



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 27, 2019 – 02:42 PM EDT

PDB ID : 6QAD
Title : Human Butyrylcholinesterase in complex with ((S)-2-(butylamino)-N-(2-(4-(dimethylamino)cyclohexyl)ethyl)-3-(1H-indol-3-yl)propanamide
Authors : Brazzolotto, X.; Nachon, F.; Harst, M.; Knez, D.; Gobec, S.
Deposited on : 2018-12-19
Resolution : 2.50 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 : 1.8.0 (224370), CSD as540be (2019)
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

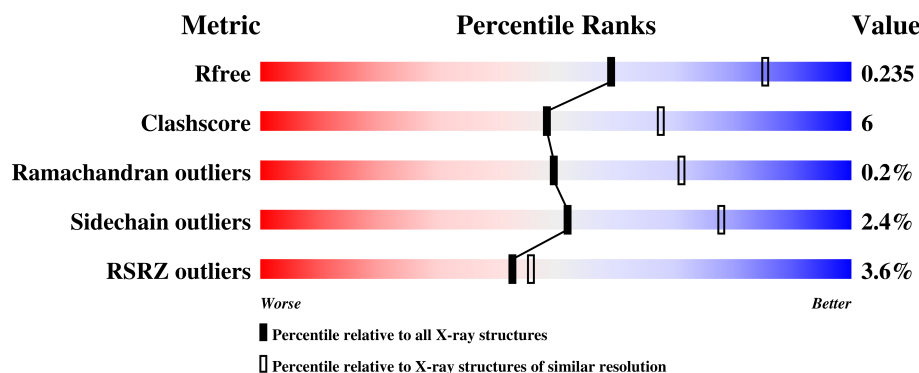
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4155 (2.50-2.50)
Clashscore	122126	4827 (2.50-2.50)
Ramachandran outliers	120053	4735 (2.50-2.50)
Sidechain outliers	120020	4737 (2.50-2.50)
RSRZ outliers	108989	4058 (2.50-2.50)

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	557	

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
3	FUC	A	602	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	GOL	A	613	-	-	-	X
5	GOL	A	616	-	-	-	X
6	MES	A	617	-	-	-	X

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 4530 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.

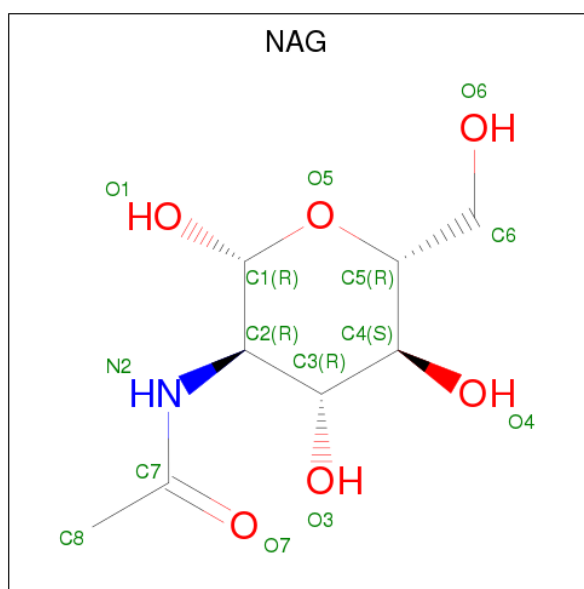
- Molecule 1 is a protein called Cholinesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	526	4233	2734	711	772	16	0	6	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-26	ASP	HIS	conflict	UNP P06276
A	17	GLN	ASN	engineered mutation	UNP P06276
A	455	GLN	ASN	engineered mutation	UNP P06276
A	481	GLN	ASN	engineered mutation	UNP P06276
A	486	GLN	ASN	engineered mutation	UNP P06276

- Molecule 2 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



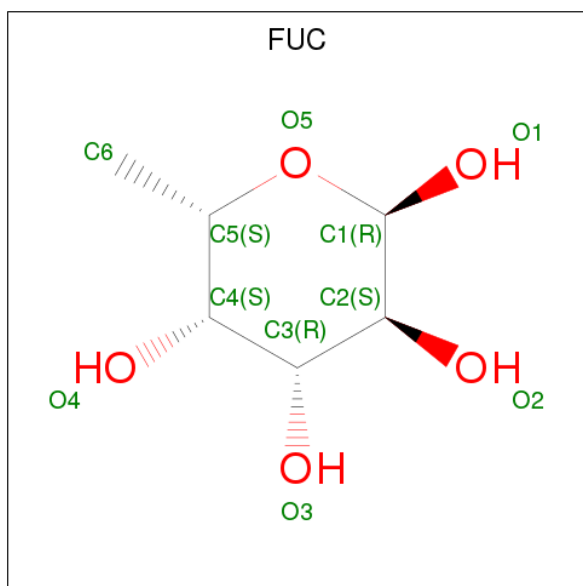
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	14	8	1	5	0	0

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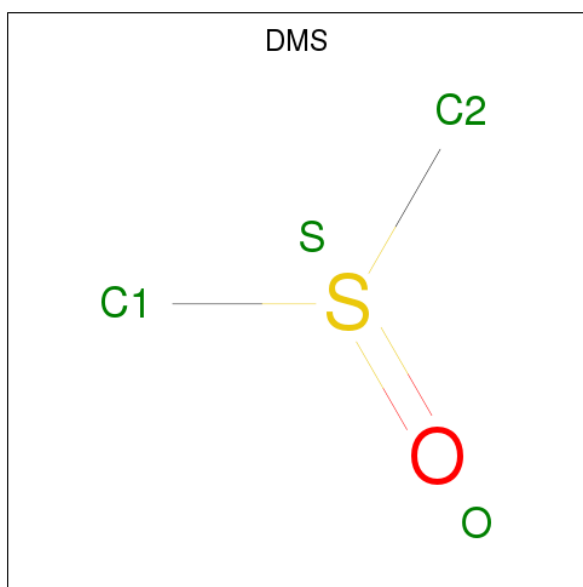
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is ALPHA-L-FUCOSE (three-letter code: FUC) (formula: C₆H₁₂O₅).



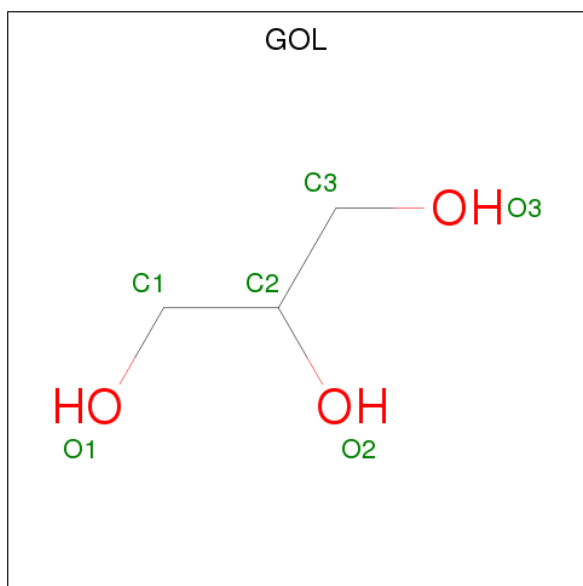
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			10	6	4		
3	A	1	Total	C	O	0	0
			10	6	4		
3	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 4 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C₂H₆OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	S	0	0
			4	2	1	1		

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



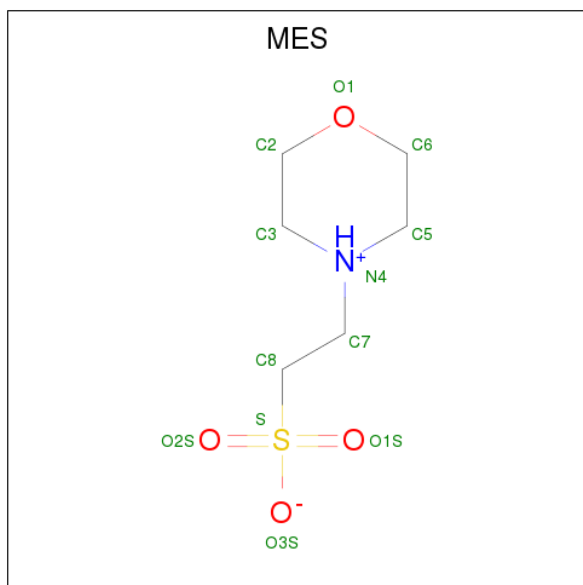
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		

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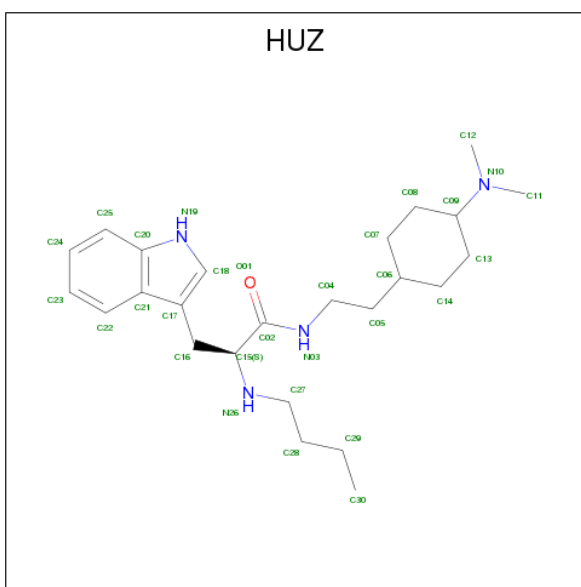
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



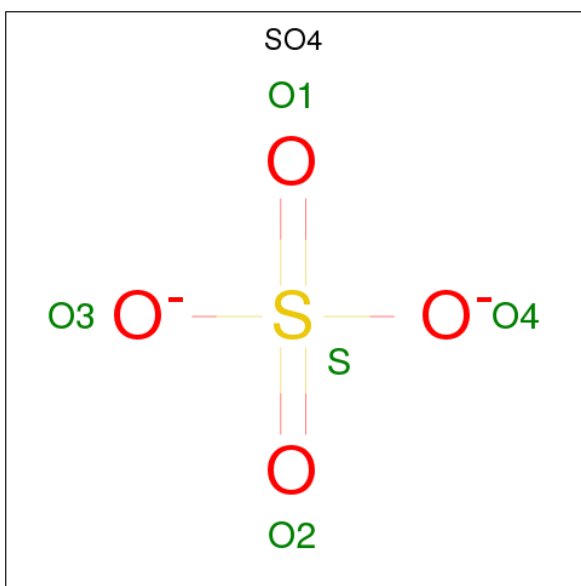
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 7 is (2 {S})-2-(butylamino)- {N}-[2-[4-(dimethylamino)cyclohexyl]ethyl]-3-(1 {H})-indol-3-yl)propanamide (three-letter code: HUZ) (formula: C₂₅H₄₀N₄O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			30	25	4	1		

- Molecule 8 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	O	S	0	0
			5	4	1		
8	A	1	Total	O	S	0	0
			5	4	1		
8	A	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	O	S	0	0
			5	4	1		

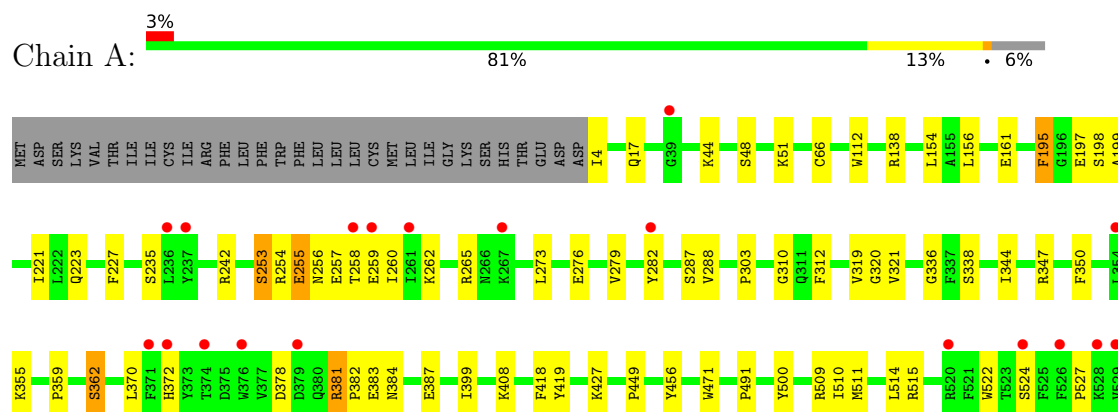
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	65	Total	O	0	0
			65	65		

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: Cholinesterase



4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	155.27Å 155.27Å 129.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	99.64 – 2.50 99.64 – 2.50	Depositor EDS
% Data completeness (in resolution range)	90.7 (99.64-2.50) 90.7 (99.64-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.48Å)	Xtriage
Refinement program	PHENIX (1.14_3260: ???)	Depositor
R, R_{free}	0.200 , 0.234 0.199 , 0.235	Depositor DCC
R_{free} test set	1248 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	49.7	Xtriage
Anisotropy	0.705	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4530	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, NAG, DMS, FUC, MES, HUZ, SO4

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	0.28	0/4363	0.48	0/5922

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	255	GLU	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4233	0	4133	56	0
2	A	112	0	99	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	30	0	30	0	0
4	A	4	0	6	0	0
5	A	24	0	32	4	0
6	A	12	0	12	2	0
7	A	30	0	0	1	0
8	A	20	0	0	1	0
9	A	65	0	0	2	0
All	All	4530	0	4312	57	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 6.

All (57) 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:256:ASN:HB3	1:A:259:GLU:HB2	1.56	0.88
1:A:255:GLU:N	1:A:259:GLU:OE2	2.10	0.84
1:A:427:LYS:HB2	5:A:613:GOL:H11	1.72	0.71
1:A:51:LYS:HD2	1:A:51:LYS:H	1.56	0.71
1:A:254:ARG:HB3	1:A:259:GLU:CD	2.12	0.70
1:A:254:ARG:HB3	1:A:259:GLU:OE2	1.94	0.66
1:A:198:SER:OG	9:A:701:HOH:O	2.15	0.64
1:A:161:GLU:HG3	1:A:258:THR:HG23	1.80	0.63
1:A:242:ARG:HH22	5:A:615:GOL:H32	1.65	0.62
1:A:344:ILE:HD13	1:A:382:PRO:HB2	1.84	0.60
1:A:254:ARG:CB	1:A:259:GLU:OE2	2.50	0.60
1:A:288:VAL:H	5:A:615:GOL:H31	1.65	0.60
1:A:287:SER:HA	5:A:615:GOL:H2	1.86	0.58
1:A:282:TYR:HB3	1:A:359:PRO:HG2	1.87	0.57
1:A:156:LEU:HD13	1:A:257:GLU:HB3	1.87	0.57
1:A:514:LEU:HG	1:A:515:ARG:HG3	1.86	0.56
1:A:509:ARG:NH1	9:A:702:HOH:O	2.23	0.56
1:A:319:VAL:O	1:A:418:PHE:HA	2.07	0.55
1:A:378:ASP:O	1:A:384:ASN:ND2	2.34	0.54
1:A:227:PHE:CE2	1:A:303:PRO:HB2	2.43	0.53
1:A:287:SER:O	7:A:618:HUZ:N26	2.44	0.51
1:A:522:TRP:O	1:A:527:PRO:HD3	2.12	0.50
1:A:500:TYR:CZ	1:A:511[B]:MET:HB2	2.46	0.50
1:A:255:GLU:H	1:A:259:GLU:CD	2.15	0.49
1:A:66:CYS:HB3	1:A:273:LEU:HD11	1.94	0.49
1:A:254:ARG:HB2	1:A:260:ILE:HG13	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:514:LEU:HD12	6:A:617:MES:H52	1.96	0.47
1:A:227:PHE:CD2	1:A:303:PRO:HB2	2.50	0.46
1:A:138:ARG:HH11	1:A:138:ARG:HG2	1.80	0.46
1:A:321:VAL:HG11	1:A:399:ILE:HA	1.98	0.46
1:A:198:SER:OG	1:A:199:ALA:N	2.49	0.45
1:A:500:TYR:CZ	1:A:511[A]:MET:HB2	2.50	0.45
1:A:383:GLU:CD	1:A:383:GLU:H	2.21	0.44
1:A:310:GLY:HA2	1:A:312:PHE:CE2	2.53	0.44
1:A:282:TYR:HD2	1:A:359:PRO:HB2	1.82	0.44
1:A:355:LYS:HE3	1:A:355:LYS:HB3	1.67	0.43
1:A:372[B]:HIS:NE2	8:A:622:SO4:O3	2.41	0.43
1:A:254:ARG:HB3	1:A:259:GLU:CG	2.48	0.43
1:A:276:GLU:O	1:A:279:VAL:HG22	2.18	0.43
1:A:253:SER:C	1:A:254:ARG:HG2	2.39	0.42
1:A:320:GLY:HA3	1:A:419:TYR:CE1	2.54	0.42
1:A:44:LYS:HD3	1:A:265:ARG:NH2	2.34	0.42
1:A:350:PHE:CD2	1:A:370:LEU:HD12	2.55	0.42
1:A:138:ARG:NH1	1:A:138:ARG:HG2	2.35	0.42
1:A:336:GLY:H	2:A:604:NAG:C8	2.33	0.42
1:A:491:PRO:HD3	1:A:510:ILE:HD12	2.02	0.42
1:A:197:GLU:HA	1:A:223:GLN:O	2.19	0.42
1:A:449:PRO:HA	1:A:456:TYR:CD1	2.55	0.41
1:A:112:TRP:HA	1:A:195:PHE:O	2.21	0.41
1:A:195:PHE:CB	1:A:221:ILE:HB	2.51	0.41
1:A:336:GLY:H	2:A:604:NAG:H83	1.85	0.41
1:A:256:ASN:O	1:A:259:GLU:OE2	2.38	0.41
1:A:4:ILE:HD12	1:A:17:GLN:HG2	2.03	0.41
6:A:617:MES:H32	6:A:617:MES:H82	1.47	0.40
1:A:347:ARG:HH12	1:A:370:LEU:HD21	1.86	0.40
1:A:262:LYS:HA	1:A:262:LYS:HD2	1.95	0.40
1:A:381:ARG:NH1	1:A:387:GLU:OE1	2.54	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	530/557 (95%)	507 (96%)	22 (4%)	1 (0%)	49 71

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	362	SER

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	456/481 (95%)	445 (98%)	11 (2%)	52 78

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	SER
1	A	154	LEU
1	A	195	PHE
1	A	235	SER
1	A	253	SER
1	A	338	SER
1	A	362	SER
1	A	381	ARG
1	A	408	LYS
1	A	471	TRP
1	A	524	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	A	517	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 ⓘ

22 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	A	601	1,3	14,14,15	0.28	0	17,19,21	0.46	0
3	FUC	A	602	2	9,10,11	0.82	0	13,14,16	0.93	0
2	NAG	A	603	1	14,14,15	0.46	0	17,19,21	0.53	0
2	NAG	A	604	2	14,14,15	0.20	0	17,19,21	0.58	0
2	NAG	A	605	1,3,2	14,14,15	0.41	0	17,19,21	0.40	0
3	FUC	A	606	2	9,10,11	0.78	0	13,14,16	0.83	0
2	NAG	A	607	1,3,2	14,14,15	0.88	1 (7%)	17,19,21	1.53	3 (17%)
2	NAG	A	608	2	14,14,15	1.21	1 (7%)	17,19,21	0.84	1 (5%)
3	FUC	A	609	2	9,10,11	0.59	0	13,14,16	0.90	1 (7%)
2	NAG	A	610	1	14,14,15	0.60	1 (7%)	17,19,21	0.54	0
2	NAG	A	611	1	14,14,15	0.44	0	17,19,21	0.56	0
4	DMS	A	612	-	3,3,3	0.64	0	3,3,3	0.57	0
5	GOL	A	613	-	5,5,5	0.87	0	5,5,5	1.12	1 (20%)
5	GOL	A	614	-	5,5,5	0.90	0	5,5,5	0.99	0
5	GOL	A	615	-	5,5,5	0.88	0	5,5,5	1.07	0
5	GOL	A	616	-	5,5,5	0.88	0	5,5,5	1.01	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	MES	A	617	-	12,12,12	2.26	1 (8%)	14,16,16	2.23	7 (50%)
7	HUZ	A	618	-	30,32,32	1.20	3 (10%)	35,42,42	1.01	1 (2%)
8	SO4	A	619	-	4,4,4	0.16	0	6,6,6	0.09	0
8	SO4	A	620	-	4,4,4	0.17	0	6,6,6	0.06	0
8	SO4	A	621	-	4,4,4	0.17	0	6,6,6	0.05	0
8	SO4	A	622	-	4,4,4	0.17	0	6,6,6	0.06	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	601	1,3	-	0/6/23/26	0/1/1/1
3	FUC	A	602	2	-	0/0/17/20	0/1/1/1
2	NAG	A	603	1	-	0/6/23/26	0/1/1/1
2	NAG	A	604	2	-	0/6/23/26	0/1/1/1
2	NAG	A	605	1,3,2	-	0/6/23/26	0/1/1/1
3	FUC	A	606	2	-	0/0/17/20	0/1/1/1
2	NAG	A	607	1,3,2	-	0/6/23/26	0/1/1/1
2	NAG	A	608	2	-	0/6/23/26	0/1/1/1
3	FUC	A	609	2	-	0/0/17/20	0/1/1/1
2	NAG	A	610	1	-	0/6/23/26	0/1/1/1
2	NAG	A	611	1	-	0/6/23/26	0/1/1/1
4	DMS	A	612	-	-	0/0/0/0	0/0/0/0
5	GOL	A	613	-	-	0/4/4/4	0/0/0/0
5	GOL	A	614	-	-	0/4/4/4	0/0/0/0
5	GOL	A	615	-	-	0/4/4/4	0/0/0/0
5	GOL	A	616	-	-	0/4/4/4	0/0/0/0
6	MES	A	617	-	-	0/6/14/14	0/1/1/1
7	HUZ	A	618	-	-	0/22/33/33	0/3/3/3
8	SO4	A	619	-	-	0/0/0/0	0/0/0/0
8	SO4	A	620	-	-	0/0/0/0	0/0/0/0
8	SO4	A	621	-	-	0/0/0/0	0/0/0/0
8	SO4	A	622	-	-	0/0/0/0	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	617	MES	C8-S	-7.58	1.66	1.77
2	A	608	NAG	O5-C1	-4.36	1.36	1.43
2	A	610	NAG	C1-C2	2.04	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	618	HUZ	C18-N19	2.08	1.40	1.36
7	A	618	HUZ	C15-C02	2.31	1.59	1.52
2	A	607	NAG	O5-C1	2.89	1.48	1.43
7	A	618	HUZ	C02-N03	3.73	1.41	1.33

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	617	MES	C2-C3-N4	-3.11	105.33	110.09
2	A	607	NAG	C4-C3-C2	-2.25	107.72	111.02
5	A	613	GOL	C3-C2-C1	-2.11	103.51	111.75
7	A	618	HUZ	C24-C25-C20	-2.09	117.03	120.08
6	A	617	MES	C6-C5-N4	-2.04	106.97	110.09
3	A	609	FUC	O5-C1-C2	2.01	113.89	110.79
6	A	617	MES	O3S-S-C8	2.07	109.12	105.77
2	A	608	NAG	C3-C4-C5	2.37	114.48	110.23
6	A	617	MES	C7-N4-C5	2.55	117.77	111.23
6	A	617	MES	O1S-S-C8	2.86	110.36	106.92
2	A	607	NAG	C1-O5-C5	3.23	116.60	112.20
6	A	617	MES	C7-N4-C3	3.27	119.60	111.23
2	A	607	NAG	O4-C4-C5	3.98	119.25	109.29
6	A	617	MES	C5-N4-C3	4.20	118.13	108.86

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	604	NAG	2	0
5	A	613	GOL	1	0
5	A	615	GOL	3	0
6	A	617	MES	2	0
7	A	618	HUZ	1	0
8	A	622	SO4	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	526/557 (94%)	0.51	19 (3%)	42 46	37, 52, 76, 111	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	261	ILE	4.6
1	A	529	VAL	4.6
1	A	258	THR	4.2
1	A	376	TRP	3.8
1	A	528	LYS	3.8
1	A	379	ASP	3.8
1	A	374	THR	3.5
1	A	259	GLU	3.2
1	A	236	LEU	3.2
1	A	524	SER	3.1
1	A	526	PHE	2.9
1	A	267	LYS	2.8
1	A	282	TYR	2.4
1	A	372[A]	HIS	2.4
1	A	237	TYR	2.3
1	A	520	ARG	2.3
1	A	39	GLY	2.2
1	A	371	PHE	2.1
1	A	354	LEU	2.0

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

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
5	GOL	A	613	6/6	0.48	0.47	68,70,78,81	0
6	MES	A	617	12/12	0.59	0.40	59,96,123,143	0
3	FUC	A	602	10/11	0.62	0.45	119,126,134,138	0
2	NAG	A	608	14/15	0.65	0.35	102,118,128,128	0
2	NAG	A	610	14/15	0.67	0.26	100,110,117,119	0
5	GOL	A	616	6/6	0.76	0.45	72,79,92,107	0
5	GOL	A	614	6/6	0.76	0.39	61,76,79,84	0
2	NAG	A	601	14/15	0.77	0.28	93,104,112,122	0
2	NAG	A	604	14/15	0.79	0.22	67,82,91,91	0
2	NAG	A	611	14/15	0.81	0.23	66,79,89,90	0
2	NAG	A	603	14/15	0.81	0.18	55,77,83,86	0
3	FUC	A	606	10/11	0.83	0.48	71,84,91,92	0
5	GOL	A	615	6/6	0.85	0.24	60,67,74,81	0
8	SO4	A	622	5/5	0.86	0.64	81,97,114,128	0
3	FUC	A	609	10/11	0.87	0.20	96,108,115,118	0
7	HUZ	A	618	30/30	0.88	0.22	44,55,68,77	0
8	SO4	A	620	5/5	0.89	0.38	83,94,104,109	0
2	NAG	A	607	14/15	0.89	0.14	85,94,109,109	0
8	SO4	A	621	5/5	0.90	0.15	89,98,110,112	0
2	NAG	A	605	14/15	0.92	0.18	51,62,77,84	0
8	SO4	A	619	5/5	0.93	0.31	70,79,91,94	0
4	DMS	A	612	4/4	0.95	0.22	42,53,75,77	0

6.5 Other polymers [i](#)

There are no such residues in this entry.