



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:15 PM GMT

PDB ID : 4OC5
Title : X-ray structure of of human glutamate carboxypeptidase II (GCPII) in a complex with CHIBzL, a urea-based inhibitor N²-{[(S)-carboxy(4-hydroxyphenyl)methyl]carbamoyl}-N⁶-(4-iodobenzoyl)-L-lysine
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Deposited on : 2014-01-08
Resolution : 1.70 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 (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

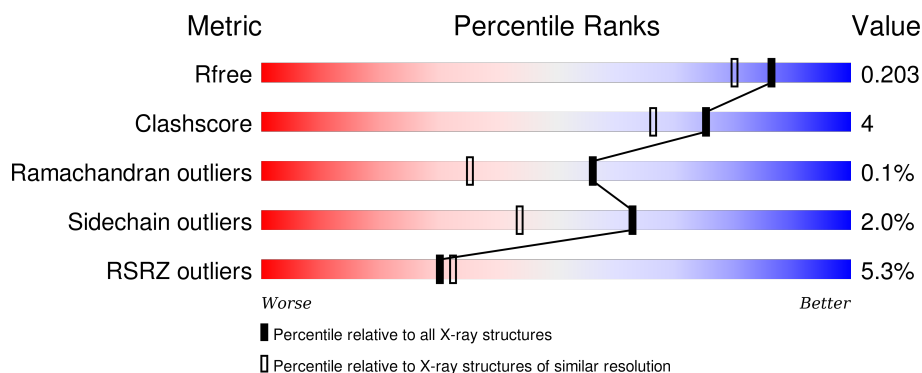
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	3190 (1.70-1.70)
Clashscore	102246	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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	

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	NAG	A	802	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	NAG	A	810	-	-	-	X

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 6466 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
1	A	694	Total	C	N	O	S	0	44	0
			5754	3700	962	1069	23			

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 a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	2	Total	C	N	O	0	0
			28	16	2	10		
2	A	2	Total	C	N	O	0	0
			28	16	2	10		
2	A	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 3 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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		

- Molecule 4 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	4	Total	C	N	O	0	0
			50	28	2	20		

- 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		

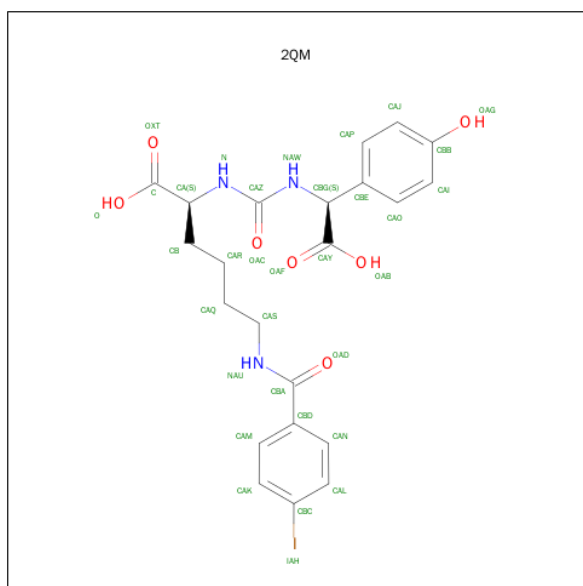
- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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 1 1	0	0

- Molecule 8 is N 2 -{[(S)-CARBOXY(4-HYDROXYPHENYL)METHYL]CARBAMOYL}-N 6 -(4-IODOBENZOYL)-L-LYSINE (three-letter code: 2QM) (formula: C₂₂H₂₄IN₃O₇).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C I N O 33 22 1 3 7	0	0

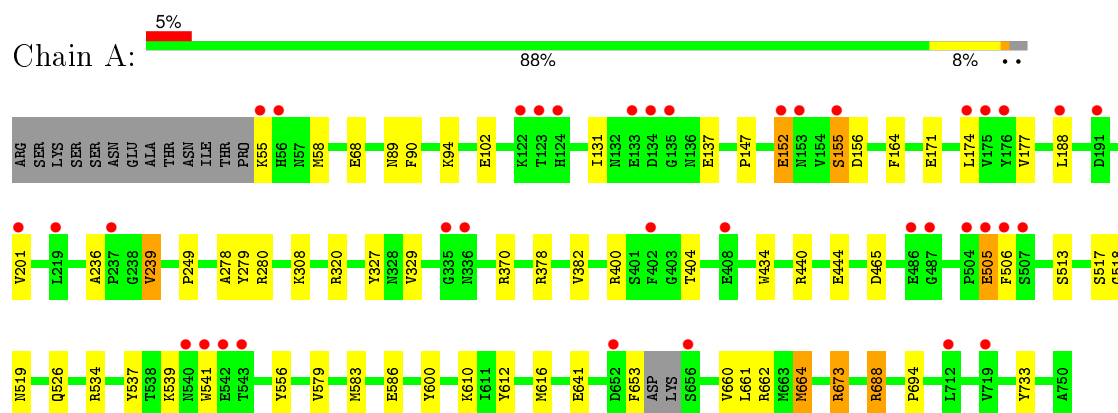
- Molecule 9 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	499	Total O 499 499	0	0

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 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: 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	101.58Å 130.45Å 159.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.70 20.00 – 1.70	Depositor EDS
% Data completeness (in resolution range)	97.7 (20.00-1.70) 97.7 (20.00-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.47 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.169 , 0.204 0.169 , 0.203	Depositor DCC
R_{free} test set	1139 reflections (1.02%)	DCC
Wilson B-factor (Å ²)	25.3	Xtriage
Anisotropy	0.067	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 45.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 113015 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6466	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.06% 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, 0.375 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: ZN, BMA, NAG, CL, CA, 2QM, 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	0.95	2/6022 (0.0%)	0.86	9/8154 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	434	TRP	CE3-CZ3	5.36	1.47	1.38
1	A	164	PHE	CE2-CZ	5.26	1.47	1.37

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	688[A]	ARG	NE-CZ-NH2	-6.63	116.99	120.30
1	A	688[B]	ARG	NE-CZ-NH2	-6.63	116.99	120.30
1	A	673	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	A	673	ARG	NE-CZ-NH2	-6.15	117.22	120.30
1	A	174[A]	LEU	CA-CB-CG	5.75	128.52	115.30
1	A	174[B]	LEU	CA-CB-CG	5.75	128.52	115.30
1	A	664	MET	CG-SD-CE	-5.25	91.79	100.20
1	A	370	ARG	NE-CZ-NH1	5.24	122.92	120.30
1	A	440	ARG	NE-CZ-NH2	-5.04	117.78	120.30

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	5754	0	5633	45	0
2	A	84	0	75	1	0
3	A	42	0	39	0	0
4	A	50	0	43	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	33	0	21	1	0
9	A	499	0	0	8	2
All	All	6466	0	5811	46	2

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 4.

All (46) 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:MET:HG2	1.38	1.22
1:A:147:PRO:HG2	1:A:152:GLU:HG2	1.46	0.97
1:A:89[B]:ASN:HD21	1:A:378:ARG:HH21	1.08	0.97
1:A:89[B]:ASN:ND2	1:A:378:ARG:HH21	1.79	0.78
1:A:90:PHE:CE2	1:A:94:LYS:HE2	2.20	0.76
1:A:733:TYR:HE2	9:A:1244:HOH:O	1.70	0.74
1:A:58[A]:MET:CE	1:A:586:GLU:HG2	2.18	0.72
1:A:131:ILE:HG22	1:A:137:GLU:HG2	1.72	0.72
1:A:641:GLU:HG3	9:A:1253:HOH:O	1.91	0.69
1:A:58[A]:MET:HE1	1:A:586:GLU:HG2	1.74	0.68
1:A:89[B]:ASN:HD21	1:A:378:ARG:NH2	1.87	0.66
1:A:400:ARG:O	1:A:404[B]:THR:HG23	1.96	0.65
1:A:539:LYS:HD2	1:A:541:TRP:HB3	1.80	0.63
1:A:278:ALA:HB3	1:A:280[A]:ARG:NH1	2.15	0.62
1:A:308:LYS:HB2	9:A:1397:HOH:O	2.01	0.60
1:A:58[A]:MET:HE2	1:A:586:GLU:HG2	1.86	0.57
1:A:177:VAL:HG12	1:A:188[A]:LEU:HD11	1.89	0.55
1:A:188[B]:LEU:HD11	1:A:329:VAL:HG11	1.89	0.54
1:A:465:ASP:OD1	1:A:513:SER:HB2	2.08	0.53
1:A:579:VAL:O	1:A:583[B]:MET:HG2	2.08	0.53
1:A:155:SER:O	1:A:156:ASP:HB2	2.13	0.49
1:A:58[A]:MET:HE1	1:A:586:GLU:CG	2.45	0.45
1:A:177:VAL:CG2	1:A:201[B]:VAL:HG13	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:VAL:HG22	1:A:201[B]:VAL:HG13	1.98	0.45
1:A:444:GLU:HA	1:A:444:GLU:OE1	2.17	0.44
1:A:249:PRO:HG2	1:A:556:TYR:CE2	2.53	0.43
1:A:612:TYR:CZ	1:A:616:MET:HG3	2.53	0.43
1:A:610:LYS:HE2	9:A:1080:HOH:O	2.18	0.43
1:A:131:ILE:CG2	1:A:137:GLU:HG2	2.43	0.43
1:A:517:SER:HB2	1:A:694:PRO:HG3	2.00	0.43
1:A:505:GLU:HG3	1:A:506:PHE:CE1	2.54	0.42
1:A:90:PHE:CZ	1:A:94:LYS:HE2	2.54	0.42
1:A:236:ALA:O	1:A:239[A]:VAL:HG13	2.19	0.42
1:A:539:LYS:CD	1:A:541:TRP:HB3	2.49	0.41
1:A:534:ARG:HG3	8:A:818:2QM:IAH	2.91	0.41
1:A:733:TYR:CE2	9:A:1244:HOH:O	2.55	0.41
1:A:526[A]:GLN:NE2	9:A:1374:HOH:O	2.37	0.41
2:A:802:NAG:H81	9:A:1144:HOH:O	2.21	0.41
1:A:320[A]:ARG:HG2	1:A:327:TYR:HB2	2.02	0.41
1:A:688[B]:ARG:NH1	9:A:970:HOH:O	2.32	0.40
1:A:653:PHE:HE1	1:A:661[A]:LEU:HB2	1.86	0.40
1:A:517:SER:OG	1:A:518:GLY:N	2.52	0.40

All (2) 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:1115:HOH:O	9:A:1364:HOH:O[2_565]	1.97	0.23
9:A:1255:HOH:O	9:A:1399:HOH:O[2_565]	2.10	0.10

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	731/709 (103%)	714 (98%)	16 (2%)	1 (0%)	56 35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	382	VAL

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	633/605 (105%)	618 (98%)	15 (2%)	57	36

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	LYS
1	A	68[A]	GLU
1	A	68[B]	GLU
1	A	152	GLU
1	A	155	SER
1	A	171	GLU
1	A	239[A]	VAL
1	A	239[B]	VAL
1	A	505	GLU
1	A	519	ASN
1	A	537	TYR
1	A	600	TYR
1	A	662[A]	ARG
1	A	662[B]	ARG
1	A	673	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	HIS
1	A	124	HIS
1	A	347	HIS
1	A	443	GLN

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Mol	Chain	Res	Type
1	A	618	HIS

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 ⓘ

10 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$
2	NAG	A	801	1,2	14,14,15	0.50	0	15,19,21	1.02	1 (6%)
2	NAG	A	802	2	14,14,15	0.71	0	15,19,21	1.06	1 (6%)
2	NAG	A	804	1,2	14,14,15	0.79	0	15,19,21	0.78	0
2	NAG	A	805	2	14,14,15	0.41	0	15,19,21	1.19	1 (6%)
2	NAG	A	808	1,2	14,14,15	0.93	1 (7%)	15,19,21	1.24	1 (6%)
2	NAG	A	809	2	14,14,15	0.52	0	15,19,21	1.27	3 (20%)
4	NAG	A	810	1,4	14,14,15	0.49	0	15,19,21	1.03	1 (6%)
4	NAG	A	811	4	14,14,15	0.69	0	15,19,21	1.47	2 (13%)
4	BMA	A	812	4	11,11,12	0.47	0	14,15,17	0.93	1 (7%)
4	MAN	A	813	4	11,11,12	0.55	0	14,15,17	1.04	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
2	NAG	A	801	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	802	2	-	0/6/23/26	0/1/1/1
2	NAG	A	804	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	805	2	-	0/6/23/26	0/1/1/1
2	NAG	A	808	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	809	2	-	0/6/23/26	0/1/1/1
4	NAG	A	810	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	811	4	-	0/6/23/26	0/1/1/1
4	BMA	A	812	4	-	0/2/19/22	0/1/1/1
4	MAN	A	813	4	-	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)
2	A	808	NAG	O7-C7	2.46	1.28	1.23

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	811	NAG	C3-C4-C5	-3.01	104.96	110.20
2	A	808	NAG	O4-C4-C3	-2.87	103.87	110.34
2	A	801	NAG	O3-C3-C4	-2.35	105.05	110.34
4	A	812	BMA	O3-C3-C2	-2.34	105.76	110.00
2	A	809	NAG	O7-C7-C8	-2.34	117.77	122.06
4	A	810	NAG	C6-C5-C4	-2.15	107.72	113.02
4	A	813	MAN	O5-C5-C6	2.07	111.82	107.35
4	A	811	NAG	C8-C7-N2	2.08	120.08	116.11
2	A	809	NAG	C8-C7-N2	2.25	120.41	116.11
2	A	805	NAG	O5-C5-C6	2.33	112.38	107.35
2	A	809	NAG	C1-O5-C5	2.33	115.21	112.25
2	A	802	NAG	C2-N2-C7	2.53	126.30	123.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	802	NAG	1	0

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	803	1	14,14,15	0.67	0	15,19,21	1.20	1 (6%)
3	NAG	A	806	1	14,14,15	0.48	0	15,19,21	1.16	2 (13%)
3	NAG	A	807	1	14,14,15	0.64	0	15,19,21	1.50	4 (26%)
8	2QM	A	818	5	28,34,34	1.31	4 (14%)	35,45,45	1.84	10 (28%)

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	803	1	-	0/6/23/26	0/1/1/1
3	NAG	A	806	1	-	0/6/23/26	0/1/1/1
3	NAG	A	807	1	-	0/6/23/26	0/1/1/1
8	2QM	A	818	5	-	0/24/32/32	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	818	2QM	CAK-CAM	2.31	1.42	1.38
8	A	818	2QM	CAL-CAN	2.55	1.43	1.38
8	A	818	2QM	CBC-IAH	3.26	2.18	2.10
8	A	818	2QM	CBD-CBA	3.39	1.57	1.50

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	818	2QM	CAL-CBC-CAK	-5.38	112.39	120.66
8	A	818	2QM	CAY-CBG-CBE	-3.10	108.48	112.97
8	A	818	2QM	CAO-CBE-CBG	-2.86	115.90	120.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	803	NAG	O3-C3-C2	-2.76	103.64	109.11
8	A	818	2QM	CAP-CAJ-CBB	-2.43	117.06	119.87
3	A	807	NAG	O7-C7-C8	-2.41	117.63	122.06
8	A	818	2QM	CAN-CBD-CAM	-2.37	115.08	118.60
3	A	807	NAG	C3-C4-C5	-2.12	106.49	110.20
8	A	818	2QM	CAI-CAO-CBE	-2.00	119.15	121.20
3	A	806	NAG	O5-C5-C6	2.01	111.69	107.35
8	A	818	2QM	CAL-CBC-IAH	2.30	123.28	119.67
3	A	807	NAG	C1-O5-C5	2.65	115.62	112.25
3	A	807	NAG	O5-C5-C6	2.77	113.33	107.35
3	A	806	NAG	C1-O5-C5	2.79	115.80	112.25
8	A	818	2QM	CAN-CAL-CBC	2.86	123.77	119.53
8	A	818	2QM	CAK-CBC-IAH	2.96	124.32	119.67
8	A	818	2QM	CAM-CAK-CBC	3.32	124.46	119.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	818	2QM	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	694/709 (97%)	0.03	37 (5%)	30 32	17, 28, 48, 65	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	152	GLU	6.1
1	A	543	THR	5.0
1	A	541	TRP	4.6
1	A	719	VAL	4.5
1	A	133	GLU	4.1
1	A	134	ASP	4.1
1	A	155	SER	4.1
1	A	174[A]	LEU	4.0
1	A	335	GLY	3.8
1	A	505	GLU	3.5
1	A	506	PHE	3.3
1	A	135	GLY	3.3
1	A	55	LYS	3.2
1	A	507	SER	3.1
1	A	153	ASN	3.1
1	A	237	PRO	3.1
1	A	652	ASP	3.1
1	A	191	ASP	2.8
1	A	123	THR	2.7
1	A	540	ASN	2.6
1	A	201[A]	VAL	2.5
1	A	504	PRO	2.5
1	A	408	GLU	2.5
1	A	336	ASN	2.4
1	A	219	LEU	2.4
1	A	712	LEU	2.4
1	A	176	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	487	GLY	2.3
1	A	188[A]	LEU	2.2
1	A	402	PHE	2.2
1	A	56	HIS	2.2
1	A	124	HIS	2.1
1	A	122	LYS	2.1
1	A	175	VAL	2.1
1	A	486[A]	GLU	2.1
1	A	656[A]	SER	2.0
1	A	542	GLU	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 ⓘ

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
2	NAG	A	802	14/15	0.71	0.28	4.76	45,53,57,58	0
4	NAG	A	810	14/15	0.91	0.10	2.20	25,29,39,45	0
4	MAN	A	813	11/12	0.93	0.18	1.49	50,52,56,57	0
2	NAG	A	808	14/15	0.93	0.13	-	31,34,42,44	0
4	BMA	A	812	11/12	0.93	0.14	-	42,44,45,47	0
4	NAG	A	811	14/15	0.88	0.27	-	39,43,52,52	0
2	NAG	A	805	14/15	0.77	0.28	-	60,62,67,69	0
2	NAG	A	801	14/15	0.92	0.10	-	35,43,48,52	0
2	NAG	A	804	14/15	0.89	0.14	-	40,44,48,54	0
2	NAG	A	809	14/15	0.91	0.21	-	36,43,49,51	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, 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
8	2QM	A	818	33/33	0.98	0.09	0.31	22,29,38,45	0
6	CA	A	816	1/1	0.99	0.07	-0.64	18,18,18,18	0
7	CL	A	817	1/1	0.98	0.03	-3.03	29,29,29,29	0
5	ZN	A	814	1/1	1.00	0.03	-3.06	23,23,23,23	0
5	ZN	A	815	1/1	1.00	0.01	-6.79	23,23,23,23	0
3	NAG	A	807	14/15	0.84	0.16	-	33,49,55,59	0
3	NAG	A	806	14/15	0.90	0.23	-	60,68,70,71	0
3	NAG	A	803	14/15	0.71	0.26	-	46,53,56,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.