAG	UNP P28472
E	456	VAL	-	EXPRESSION TAG	UNP P28472
E	457	ALA	-	EXPRESSION TAG	UNP P28472
E	458	PRO	-	EXPRESSION TAG	UNP P28472
E	459	ALA	-	EXPRESSION TAG	UNP P28472
E	308	SER	-	LINKER	UNP P28472
E	309	GLN	-	LINKER	UNP P28472
E	310	PRO	-	LINKER	UNP P28472
E	311	ALA	-	LINKER	UNP P28472

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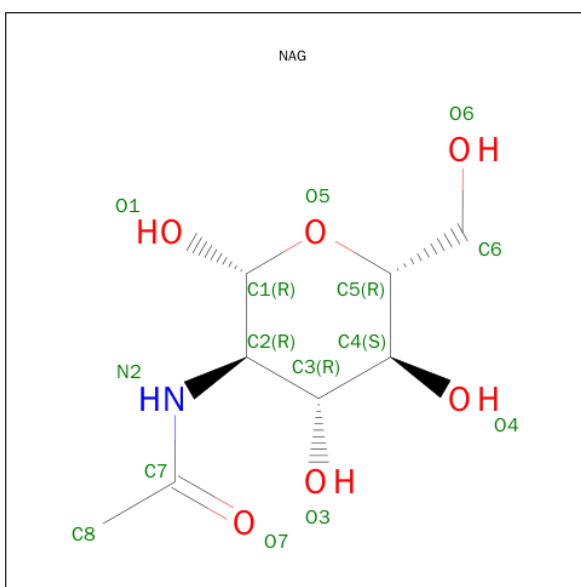
Chain	Residue	Modelled	Actual	Comment	Reference
E	312	ARG	-	LINKER	UNP P28472
E	313	ALA	-	LINKER	UNP P28472
E	314	ALA	-	LINKER	UNP P28472

- Molecule 2 is BENZAMIDINE (three-letter code: BEN) (formula:  $C_7H_8N_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			9	7	2		
2	B	1	Total	C	N	0	0
			9	7	2		
2	C	1	Total	C	N	0	0
			9	7	2		
2	D	1	Total	C	N	0	0
			9	7	2		
2	E	1	Total	C	N	0	0
			9	7	2		

- 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	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	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	E	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Cl	0	0
			1	1		
4	A	1	Total	Cl	0	0
			1	1		
4	D	1	Total	Cl	0	0
			1	1		
4	C	1	Total	Cl	0	0
			1	1		
4	E	1	Total	Cl	0	0
			1	1		

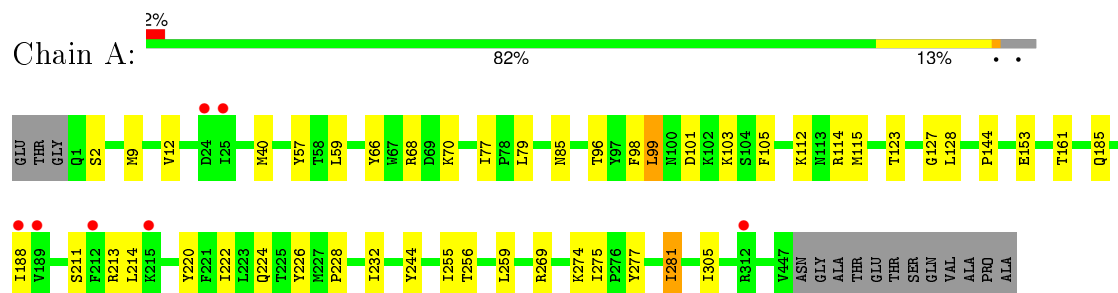
- Molecule 5 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	3	Total	C	N	O	0	0
			39	22	2	15		
5	B	3	Total	C	N	O	0	0
			39	22	2	15		
5	C	3	Total	C	N	O	0	0
			39	22	2	15		
5	D	3	Total	C	N	O	0	0
			39	22	2	15		
5	E	3	Total	C	N	O	0	0
			39	22	2	15		

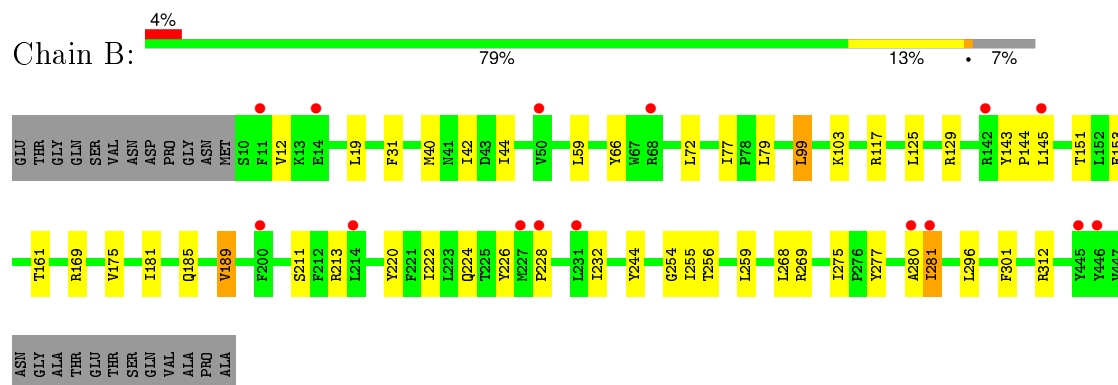
### 3 Residue-property plots

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

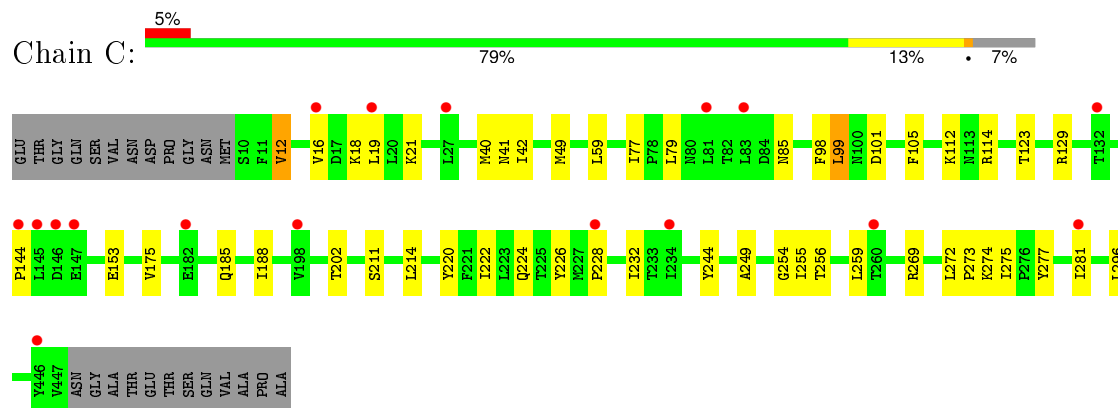
#### • Molecule 1: GAMMA-AMINOBUTYRIC ACID RECEPTOR SUBUNIT BETA-3



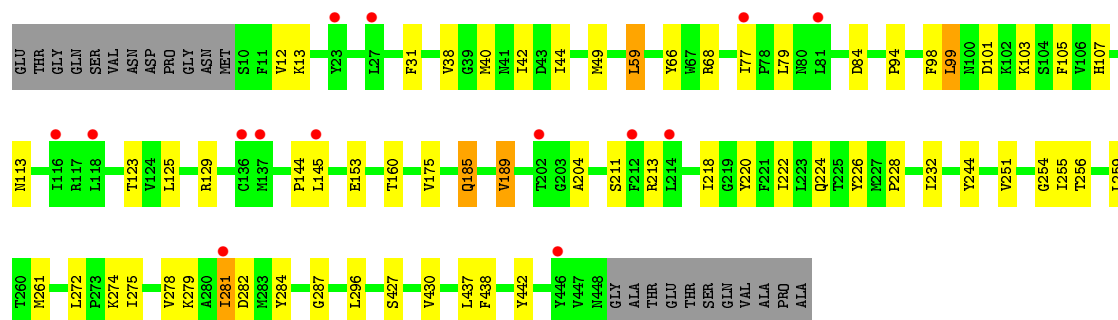
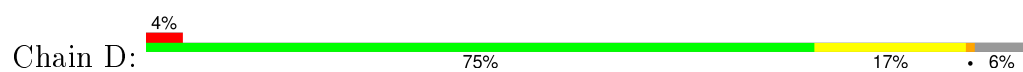
#### • Molecule 1: GAMMA-AMINOBUTYRIC ACID RECEPTOR SUBUNIT BETA-3



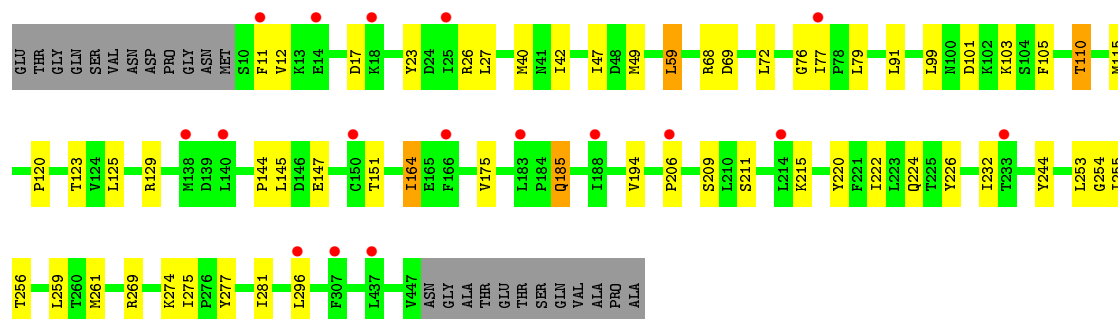
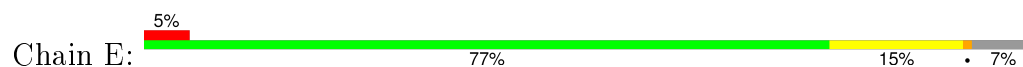
#### • Molecule 1: GAMMA-AMINOBUTYRIC ACID RECEPTOR SUBUNIT BETA-3



#### • Molecule 1: GAMMA-AMINOBUTYRIC ACID RECEPTOR SUBUNIT BETA-3



• Molecule 1: GAMMA-AMINOBUTYRIC ACID RECEPTOR SUBUNIT BETA-3



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	174.10Å 108.90Å 207.44Å 90.00° 107.43° 90.00°	Depositor
Resolution (Å)	40.00 – 2.97 98.96 – 2.97	Depositor EDS
% Data completeness (in resolution range)	99.8 (40.00-2.97) 99.8 (98.96-2.97)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.67 (at 2.96Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, $R_{free}$	0.205 , 0.226 0.212 , 0.234	Depositor DCC
$R_{free}$ test set	3835 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	81.8	Xtriage
Anisotropy	0.684	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 67.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	3 of 76360 reflections (0.004%)	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	13972	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	102.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

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

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, BEN, NAG, CL

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	0.51	0/2854	0.68	0/3886
1	B	0.52	0/2788	0.68	0/3796
1	C	0.51	0/2788	0.68	0/3796
1	D	0.52	0/2796	0.69	0/3807
1	E	0.51	0/2788	0.68	0/3796
All	All	0.51	0/14014	0.68	0/19081

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 [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	2779	0	2769	35	0
1	B	2714	0	2710	36	0
1	C	2714	0	2710	32	0
1	D	2722	0	2716	48	0
1	E	2714	0	2710	43	0
2	A	9	0	7	0	0
2	B	9	0	7	0	0
2	C	9	0	7	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	9	0	7	0	0
2	E	9	0	7	0	0
3	A	28	0	26	0	0
3	B	14	0	13	0	0
3	C	14	0	13	0	0
3	D	14	0	13	0	0
3	E	14	0	13	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
5	A	39	0	34	0	0
5	B	39	0	34	0	0
5	C	39	0	34	0	0
5	D	39	0	34	0	0
5	E	39	0	34	0	0
All	All	13972	0	13898	148	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 5.

All (148) 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:E:68:ARG:HG2	1:E:123:THR:HG22	1.41	1.01
1:D:144:PRO:HD3	1:D:281:ILE:HG12	1.52	0.90
1:A:85:ASN:HB2	1:A:114:ARG:HG3	1.61	0.83
1:D:66:TYR:CE2	1:D:125:LEU:HD13	2.18	0.78
1:A:256:THR:HG23	1:B:259:LEU:HD11	1.65	0.78
1:B:66:TYR:CE2	1:B:125:LEU:HD13	2.19	0.78
1:D:256:THR:HG23	1:E:259:LEU:HD11	1.67	0.76
1:A:259:LEU:HD11	1:E:256:THR:HG23	1.66	0.76
1:D:44:ILE:HD12	1:D:59:LEU:HD11	1.70	0.74
1:D:129:ARG:HH22	1:E:101:ASP:HB3	1.53	0.73
1:C:256:THR:HG23	1:D:259:LEU:HD11	1.70	0.71
1:B:228:PRO:O	1:B:232:ILE:HG12	1.91	0.70
1:B:256:THR:HG23	1:C:259:LEU:HD11	1.72	0.70
1:D:66:TYR:CZ	1:D:125:LEU:HD13	2.28	0.68
1:A:224:GLN:HG2	1:B:269:ARG:HD2	1.76	0.65
1:D:256:THR:HG21	1:E:255:ILE:HG23	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:272:LEU:HD13	1:D:279:LYS:HE3	1.77	0.65
1:E:151:THR:HG22	1:E:211:SER:HB3	1.80	0.64
1:C:256:THR:HG21	1:D:255:ILE:HG23	1.82	0.62
1:C:254:GLY:HA3	1:C:296:LEU:HD13	1.82	0.60
1:A:259:LEU:HD11	1:E:256:THR:CG2	2.32	0.60
1:A:256:THR:HG21	1:B:255:ILE:HG23	1.84	0.59
1:D:144:PRO:HD2	1:D:145:LEU:HD12	1.84	0.58
1:D:224:GLN:HG2	1:E:269:ARG:HD2	1.85	0.58
1:A:269:ARG:HD2	1:E:224:GLN:HG2	1.84	0.58
1:E:144:PRO:HD3	1:E:281:ILE:HB	1.85	0.57
1:C:256:THR:HG21	1:D:255:ILE:CG2	2.34	0.57
1:A:57:TYR:HE2	1:A:59:LEU:HD23	1.69	0.57
1:D:275:ILE:HD11	1:D:279:LYS:HE2	1.87	0.56
1:B:66:TYR:CZ	1:B:125:LEU:HD13	2.41	0.56
1:A:96:THR:O	1:E:110:THR:HB	2.05	0.56
1:E:11:PHE:HZ	1:E:76:GLY:HA3	1.71	0.56
1:D:256:THR:HG21	1:E:255:ILE:CG2	2.36	0.56
1:B:256:THR:HG21	1:C:255:ILE:HG23	1.88	0.55
1:A:114:ARG:HB3	1:A:128:LEU:HD23	1.88	0.55
1:E:11:PHE:CZ	1:E:76:GLY:HA3	2.42	0.54
1:D:254:GLY:HA3	1:D:296:LEU:HD13	1.89	0.54
1:E:164:ILE:HG23	1:E:206:PRO:HG3	1.89	0.54
1:C:16:VAL:O	1:C:19:LEU:HB2	2.07	0.53
1:B:254:GLY:HA3	1:B:296:LEU:HD13	1.90	0.53
1:E:72:LEU:HD21	1:E:91:LEU:HD22	1.89	0.53
1:A:256:THR:CG2	1:B:259:LEU:HD11	2.34	0.53
1:D:256:THR:CG2	1:E:259:LEU:HD11	2.37	0.53
1:C:144:PRO:HD3	1:C:281:ILE:HB	1.91	0.52
1:E:254:GLY:HA3	1:E:296:LEU:HD13	1.91	0.52
1:A:144:PRO:HD3	1:A:281:ILE:HB	1.92	0.52
1:D:272:LEU:HD13	1:D:279:LYS:CE	2.41	0.51
1:E:220:TYR:CZ	1:E:224:GLN:HG3	2.45	0.51
1:A:220:TYR:CZ	1:A:224:GLN:HG3	2.45	0.51
1:A:256:THR:HG21	1:B:255:ILE:CG2	2.40	0.51
1:D:220:TYR:CZ	1:D:224:GLN:HG3	2.46	0.51
1:B:144:PRO:HD2	1:B:145:LEU:HD12	1.92	0.51
1:E:232:ILE:HB	1:E:261:MET:HE2	1.92	0.51
1:B:256:THR:HG21	1:C:255:ILE:CG2	2.41	0.50
1:B:220:TYR:CZ	1:B:224:GLN:HG3	2.46	0.50
1:B:129:ARG:HH22	1:C:101:ASP:HB3	1.76	0.50
1:A:66:TYR:CZ	1:A:123:THR:HG21	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:228:PRO:O	1:C:232:ILE:HG12	2.12	0.50
1:A:255:ILE:HG23	1:E:256:THR:HG21	1.92	0.50
1:B:224:GLN:HG2	1:C:269:ARG:HD2	1.94	0.50
1:D:228:PRO:O	1:D:232:ILE:HG12	2.12	0.49
1:A:255:ILE:CG2	1:E:256:THR:HG21	2.42	0.49
1:B:66:TYR:CD1	1:B:169:ARG:HD3	2.48	0.49
1:C:220:TYR:CZ	1:C:224:GLN:HG3	2.47	0.49
1:C:98:PHE:HB2	1:C:101:ASP:HB2	1.93	0.49
1:A:274:LYS:HD2	1:E:185:GLN:NE2	2.28	0.49
1:A:228:PRO:O	1:A:232:ILE:HG12	2.13	0.48
1:D:68:ARG:HA	1:D:123:THR:HA	1.94	0.48
1:C:256:THR:CG2	1:D:259:LEU:HD11	2.40	0.48
1:D:189:VAL:HG12	1:D:213:ARG:HB3	1.95	0.48
1:D:145:LEU:HG	1:D:218:ILE:HD13	1.95	0.48
1:A:68:ARG:HD2	1:A:70:LYS:NZ	2.29	0.48
1:B:117:ARG:NH2	1:C:202:THR:HB	2.29	0.48
1:A:68:ARG:HA	1:A:123:THR:HA	1.95	0.47
1:D:98:PHE:HB2	1:D:101:ASP:HB2	1.97	0.47
1:A:101:ASP:HB3	1:E:129:ARG:HH22	1.79	0.47
1:A:98:PHE:HB2	1:A:101:ASP:HB2	1.96	0.47
1:B:189:VAL:HG12	1:B:213:ARG:HB3	1.96	0.47
1:B:268:LEU:HD21	1:B:281:ILE:HG21	1.96	0.46
1:C:129:ARG:HH22	1:D:101:ASP:HB3	1.79	0.46
1:A:188:ILE:HD13	1:A:214:LEU:HD23	1.98	0.46
1:C:42:ILE:HB	1:C:175:VAL:HG22	1.97	0.46
1:C:99:LEU:HB3	1:C:153:GLU:HB2	1.97	0.46
1:E:144:PRO:HD2	1:E:145:LEU:HD12	1.98	0.46
1:B:275:ILE:HD12	1:B:277:TYR:CE1	2.51	0.46
1:D:77:ILE:HG22	1:D:79:LEU:H	1.81	0.46
1:D:185:GLN:NE2	1:E:274:LYS:HD2	2.30	0.46
1:C:18:LYS:HD3	1:C:21:LYS:HD2	1.97	0.46
1:D:42:ILE:HB	1:D:175:VAL:HG22	1.97	0.45
1:E:275:ILE:HD12	1:E:277:TYR:CE1	2.51	0.45
1:C:105:PHE:CZ	1:D:103:LYS:HD2	2.51	0.45
1:C:272:LEU:HB3	1:C:273:PRO:HD2	1.99	0.45
1:D:278:VAL:HG23	1:D:282:ASP:HB3	1.99	0.45
1:E:23:TYR:CE2	1:E:72:LEU:HD11	2.52	0.45
1:E:77:ILE:HG22	1:E:79:LEU:H	1.82	0.45
1:E:42:ILE:HB	1:E:175:VAL:HG22	1.98	0.44
1:E:115:MET:HE1	1:E:125:LEU:HG	1.99	0.44
1:A:103:LYS:HD2	1:E:105:PHE:CZ	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:222:ILE:HA	1:C:226:TYR:HB2	1.98	0.44
1:E:47:ILE:HG12	1:E:59:LEU:HD12	2.00	0.44
1:C:188:ILE:HD13	1:C:214:LEU:HD23	1.98	0.44
1:D:84:ASP:HB3	1:E:26:ARG:HA	1.98	0.44
1:E:222:ILE:HA	1:E:226:TYR:HB2	1.98	0.44
1:D:107:HIS:HB2	1:D:113:ASN:OD1	2.18	0.44
1:A:222:ILE:HA	1:A:226:TYR:HB2	2.00	0.44
1:D:13:LYS:HG2	1:E:27:LEU:HD22	2.00	0.44
1:A:105:PHE:CZ	1:B:103:LYS:HD2	2.53	0.44
1:A:99:LEU:HB3	1:A:153:GLU:HB2	2.00	0.44
1:C:275:ILE:HD12	1:C:277:TYR:CE1	2.52	0.44
1:E:68:ARG:HA	1:E:123:THR:HA	1.99	0.44
1:B:77:ILE:HG22	1:B:79:LEU:H	1.82	0.44
1:D:105:PHE:CZ	1:E:103:LYS:HD2	2.53	0.44
1:B:99:LEU:HB3	1:B:153:GLU:HB2	2.00	0.44
1:D:222:ILE:HA	1:D:226:TYR:HB2	2.00	0.43
1:B:256:THR:CG2	1:C:259:LEU:HD11	2.43	0.43
1:A:9:MET:HG2	1:B:31:PHE:HB3	1.99	0.43
1:A:68:ARG:HH11	1:A:70:LYS:HE2	1.83	0.43
1:A:275:ILE:HD12	1:A:277:TYR:CE1	2.54	0.43
1:A:114:ARG:HA	1:A:127:GLY:O	2.18	0.43
1:D:129:ARG:NH2	1:E:101:ASP:HB3	2.27	0.43
1:C:12:VAL:HG11	1:D:31:PHE:HE2	1.84	0.43
1:A:77:ILE:HG22	1:A:79:LEU:H	1.83	0.42
1:D:287:GLY:HA3	1:D:438:PHE:CZ	2.54	0.42
1:B:42:ILE:HB	1:B:175:VAL:HG22	2.01	0.42
1:D:427:SER:HA	1:D:430:VAL:HG12	2.01	0.42
1:D:99:LEU:HB3	1:D:153:GLU:HB2	2.00	0.42
1:C:220:TYR:CZ	1:D:274:LYS:HG2	2.54	0.42
1:B:222:ILE:HA	1:B:226:TYR:HB2	2.00	0.42
1:D:160:THR:HG23	1:D:204:ALA:O	2.20	0.42
1:E:69:ASP:HB3	1:E:72:LEU:HD13	2.01	0.42
1:E:194:VAL:HB	1:E:209:SER:HB3	2.02	0.42
1:B:151:THR:HG22	1:B:211:SER:HB3	2.02	0.42
1:C:77:ILE:HG22	1:C:79:LEU:H	1.83	0.42
1:B:44:ILE:HG22	1:B:181:ILE:HD11	2.01	0.41
1:B:220:TYR:CZ	1:C:274:LYS:HG2	2.55	0.41
1:C:85:ASN:HB2	1:C:114:ARG:HB3	2.02	0.41
1:A:255:ILE:HD11	1:E:253:LEU:HG	2.03	0.41
1:B:19:LEU:HD22	1:B:72:LEU:HA	2.02	0.41
1:D:284:TYR:HB2	1:D:442:TYR:CE1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:38:VAL:HG11	1:D:94:PRO:HG2	2.03	0.41
1:B:301:PHE:HZ	1:B:312:ARG:HH21	1.67	0.41
1:E:147:GLU:HG3	1:E:215:LYS:HD2	2.02	0.41
1:B:268:LEU:HD21	1:B:281:ILE:CG2	2.52	0.41
1:D:437:LEU:HA	1:D:437:LEU:HD12	1.95	0.40
1:B:143:TYR:O	1:B:280:ALA:HB3	2.22	0.40
1:A:12:VAL:HG11	1:B:31:PHE:CE2	2.56	0.40
1:C:249:ALA:HB1	1:D:251:VAL:HG21	2.03	0.40
1:D:272:LEU:HD13	1:D:279:LYS:NZ	2.37	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	338/355 (95%)	329 (97%)	9 (3%)	0	100	100
1	B	329/355 (93%)	318 (97%)	11 (3%)	0	100	100
1	C	329/355 (93%)	319 (97%)	10 (3%)	0	100	100
1	D	330/355 (93%)	319 (97%)	11 (3%)	0	100	100
1	E	329/355 (93%)	319 (97%)	10 (3%)	0	100	100
All	All	1655/1775 (93%)	1604 (97%)	51 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	304/314 (97%)	292 (96%)	12 (4%)	39	76
1	B	296/314 (94%)	287 (97%)	9 (3%)	48	82
1	C	296/314 (94%)	285 (96%)	11 (4%)	41	77
1	D	297/314 (95%)	286 (96%)	11 (4%)	41	77
1	E	296/314 (94%)	285 (96%)	11 (4%)	41	77
All	All	1489/1570 (95%)	1435 (96%)	54 (4%)	42	78

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	SER
1	A	40	MET
1	A	99	LEU
1	A	112	LYS
1	A	115	MET
1	A	161	THR
1	A	185	GLN
1	A	211	SER
1	A	213	ARG
1	A	244	TYR
1	A	281	ILE
1	A	305	ILE
1	B	12	VAL
1	B	40	MET
1	B	59	LEU
1	B	99	LEU
1	B	161	THR
1	B	185	GLN
1	B	189	VAL
1	B	244	TYR
1	B	281	ILE
1	C	12	VAL
1	C	40	MET
1	C	41	ASN
1	C	49	MET
1	C	59	LEU
1	C	99	LEU
1	C	112	LYS

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Mol	Chain	Res	Type
1	C	123	THR
1	C	185	GLN
1	C	211	SER
1	C	244	TYR
1	D	12	VAL
1	D	40	MET
1	D	49	MET
1	D	59	LEU
1	D	99	LEU
1	D	185	GLN
1	D	189	VAL
1	D	211	SER
1	D	244	TYR
1	D	261	MET
1	D	281	ILE
1	E	12	VAL
1	E	17	ASP
1	E	40	MET
1	E	49	MET
1	E	59	LEU
1	E	99	LEU
1	E	110	THR
1	E	120	PRO
1	E	164	ILE
1	E	185	GLN
1	E	244	TYR

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

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

15 carbohydrates are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
5	NAG	A	3000	1,5	14,14,15	0.54	0	15,19,21	1.39	2 (13%)
5	NAG	A	3001	5	14,14,15	0.56	0	15,19,21	1.43	2 (13%)
5	BMA	A	3002	5	11,11,12	0.42	0	14,15,17	0.73	0
5	NAG	B	3000	1,5	14,14,15	0.50	0	15,19,21	1.61	3 (20%)
5	NAG	B	3001	5	14,14,15	0.62	0	15,19,21	1.33	1 (6%)
5	BMA	B	3002	5	11,11,12	0.40	0	14,15,17	0.72	0
5	NAG	C	3000	1,5	14,14,15	0.55	0	15,19,21	1.69	3 (20%)
5	NAG	C	3001	5	14,14,15	0.62	0	15,19,21	1.46	2 (13%)
5	BMA	C	3002	5	11,11,12	0.40	0	14,15,17	0.64	0
5	NAG	D	3000	1,5	14,14,15	0.54	0	15,19,21	1.61	3 (20%)
5	NAG	D	3001	5	14,14,15	0.59	0	15,19,21	1.44	2 (13%)
5	BMA	D	3002	5	11,11,12	0.41	0	14,15,17	0.75	0
5	NAG	E	3000	1,5	14,14,15	0.50	0	15,19,21	1.47	3 (20%)
5	NAG	E	3001	5	14,14,15	0.58	0	15,19,21	1.49	1 (6%)
5	BMA	E	3002	5	11,11,12	0.43	0	14,15,17	0.84	1 (7%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	3000	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	3001	5	-	0/6/23/26	0/1/1/1
5	BMA	A	3002	5	-	0/2/19/22	0/1/1/1
5	NAG	B	3000	1,5	-	0/6/23/26	0/1/1/1
5	NAG	B	3001	5	-	0/6/23/26	0/1/1/1
5	BMA	B	3002	5	-	0/2/19/22	0/1/1/1
5	NAG	C	3000	1,5	-	0/6/23/26	0/1/1/1
5	NAG	C	3001	5	-	0/6/23/26	0/1/1/1
5	BMA	C	3002	5	-	0/2/19/22	0/1/1/1
5	NAG	D	3000	1,5	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	D	3001	5	-	0/6/23/26	0/1/1/1
5	BMA	D	3002	5	-	0/2/19/22	0/1/1/1
5	NAG	E	3000	1,5	-	0/6/23/26	0/1/1/1
5	NAG	E	3001	5	-	0/6/23/26	0/1/1/1
5	BMA	E	3002	5	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	3001	NAG	C4-C3-C2	2.08	114.47	111.23
5	E	3000	NAG	C2-N2-C7	2.09	125.72	123.04
5	E	3000	NAG	C4-C3-C2	2.12	114.52	111.23
5	A	3000	NAG	C4-C3-C2	2.12	114.53	111.23
5	A	3001	NAG	C4-C3-C2	2.18	114.62	111.23
5	C	3001	NAG	C4-C3-C2	2.32	114.83	111.23
5	D	3000	NAG	C2-N2-C7	2.35	126.05	123.04
5	E	3002	BMA	C1-O5-C5	2.48	115.40	112.25
5	B	3000	NAG	C4-C3-C2	2.52	115.14	111.23
5	C	3000	NAG	C2-N2-C7	2.68	126.48	123.04
5	D	3000	NAG	C4-C3-C2	2.70	115.42	111.23
5	C	3000	NAG	C4-C3-C2	2.82	115.61	111.23
5	B	3000	NAG	C2-N2-C7	2.96	126.84	123.04
5	E	3000	NAG	C1-O5-C5	3.26	116.39	112.25
5	A	3000	NAG	C1-O5-C5	3.33	116.47	112.25
5	D	3000	NAG	C1-O5-C5	3.35	116.50	112.25
5	B	3000	NAG	C1-O5-C5	3.52	116.72	112.25
5	C	3000	NAG	C1-O5-C5	3.74	116.99	112.25
5	B	3001	NAG	C1-O5-C5	3.84	117.12	112.25
5	C	3001	NAG	C1-O5-C5	3.93	117.23	112.25
5	A	3001	NAG	C1-O5-C5	4.22	117.61	112.25
5	D	3001	NAG	C1-O5-C5	4.23	117.61	112.25
5	E	3001	NAG	C1-O5-C5	4.53	118.00	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.6 Ligand geometry

Of 16 ligands modelled in this entry, 5 are monoatomic - leaving 11 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
3	NAG	A	1000	1	14,14,15	0.57	0	15,19,21	1.86	5 (33%)
3	NAG	A	2000	1	14,14,15	0.57	0	15,19,21	1.76	4 (26%)
2	BEN	A	500	-	9,9,9	2.40	4 (44%)	9,11,11	1.23	1 (11%)
3	NAG	B	2000	1	14,14,15	0.52	0	15,19,21	1.68	3 (20%)
2	BEN	B	500	-	9,9,9	2.38	4 (44%)	9,11,11	0.90	0
3	NAG	C	2000	1	14,14,15	0.57	0	15,19,21	1.48	2 (13%)
2	BEN	C	500	-	9,9,9	2.17	4 (44%)	9,11,11	0.93	0
3	NAG	D	2000	1	14,14,15	0.54	0	15,19,21	1.59	2 (13%)
2	BEN	D	500	-	9,9,9	2.04	3 (33%)	9,11,11	0.87	0
3	NAG	E	2000	1	14,14,15	0.61	0	15,19,21	1.54	3 (20%)
2	BEN	E	500	-	9,9,9	1.99	3 (33%)	9,11,11	1.01	1 (11%)

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	1000	1	-	0/6/23/26	0/1/1/1
3	NAG	A	2000	1	-	0/6/23/26	0/1/1/1
2	BEN	A	500	-	-	0/4/4/4	0/1/1/1
3	NAG	B	2000	1	-	0/6/23/26	0/1/1/1
2	BEN	B	500	-	-	0/4/4/4	0/1/1/1
3	NAG	C	2000	1	-	0/6/23/26	0/1/1/1
2	BEN	C	500	-	-	0/4/4/4	0/1/1/1
3	NAG	D	2000	1	-	0/6/23/26	0/1/1/1
2	BEN	D	500	-	-	0/4/4/4	0/1/1/1
3	NAG	E	2000	1	-	0/6/23/26	0/1/1/1
2	BEN	E	500	-	-	0/4/4/4	0/1/1/1

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	BEN	C5-C4	2.01	1.43	1.38
2	C	500	BEN	C5-C6	2.03	1.43	1.38
2	A	500	BEN	C-N2	2.26	1.39	1.33
2	E	500	BEN	C-N2	2.33	1.39	1.33
2	A	500	BEN	C1-C	2.39	1.51	1.47
2	C	500	BEN	C6-C1	2.48	1.43	1.39
2	D	500	BEN	C6-C1	2.67	1.43	1.39
2	D	500	BEN	C-N2	2.68	1.40	1.33
2	C	500	BEN	C-N2	2.86	1.40	1.33
2	E	500	BEN	C2-C1	3.11	1.44	1.39
2	B	500	BEN	C-N2	3.11	1.41	1.33
2	E	500	BEN	C6-C1	3.13	1.44	1.39
2	B	500	BEN	C6-C1	3.43	1.45	1.39
2	A	500	BEN	C2-C1	3.45	1.45	1.39
2	C	500	BEN	C2-C1	3.59	1.45	1.39
2	D	500	BEN	C2-C1	3.64	1.45	1.39
2	B	500	BEN	C2-C1	3.99	1.46	1.39
2	A	500	BEN	C6-C1	4.05	1.46	1.39

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2000	NAG	O7-C7-N2	2.01	125.96	121.86
3	A	1000	NAG	C3-C2-N2	2.14	115.69	110.56
3	E	2000	NAG	C2-N2-C7	2.23	125.90	123.04
3	A	1000	NAG	C4-C3-C2	2.23	114.70	111.23
3	B	2000	NAG	C3-C4-C5	2.24	114.10	110.20
3	C	2000	NAG	C4-C3-C2	2.37	114.91	111.23
2	E	500	BEN	C1-C-N2	2.37	121.90	118.11
3	D	2000	NAG	C4-C3-C2	2.46	115.06	111.23
3	A	2000	NAG	C3-C4-C5	2.47	114.50	110.20
3	A	2000	NAG	C4-C3-C2	2.71	115.45	111.23
3	B	2000	NAG	C4-C3-C2	2.76	115.52	111.23
3	E	2000	NAG	C1-O5-C5	2.86	115.88	112.25
3	E	2000	NAG	C4-C3-C2	2.96	115.84	111.23
3	A	1000	NAG	C3-C4-C5	3.22	115.80	110.20
2	A	500	BEN	C1-C-N2	3.63	123.91	118.11
3	A	1000	NAG	C1-O5-C5	3.68	116.92	112.25
3	C	2000	NAG	C1-O5-C5	3.71	116.96	112.25
3	A	1000	NAG	C2-N2-C7	3.88	128.02	123.04
3	D	2000	NAG	C1-O5-C5	4.06	117.40	112.25
3	B	2000	NAG	C1-O5-C5	4.33	117.75	112.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2000	NAG	C1-O5-C5	4.51	117.97	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	340/355 (95%)	0.40	7 (2%) 67 44	68, 99, 127, 158	0
1	B	331/355 (93%)	0.59	15 (4%) 37 20	69, 93, 135, 158	0
1	C	331/355 (93%)	0.67	17 (5%) 32 17	67, 94, 130, 168	0
1	D	332/355 (93%)	0.59	14 (4%) 40 22	70, 102, 139, 191	0
1	E	331/355 (93%)	0.60	17 (5%) 32 17	66, 103, 138, 161	0
All	All	1665/1775 (93%)	0.57	70 (4%) 40 22	66, 99, 136, 191	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	11	PHE	4.4
1	B	446	TYR	4.3
1	B	145	LEU	3.8
1	C	16	VAL	3.8
1	E	77	ILE	3.7
1	B	14	GLU	3.7
1	C	81	LEU	3.5
1	E	206	PRO	3.4
1	A	25	ILE	3.3
1	E	25	ILE	3.2
1	D	202	THR	3.2
1	C	145	LEU	3.1
1	E	14	GLU	2.9
1	B	228	PRO	2.9
1	B	11	PHE	2.8
1	A	188	ILE	2.8
1	C	147	GLU	2.8
1	D	77	ILE	2.7
1	D	446	TYR	2.7
1	B	142	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	189	VAL	2.6
1	A	212	PHE	2.6
1	E	138	MET	2.6
1	C	132	THR	2.5
1	C	19	LEU	2.5
1	D	23	TYR	2.4
1	B	68	ARG	2.4
1	C	182	GLU	2.4
1	B	281	ILE	2.4
1	C	83	LEU	2.4
1	B	227	MET	2.4
1	C	144	PRO	2.3
1	E	166	PHE	2.3
1	E	307	PHE	2.3
1	C	146	ASP	2.3
1	B	214	LEU	2.3
1	D	116	ILE	2.3
1	D	145	LEU	2.3
1	E	140	LEU	2.3
1	E	214	LEU	2.3
1	B	445	TYR	2.2
1	A	312	ARG	2.2
1	D	212	PHE	2.2
1	C	446	TYR	2.2
1	E	150	CYS	2.2
1	C	281	ILE	2.2
1	E	183	LEU	2.2
1	D	136	CYS	2.2
1	D	81	LEU	2.2
1	E	296	LEU	2.2
1	B	280	ALA	2.2
1	D	137	MET	2.1
1	B	200	PHE	2.1
1	E	233	THR	2.1
1	D	27	LEU	2.1
1	A	24	ASP	2.1
1	C	234	ILE	2.1
1	D	281	ILE	2.1
1	E	188	ILE	2.1
1	D	118	LEU	2.1
1	E	18	LYS	2.1
1	C	27	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	198	VAL	2.0
1	C	260	THR	2.0
1	B	50	VAL	2.0
1	B	231	LEU	2.0
1	D	214	LEU	2.0
1	C	228	PRO	2.0
1	E	437	LEU	2.0
1	A	215	LYS	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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
5	NAG	C	3000	14/15	0.92	0.26	-0.08	123,125,130,138	0
5	NAG	A	3000	14/15	0.94	0.18	-0.48	113,115,118,121	0
5	NAG	B	3000	14/15	0.96	0.20	-0.49	107,112,118,124	0
5	NAG	E	3000	14/15	0.94	0.17	-0.55	108,114,119,124	0
5	NAG	D	3000	14/15	0.96	0.17	-0.64	101,111,117,120	0
5	BMA	A	3002	11/12	0.58	0.19	-	155,158,160,160	0
5	BMA	C	3002	11/12	0.66	0.45	-	167,169,171,172	0
5	NAG	D	3001	14/15	0.89	0.15	-	124,132,141,151	0
5	NAG	C	3001	14/15	0.83	0.33	-	134,147,153,161	0
5	NAG	E	3001	14/15	0.91	0.17	-	120,132,140,150	0
5	NAG	B	3001	14/15	0.88	0.30	-	115,131,140,149	0
5	NAG	A	3001	14/15	0.88	0.18	-	120,128,137,148	0
5	BMA	B	3002	11/12	0.62	0.27	-	155,157,160,161	0
5	BMA	D	3002	11/12	0.63	0.16	-	159,163,163,164	0
5	BMA	E	3002	11/12	0.60	0.19	-	157,160,161,162	0

## 6.4 Ligands ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	BEN	B	500	9/9	0.73	0.39	2.29	125,126,126,126	0
2	BEN	A	500	9/9	0.78	0.28	2.03	135,136,136,136	0
2	BEN	C	500	9/9	0.74	0.23	0.45	129,129,129,129	0
3	NAG	A	2000	14/15	0.94	0.21	0.14	114,117,124,124	0
3	NAG	A	1000	14/15	0.86	0.17	0.13	144,147,148,149	0
4	CL	C	1448	1/1	0.87	0.26	0.09	109,109,109,109	0
4	CL	A	1448	1/1	0.91	0.23	-0.19	97,97,97,97	0
2	BEN	E	500	9/9	0.77	0.20	-0.76	138,139,139,139	0
4	CL	B	1448	1/1	0.81	0.20	-0.92	100,100,100,100	0
3	NAG	E	2000	14/15	0.91	0.14	-1.04	136,139,141,142	0
2	BEN	D	500	9/9	0.88	0.16	-1.23	117,117,118,118	0
4	CL	E	1448	1/1	0.72	0.16	-1.75	100,100,100,100	0
4	CL	D	1449	1/1	0.86	0.14	-1.77	111,111,111,111	0
3	NAG	D	2000	14/15	0.88	0.21	-	158,162,166,166	0
3	NAG	B	2000	14/15	0.89	0.20	-	152,156,157,157	0
3	NAG	C	2000	14/15	0.86	0.18	-	147,151,154,156	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.