



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2018 – 11:40 am GMT

PDB ID : 3D7F
Title : A high resolution crystal structure of human glutamate carboxypeptidase II (GCPII) in a complex with DCIT, a urea-based inhibitor
Authors : Lubkowski, J.; Barinka, C.
Deposited on : 2008-05-21
Resolution : 1.54 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : trunk30967
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) : trunk30967

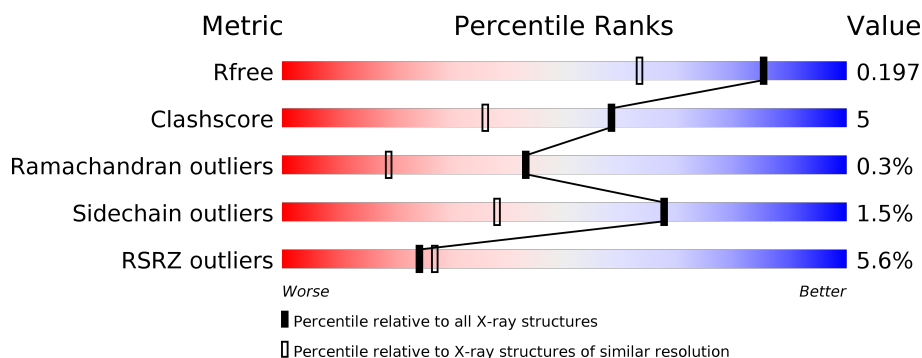
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 1.54 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	2055 (1.56-1.52)
Clashscore	122126	2124 (1.56-1.52)
Ramachandran outliers	120053	2083 (1.56-1.52)
Sidechain outliers	120020	2081 (1.56-1.52)
RSRZ outliers	108989	2027 (1.56-1.52)

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	709	

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 6603 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 Glutamate carboxypeptidase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	691	5729	3687	951	1067	24	0	61	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	42	ARG	-	EXPRESSION TAG	UNP Q04609
A	43	SER	-	EXPRESSION TAG	UNP Q04609

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



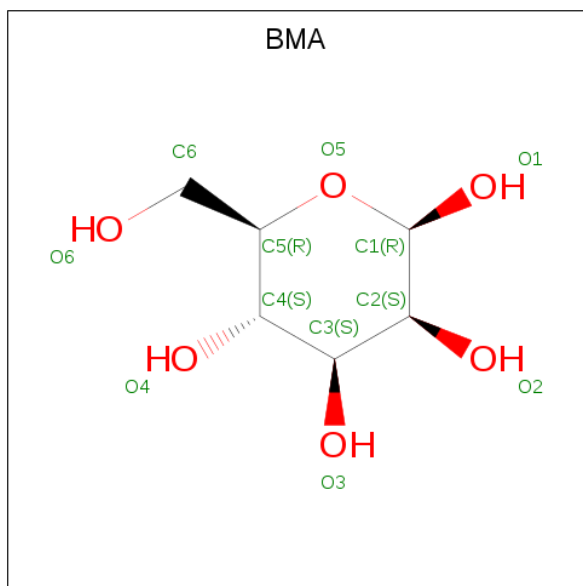
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	14	8	1	5	0	0
2	A	1	14	8	1	5	0	0

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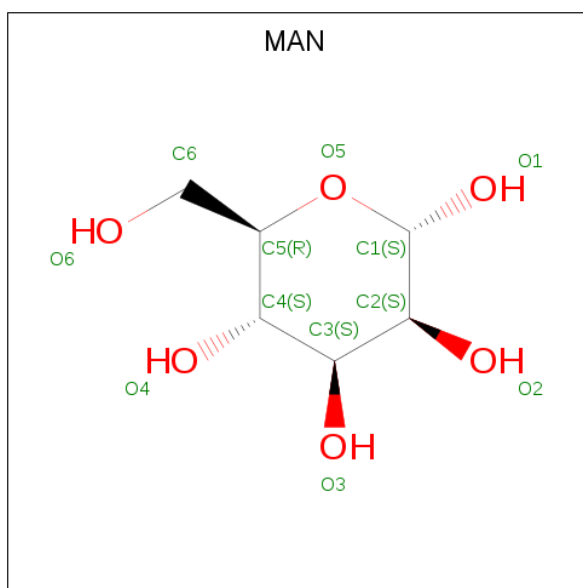
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is BETA-D-MANNOSE (three-letter code: BMA) (formula: C₆H₁₂O₆).



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

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



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

- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Zn	0	0
			2	2		

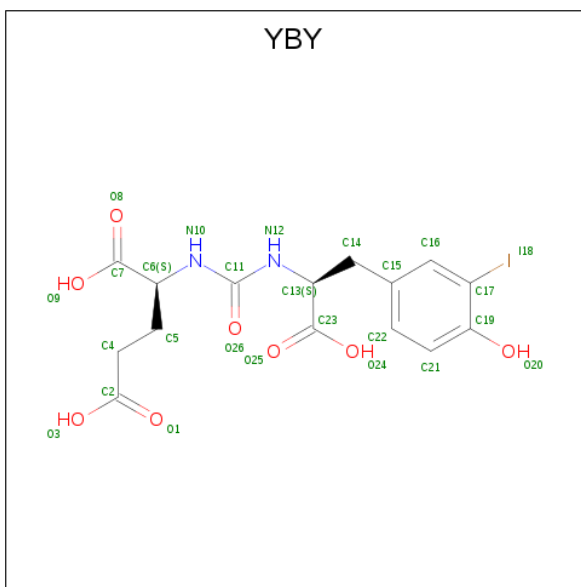
- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Ca	0	0
			1	1		

- Molecule 7 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Cl	0	0
			1	1		

- Molecule 8 is N-[(1S)-1-carboxy-2-(4-hydroxy-3-iodophenyl)ethyl]carbamoyl-L-glutamic acid (three-letter code: YBY) (formula: $C_{15}H_{17}IN_2O_8$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	A	1	Total	C	I	N	O	0	0
			26	15	1	2	8		

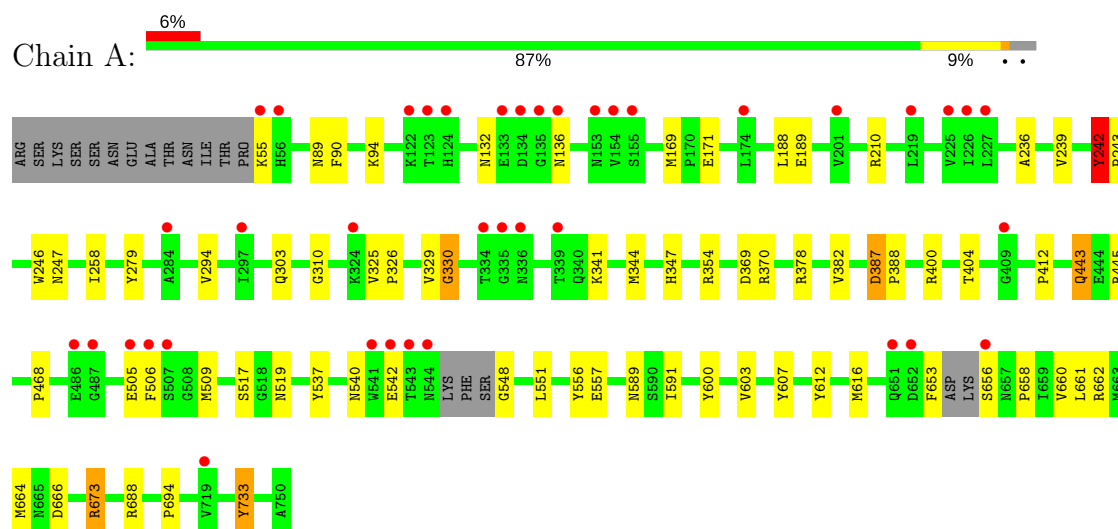
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	657	Total	O	0	3
			657	657		

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: Glutamate carboxypeptidase 2



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	101.43Å 130.06Å 158.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 1.54 29.48 – 1.54	Depositor EDS
% Data completeness (in resolution range)	94.6 (15.00-1.54) 94.5 (29.48-1.54)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.68 (at 1.54Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.177 , 0.195 0.178 , 0.197	Depositor DCC
R_{free} test set	1447 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å ²)	20.1	Xtriage
Anisotropy	0.510	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 54.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6603	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% 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: YBY, ZN, BMA, NAG, CL, CA, 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	0.89	3/6112 (0.0%)	0.89	7/8274 (0.1%)

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	5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	733	TYR	CD1-CE1	5.44	1.47	1.39
1	A	443[A]	GLN	CB-CG	-5.21	1.38	1.52
1	A	443[B]	GLN	CB-CG	-5.21	1.38	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	673	ARG	NE-CZ-NH2	-8.16	116.22	120.30
1	A	370	ARG	NE-CZ-NH1	7.27	123.93	120.30
1	A	242	TYR	C-N-CD	-7.07	105.04	120.60
1	A	370	ARG	NE-CZ-NH2	-7.03	116.79	120.30
1	A	445	ARG	NE-CZ-NH1	5.62	123.11	120.30
1	A	354	ARG	NE-CZ-NH1	-5.28	117.66	120.30
1	A	210	ARG	NE-CZ-NH1	-5.20	117.70	120.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	242	TYR	Mainchain,Peptide
1	A	330	GLY	Mainchain,Peptide
1	A	387	ASP	Mainchain

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	5729	0	5597	59	2
2	A	154	0	137	4	0
3	A	22	0	19	0	0
4	A	11	0	10	0	0
5	A	2	0	0	0	0
6	A	1	0	0	0	0
7	A	1	0	0	0	0
8	A	26	0	14	0	0
9	A	657	0	0	13	4
All	All	6603	0	5777	62	4

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 (62) 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:660[A]:VAL:O	1:A:664[A]:MET:HG2	1.29	1.26
1:A:171[B]:GLU:CD	1:A:341[B]:LYS:HE3	1.87	0.95
1:A:89[B]:ASN:HD21	1:A:378:ARG:HH21	1.03	0.95
1:A:171[B]:GLU:OE2	1:A:341[B]:LYS:CE	2.17	0.93
1:A:658[A]:PRO:HB3	9:A:2128:HOH:O	1.74	0.87
1:A:171[B]:GLU:OE2	1:A:341[B]:LYS:HE3	1.72	0.87
1:A:89[B]:ASN:ND2	1:A:378:ARG:HH21	1.75	0.84
1:A:733:TYR:HE2	9:A:2147:HOH:O	1.60	0.83
1:A:242:TYR:OH	9:A:2416:HOH:O	2.01	0.78
1:A:400:ARG:O	1:A:404[B]:THR:HG23	1.83	0.77
1:A:540:ASN:OD1	1:A:542:GLU:HB3	1.85	0.75
1:A:171[B]:GLU:OE2	1:A:341[B]:LYS:NZ	2.19	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89[B]:ASN:HD21	1:A:378:ARG:NH2	1.83	0.75
1:A:548:GLY:HA3	9:A:2387:HOH:O	1.87	0.74
1:A:188:LEU:HD11	1:A:329:VAL:HG11	1.74	0.70
1:A:603[B]:VAL:HG13	1:A:607:TYR:CE2	2.27	0.70
1:A:90[B]:PHE:HE2	1:A:94:LYS:HZ3	1.40	0.69
1:A:660[A]:VAL:O	1:A:664[A]:MET:CG	2.25	0.65
1:A:603[B]:VAL:HG13	1:A:607:TYR:CZ	2.31	0.65
1:A:506:PHE:HB2	1:A:509:MET:HE2	1.80	0.64
1:A:612:TYR:CZ	1:A:616:MET:HG3	2.39	0.58
1:A:603[B]:VAL:CG1	1:A:607:TYR:CE2	2.87	0.57
1:A:591[A]:ILE:HG23	1:A:661[A]:LEU:HD21	1.87	0.57
1:A:369:ASP:OD2	1:A:662[B]:ARG:NH2	2.28	0.55
1:A:666:ASP:OD2	9:A:1986:HOH:O	2.18	0.55
1:A:188:LEU:HD11	1:A:329:VAL:CG1	2.38	0.52
1:A:517:SER:HB2	1:A:694:PRO:HG3	1.90	0.52
1:A:661[A]:LEU:O	1:A:664[A]:MET:HB2	2.10	0.52
1:A:303:GLN:HE21	1:A:325:VAL:HG13	1.75	0.51
2:A:1756:NAG:H81	9:A:2019:HOH:O	2.12	0.50
1:A:90[B]:PHE:CE2	1:A:94:LYS:NZ	2.77	0.50
1:A:258:ILE:HD13	1:A:294:VAL:HB	1.92	0.50
1:A:171[B]:GLU:CG	1:A:341[B]:LYS:HE3	2.42	0.50
1:A:246:TRP:CD1	2:A:1760:NAG:H83	2.48	0.49
1:A:653:PHE:HZ	1:A:661[A]:LEU:HD13	1.78	0.48
1:A:310:GLY:O	1:A:330:GLY:HA2	2.14	0.47
1:A:733:TYR:CE2	9:A:2147:HOH:O	2.47	0.47
1:A:239[A]:VAL:HG22	1:A:247:ASN:ND2	2.29	0.46
1:A:303:GLN:NE2	1:A:325:VAL:HG13	2.31	0.46
1:A:412:PRO:HA	1:A:589[B]:ASN:OD1	2.15	0.46
1:A:90[B]:PHE:CZ	1:A:94:LYS:HD3	2.52	0.44
1:A:347:HIS:HE1	9:A:2112:HOH:O	1.99	0.44
1:A:540:ASN:C	1:A:542:GLU:H	2.21	0.44
1:A:589[A]:ASN:OD1	9:A:2417:HOH:O	2.17	0.44
1:A:551:LEU:HD22	1:A:556:TYR:HB2	2.00	0.43
2:A:1761:NAG:H81	9:A:2092:HOH:O	2.18	0.43
1:A:688[A]:ARG:CZ	9:A:1878:HOH:O	2.66	0.43
2:A:1761:NAG:H83	9:A:2020:HOH:O	2.18	0.43
1:A:236:ALA:O	1:A:239[A]:VAL:HG13	2.19	0.42
1:A:132:ASN:HD21	1:A:136:ASN:HD22	1.66	0.42
1:A:347:HIS:CE1	9:A:2112:HOH:O	2.72	0.42
1:A:325:VAL:HB	1:A:326:PRO:HD2	2.01	0.42
1:A:169:MET:HA	1:A:344:MET:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:468:PRO:HG2	1:A:603[A]:VAL:HG21	2.01	0.41
1:A:303:GLN:HE21	1:A:325:VAL:CG1	2.34	0.41
1:A:387:ASP:HA	1:A:388:PRO:HA	1.93	0.41
1:A:239[A]:VAL:HG22	1:A:247:ASN:CG	2.41	0.41
1:A:551:LEU:HA	1:A:551:LEU:HD23	1.88	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:A:1986:HOH:O	9:A:2331:HOH:O[2_565]	1.93	0.27
9:A:2034:HOH:O	9:A:2161:HOH:O[2_565]	1.95	0.25
1:A:656[B]:SER:OG	9:A:2128:HOH:O[4_566]	2.06	0.14
1:A:656[B]:SER:O	9:A:2128:HOH:O[4_566]	2.11	0.09

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	741/709 (104%)	718 (97%)	21 (3%)	2 (0%)	43 19

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	382	VAL
1	A	243	PRO

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	645/605 (107%)	635 (98%)	10 (2%)	65	35

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	LYS
1	A	189	GLU
1	A	443[A]	GLN
1	A	443[B]	GLN
1	A	505	GLU
1	A	519	ASN
1	A	537	TYR
1	A	557	GLU
1	A	600	TYR
1	A	673	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	136	ASN
1	A	303	GLN
1	A	618	HIS
1	A	620	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 19 ligands modelled in this entry, 4 are monoatomic - leaving 15 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
8	YBY	A	1	-	17,26,26	0.92	1 (5%)	21,35,35	2.34	6 (28%)
2	NAG	A	1755	1,2	14,14,15	0.56	0	17,19,21	1.44	1 (5%)
2	NAG	A	1756	2	14,14,15	0.61	0	17,19,21	0.86	0
2	NAG	A	1757	1	14,14,15	0.63	0	17,19,21	1.36	2 (11%)
2	NAG	A	1758	1,2	14,14,15	0.74	0	17,19,21	1.17	1 (5%)
2	NAG	A	1759	1	14,14,15	0.49	0	17,19,21	1.43	1 (5%)
2	NAG	A	1760	1	14,14,15	0.61	0	17,19,21	1.58	4 (23%)
2	NAG	A	1761	1,2	14,14,15	0.91	0	17,19,21	1.53	2 (11%)
2	NAG	A	1762	3,2	14,14,15	0.48	0	17,19,21	1.08	2 (11%)
2	NAG	A	1763	1,2	14,14,15	0.82	0	17,19,21	1.43	2 (11%)
2	NAG	A	1764	3,2	14,14,15	0.63	0	17,19,21	1.37	4 (23%)
3	BMA	A	1765	2,4	11,11,12	0.63	0	15,15,17	0.88	1 (6%)
4	MAN	A	1766	3	11,11,12	0.65	0	15,15,17	1.16	2 (13%)
2	NAG	A	1767	2	14,14,15	0.44	0	17,19,21	1.21	1 (5%)
3	BMA	A	1768	2	11,11,12	0.60	0	15,15,17	0.92	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
8	YBY	A	1	-	-	0/15/25/25	0/1/1/1
2	NAG	A	1755	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	1756	2	-	0/6/23/26	0/1/1/1
2	NAG	A	1757	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1758	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	1759	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	A	1760	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1761	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	1762	3,2	-	0/6/23/26	0/1/1/1
2	NAG	A	1763	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	1764	3,2	-	0/6/23/26	0/1/1/1
3	BMA	A	1765	2,4	-	0/2/19/22	0/1/1/1
4	MAN	A	1766	3	-	0/2/19/22	0/1/1/1
2	NAG	A	1767	2	-	0/6/23/26	0/1/1/1
3	BMA	A	1768	2	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	1	YBY	C5-C6	2.25	1.56	1.53

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	1	YBY	C5-C6-C7	-6.37	102.64	112.11
8	A	1	YBY	C7-C6-N10	-4.47	102.70	112.39
2	A	1755	NAG	O5-C1-C2	-4.22	105.69	111.52
2	A	1761	NAG	O5-C1-C2	-3.99	106.01	111.52
8	A	1	YBY	C14-C13-N12	-3.91	101.23	108.90
2	A	1763	NAG	O5-C1-C2	-3.81	106.27	111.52
2	A	1761	NAG	O5-C5-C4	-3.18	103.10	110.83
2	A	1758	NAG	O5-C1-C2	-2.93	107.47	111.52
8	A	1	YBY	O26-C11-N10	-2.72	117.65	122.56
2	A	1760	NAG	O3-C3-C4	-2.63	104.20	110.34
2	A	1762	NAG	O7-C7-C8	-2.35	117.83	122.07
3	A	1765	BMA	O3-C3-C2	-2.32	105.73	110.04
2	A	1760	NAG	O7-C7-C8	-2.30	117.91	122.07
2	A	1764	NAG	O7-C7-C8	-2.20	118.09	122.07
2	A	1764	NAG	C3-C4-C5	-2.16	106.37	110.24
2	A	1763	NAG	C6-C5-C4	-2.08	108.09	112.99
2	A	1760	NAG	C2-N2-C7	-2.06	119.94	122.94
2	A	1764	NAG	O5-C5-C6	2.02	110.34	107.15
4	A	1766	MAN	C1-C2-C3	2.12	112.34	109.66
2	A	1757	NAG	C4-C3-C2	2.17	114.20	111.02
4	A	1766	MAN	O5-C5-C6	2.26	110.73	107.15
2	A	1764	NAG	C8-C7-N2	2.41	120.32	116.10
8	A	1	YBY	O26-C11-N12	2.48	127.04	122.56
2	A	1762	NAG	C8-C7-N2	2.56	120.58	116.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1757	NAG	O5-C5-C6	2.74	111.48	107.15
2	A	1767	NAG	O5-C5-C6	2.85	111.66	107.15
8	A	1	YBY	C19-C17-I18	3.11	123.02	119.80
2	A	1760	NAG	C1-O5-C5	3.13	116.50	112.19
2	A	1759	NAG	O5-C5-C6	4.36	114.05	107.15

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1756	NAG	1	0
2	A	1760	NAG	1	0
2	A	1761	NAG	2	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	691/709 (97%)	0.03	39 (5%) 24 27	14, 23, 42, 67	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	541	TRP	6.8
1	A	55	LYS	5.8
1	A	543	THR	5.6
1	A	542	GLU	4.7
1	A	201[A]	VAL	4.2
1	A	153	ASN	4.2
1	A	719	VAL	4.1
1	A	506	PHE	4.0
1	A	544	ASN	4.0
1	A	335	GLY	3.8
1	A	507	SER	3.8
1	A	336	ASN	3.7
1	A	134	ASP	3.6
1	A	155[A]	SER	3.5
1	A	174[A]	LEU	3.4
1	A	56	HIS	3.2
1	A	652	ASP	3.2
1	A	135	GLY	3.1
1	A	136	ASN	3.0
1	A	226	ILE	3.0
1	A	505	GLU	3.0
1	A	123	THR	2.9
1	A	133	GLU	2.8
1	A	227	LEU	2.7
1	A	339	THR	2.6
1	A	656[A]	SER	2.5
1	A	324	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	486	GLU	2.4
1	A	284	ALA	2.4
1	A	122	LYS	2.3
1	A	225	VAL	2.3
1	A	409	GLY	2.3
1	A	487	GLY	2.3
1	A	651[A]	GLN	2.2
1	A	219	LEU	2.2
1	A	297	ILE	2.2
1	A	124	HIS	2.1
1	A	154	VAL	2.0
1	A	334	THR	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. 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(\AA^2)	Q<0.9
2	NAG	A	1756	14/15	0.66	0.25	43,50,53,56	0
2	NAG	A	1757	14/15	0.69	0.24	43,49,53,53	0
3	BMA	A	1768	11/12	0.72	0.29	55,57,61,63	0
2	NAG	A	1767	14/15	0.79	0.31	54,57,62,63	0
2	NAG	A	1759	14/15	0.85	0.21	59,64,67,67	0
2	NAG	A	1758	14/15	0.86	0.13	37,40,43,49	0
2	NAG	A	1762	14/15	0.86	0.17	35,42,49,50	0
2	NAG	A	1760	14/15	0.89	0.12	28,40,46,48	0
2	NAG	A	1761	14/15	0.90	0.09	27,31,35,39	0
3	BMA	A	1765	11/12	0.90	0.17	32,34,36,37	0
2	NAG	A	1764	14/15	0.90	0.20	33,36,44,47	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NAG	A	1763	14/15	0.90	0.10	21,26,33,43	0
4	MAN	A	1766	11/12	0.91	0.19	38,41,45,48	0
8	YBY	A	1	26/26	0.92	0.10	16,20,38,48	0
2	NAG	A	1755	14/15	0.93	0.09	32,39,43,47	0
5	ZN	A	1752	1/1	1.00	0.04	19,19,19,19	0
7	CL	A	1754	1/1	1.00	0.04	22,22,22,22	0
5	ZN	A	1751	1/1	1.00	0.05	18,18,18,18	0
6	CA	A	1753	1/1	1.00	0.05	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.