



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:54 AM UTC

PDB ID : 8E6G / pdb_00008e6g
Title : Co-crystal structure of Chaetomium glucosidase with compound 10
Authors : Karade, S.S.; Mariuzza, R.A.
Deposited on : 2022-08-22
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

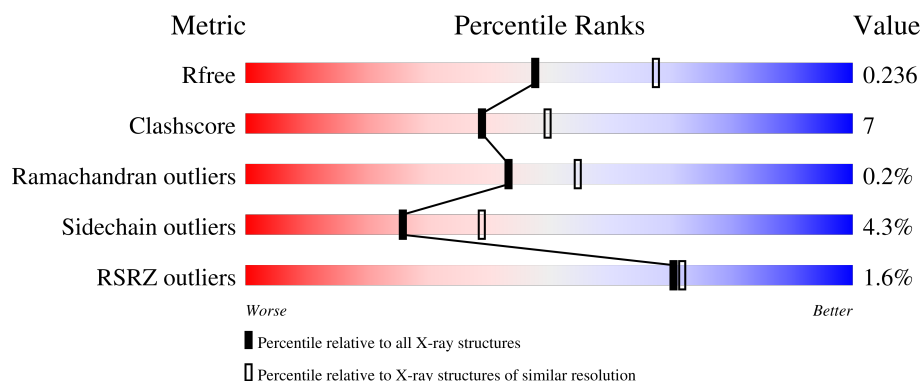
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

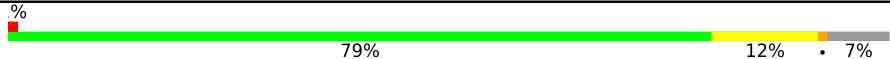
The reported resolution of this entry is 2.30 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	819	 79% 12% • 7%
1	B	819	 78% 14% • 7%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 12815 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 Chaetomium alpha glucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	763	Total	C	N	O	S	0	1	0
			6057	3893	1021	1130	13			
1	B	764	Total	C	N	O	S	0	0	0
			6079	3901	1021	1144	13			

There are 78 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP G0SFD1
A	0	GLY	-	expression tag	UNP G0SFD1
A	1	ILE	-	expression tag	UNP G0SFD1
A	2	LEU	-	expression tag	UNP G0SFD1
A	3	PRO	-	expression tag	UNP G0SFD1
A	4	SER	-	expression tag	UNP G0SFD1
A	5	PRO	-	expression tag	UNP G0SFD1
A	6	GLY	-	expression tag	UNP G0SFD1
A	7	MET	-	expression tag	UNP G0SFD1
A	8	PRO	-	expression tag	UNP G0SFD1
A	9	ALA	-	expression tag	UNP G0SFD1
A	10	LEU	-	expression tag	UNP G0SFD1
A	11	LEU	-	expression tag	UNP G0SFD1
A	12	SER	-	expression tag	UNP G0SFD1
A	13	LEU	-	expression tag	UNP G0SFD1
A	14	VAL	-	expression tag	UNP G0SFD1
A	15	SER	-	expression tag	UNP G0SFD1
A	16	LEU	-	expression tag	UNP G0SFD1
A	17	LEU	-	expression tag	UNP G0SFD1
A	18	SER	-	expression tag	UNP G0SFD1
A	19	VAL	-	expression tag	UNP G0SFD1
A	20	LEU	-	expression tag	UNP G0SFD1
A	21	LEU	-	expression tag	UNP G0SFD1
A	22	MET	-	expression tag	UNP G0SFD1
A	23	GLY	-	expression tag	UNP G0SFD1

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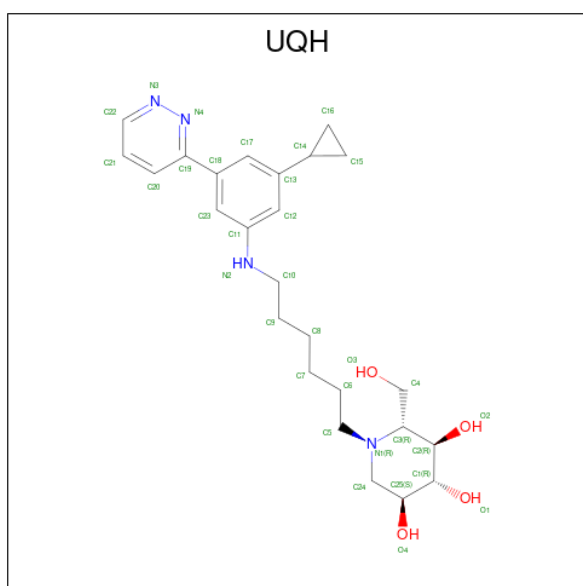
Chain	Residue	Modelled	Actual	Comment	Reference
A	24	CYS	-	expression tag	UNP G0SFD1
A	25	VAL	-	expression tag	UNP G0SFD1
A	26	ALA	-	expression tag	UNP G0SFD1
A	27	GLU	-	expression tag	UNP G0SFD1
A	28	THR	-	expression tag	UNP G0SFD1
A	29	GLY	-	expression tag	UNP G0SFD1
A	810	SER	-	expression tag	UNP G0SFD1
A	811	GLY	-	expression tag	UNP G0SFD1
A	812	HIS	-	expression tag	UNP G0SFD1
A	813	HIS	-	expression tag	UNP G0SFD1
A	814	HIS	-	expression tag	UNP G0SFD1
A	815	HIS	-	expression tag	UNP G0SFD1
A	816	HIS	-	expression tag	UNP G0SFD1
A	817	HIS	-	expression tag	UNP G0SFD1
B	-1	MET	-	initiating methionine	UNP G0SFD1
B	0	GLY	-	expression tag	UNP G0SFD1
B	1	ILE	-	expression tag	UNP G0SFD1
B	2	LEU	-	expression tag	UNP G0SFD1
B	3	PRO	-	expression tag	UNP G0SFD1
B	4	SER	-	expression tag	UNP G0SFD1
B	5	PRO	-	expression tag	UNP G0SFD1
B	6	GLY	-	expression tag	UNP G0SFD1
B	7	MET	-	expression tag	UNP G0SFD1
B	8	PRO	-	expression tag	UNP G0SFD1
B	9	ALA	-	expression tag	UNP G0SFD1
B	10	LEU	-	expression tag	UNP G0SFD1
B	11	LEU	-	expression tag	UNP G0SFD1
B	12	SER	-	expression tag	UNP G0SFD1
B	13	LEU	-	expression tag	UNP G0SFD1
B	14	VAL	-	expression tag	UNP G0SFD1
B	15	SER	-	expression tag	UNP G0SFD1
B	16	LEU	-	expression tag	UNP G0SFD1
B	17	LEU	-	expression tag	UNP G0SFD1
B	18	SER	-	expression tag	UNP G0SFD1
B	19	VAL	-	expression tag	UNP G0SFD1
B	20	LEU	-	expression tag	UNP G0SFD1
B	21	LEU	-	expression tag	UNP G0SFD1
B	22	MET	-	expression tag	UNP G0SFD1
B	23	GLY	-	expression tag	UNP G0SFD1
B	24	CYS	-	expression tag	UNP G0SFD1
B	25	VAL	-	expression tag	UNP G0SFD1
B	26	ALA	-	expression tag	UNP G0SFD1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	27	GLU	-	expression tag	UNP G0SFD1
B	28	THR	-	expression tag	UNP G0SFD1
B	29	GLY	-	expression tag	UNP G0SFD1
B	810	SER	-	expression tag	UNP G0SFD1
B	811	GLY	-	expression tag	UNP G0SFD1
B	812	HIS	-	expression tag	UNP G0SFD1
B	813	HIS	-	expression tag	UNP G0SFD1
B	814	HIS	-	expression tag	UNP G0SFD1
B	815	HIS	-	expression tag	UNP G0SFD1
B	816	HIS	-	expression tag	UNP G0SFD1
B	817	HIS	-	expression tag	UNP G0SFD1

- Molecule 2 is (2R,3R,4R,5S)-1-(6-{[(5M)-3-cyclopropyl-5-(pyridazin-3-yl)phenyl]amino}hexyl)-2-(hydroxymethyl)piperidine-3,4,5-triol (CCD ID: UQH) (formula: C₂₅H₃₆N₄O₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			33	25	4	4		
2	B	1	Total	C	N	O	0	1
			66	50	8	8		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



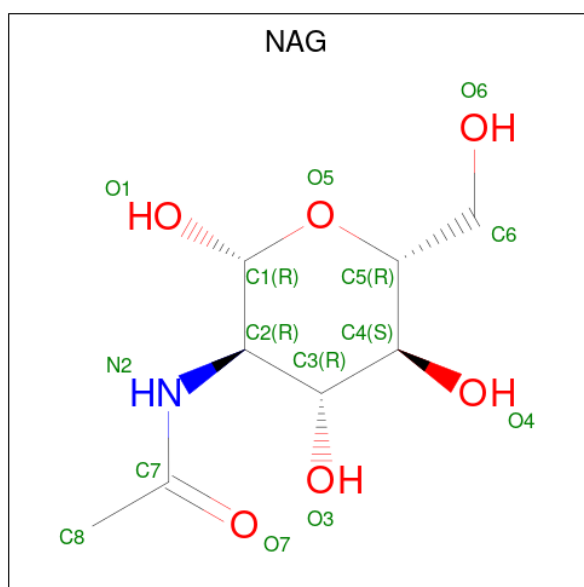
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		

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

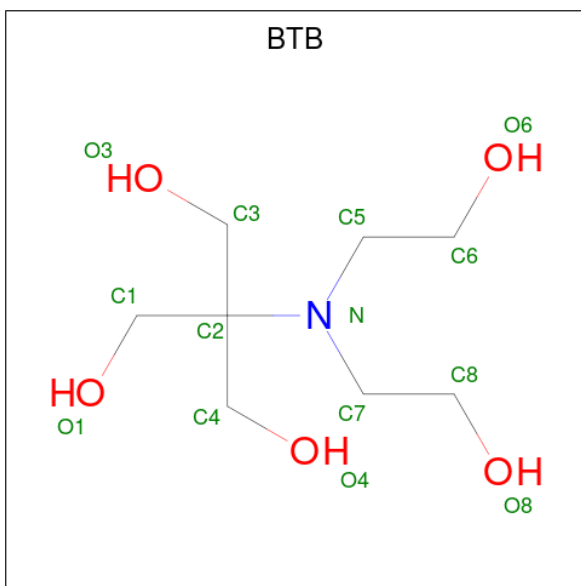
- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 6 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN

E-1,3-DIOL (CCD ID: BTB) (formula: C₈H₁₉NO₅).

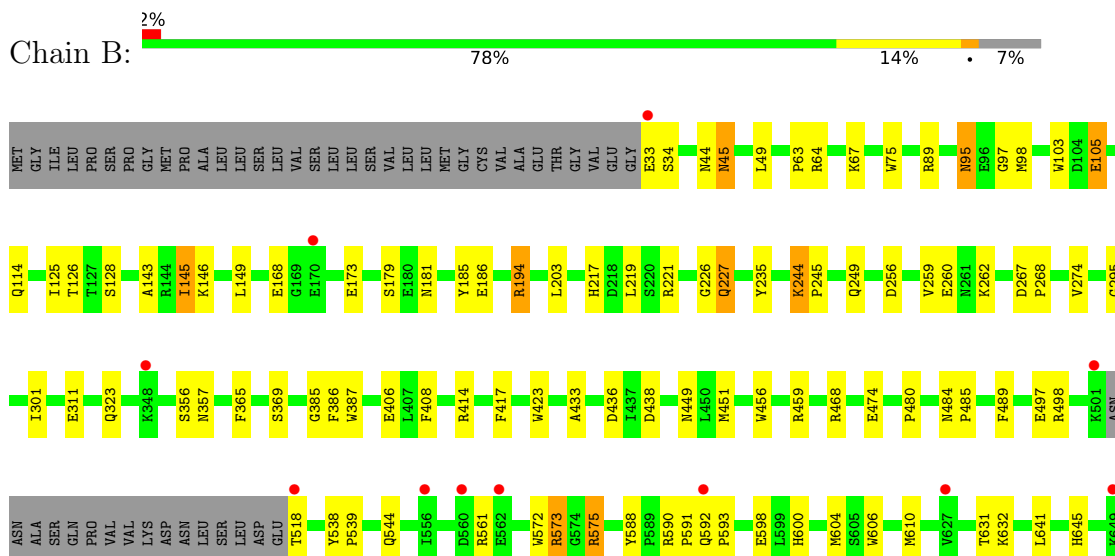


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	237	Total	O	0	0
			237	237		
7	B	247	Total	O	0	0
			247	247		

- Molecule 1: Chaetomium alpha glucosidase





4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	135.83Å 178.80Å 179.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.81 – 2.30 46.81 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.1 (46.81-2.30) 98.4 (46.81-2.30)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.192 , 0.236 0.192 , 0.236	Depositor DCC
R_{free} test set	4847 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	35.7	Xtriage
Anisotropy	0.439	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 25.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.011 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12815	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% 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: NAG, UQH, GOL, BTB, 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	0.59	0/6235	1.17	24/8487 (0.3%)
1	B	0.60	0/6255	1.13	13/8513 (0.2%)
All	All	0.60	0/12490	1.15	37/17000 (0.2%)

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	459	ARG	CG-CD-NE	-9.66	90.75	112.00
1	A	575	ARG	CB-CA-C	7.59	122.06	109.53
1	B	45	ASN	CB-CA-C	7.18	122.71	110.79
1	B	575	ARG	CB-CA-C	7.08	123.85	109.76
1	A	657	THR	CB-CA-C	-7.07	96.86	114.11
1	B	459	ARG	CG-CD-NE	-7.05	96.50	112.00
1	A	795	THR	CA-CB-OG1	-6.54	99.80	109.60
1	A	573	ARG	CG-CD-NE	-6.49	97.73	112.00
1	A	91	THR	CA-CB-OG1	-6.43	99.95	109.60
1	A	212	ILE	CB-CA-C	6.39	118.13	110.78
1	A	795	THR	CB-CA-C	6.30	120.18	109.53
1	A	459	ARG	CB-CG-CD	6.20	125.55	111.30
1	B	776	GLU	CB-CG-CD	6.17	123.10	112.60
1	A	484	ASN	CA-CB-CG	5.89	118.50	112.60
1	A	459	ARG	NE-CZ-NH2	-5.89	113.90	119.20
1	A	270	PRO	N-CA-CB	5.83	106.45	103.19
1	A	459	ARG	CD-NE-CZ	5.79	132.50	124.40
1	B	657	THR	CB-CA-C	-5.79	98.96	111.11
1	B	468	ARG	CG-CD-NE	-5.73	99.39	112.00
1	A	382	GLU	CB-CG-CD	5.64	122.18	112.60
1	A	802	THR	CA-CB-OG1	-5.60	101.20	109.60
1	B	631	THR	CB-CA-C	5.47	119.46	110.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	436	ASP	CA-CB-CG	5.45	118.05	112.60
1	A	748	GLN	CB-CA-C	-5.45	98.90	109.68
1	A	658	ILE	CA-C-N	5.42	130.60	120.95
1	A	658	ILE	C-N-CA	5.42	130.60	120.95
1	A	339	LYS	CB-CA-C	5.37	118.66	109.80
1	B	573	ARG	CB-CG-CD	5.36	123.62	111.30
1	B	793	GLN	N-CA-CB	-5.35	103.27	111.56
1	A	484	ASN	O-C-N	-5.21	117.77	121.83
1	B	518	THR	CB-CA-C	5.21	120.56	109.10
1	B	684	LYS	CB-CA-C	5.19	116.90	109.11
1	A	158	LYS	CB-CA-C	5.14	118.52	110.19
1	A	285	ASN	CB-CA-C	5.08	118.56	110.09
1	B	260	GLU	CB-CG-CD	5.05	121.19	112.60
1	A	40	ILE	CA-C-N	5.05	125.70	119.99
1	A	40	ILE	C-N-CA	5.05	125.70	119.99

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	6057	0	5738	74	0
1	B	6079	0	5739	87	0
2	A	33	0	0	0	0
2	B	66	0	0	0	0
3	A	12	0	16	0	0
3	B	6	0	8	0	0
4	A	30	0	0	1	0
4	B	20	0	0	1	0
5	B	14	0	13	0	0
6	B	14	0	19	1	0
7	A	237	0	0	4	0
7	B	247	0	0	4	0
All	All	12815	0	11533	161	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 7.

All (161) 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:451:MET:HE1	1:A:540:LEU:HB3	1.28	1.11
1:A:658:ILE:HD12	1:A:662:GLU:HA	1.31	1.08
1:A:657:THR:HG21	7:A:1097:HOH:O	1.54	1.06
1:B:572:TRP:H	1:B:600:HIS:HD2	1.07	0.99
1:A:451:MET:CE	1:A:540:LEU:HB3	1.95	0.95
1:B:561:ARG:HE	1:B:664:HIS:HD2	1.02	0.93
1:A:434:ASP:OD2	1:A:807:LYS:HD3	1.75	0.86
1:B:572:TRP:H	1:B:600:HIS:CD2	1.92	0.86
1:B:561:ARG:HE	1:B:664:HIS:CD2	1.94	0.81
6:B:903:BTB:H62	7:B:1150:HOH:O	1.83	0.77
1:A:451:MET:HE1	1:A:540:LEU:CB	2.10	0.77
1:B:114:GLN:NE2	1:B:414:ARG:HH12	1.83	0.77
1:A:658:ILE:CD1	1:A:662:GLU:HA	2.14	0.76
1:A:572:TRP:H	1:A:600:HIS:HD2	1.36	0.73
1:B:194:ARG:CG	1:B:194:ARG:HH11	2.05	0.69
1:A:659:ASP:HB3	1:A:661:PHE:H	1.57	0.69
1:B:766:ASN:O	1:B:770:THR:HG23	1.93	0.69
1:B:217:HIS:HE1	1:B:267:ASP:O	1.75	0.69
1:B:451:MET:HE1	1:B:544:GLN:HB2	1.76	0.68
1:B:786:ASN:ND2	1:B:788:GLU:H	1.95	0.64
1:B:451:MET:CE	1:B:544:GLN:HB2	2.29	0.63
1:B:727:ARG:O	1:B:728:SER:HB3	1.98	0.63
1:B:179:SER:HB3	1:B:185:TYR:CD2	2.34	0.62
1:A:685:PRO:HG3	1:A:748:GLN:HE22	1.64	0.62
1:A:783:GLU:OE1	1:A:795:THR:HB	1.99	0.62
1:B:489:PHE:CE1	1:B:610:MET:HE3	2.35	0.62
1:B:786:ASN:HD22	1:B:789:THR:H	1.48	0.62
1:B:217:HIS:HD2	1:B:219:LEU:H	1.48	0.61
1:A:569:ALA:HA	1:A:604:MET:CE	2.30	0.61
1:A:727:ARG:O	1:A:728:SER:HB3	2.01	0.60
1:B:561:ARG:NE	1:B:664:HIS:HD2	1.86	0.60
1:B:103:TRP:H	1:B:357:ASN:ND2	1.99	0.60
1:B:386:PHE:HB3	7:B:1141:HOH:O	2.02	0.60
1:B:433:ALA:O	1:B:498:ARG:NH2	2.35	0.58
1:A:578:SER:N	4:A:908:SO4:O1	2.30	0.58
1:B:194:ARG:HH11	1:B:194:ARG:HG3	1.68	0.58
1:A:244:LYS:HB3	1:A:245:PRO:HD3	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:423:TRP:H	1:A:484:ASN:ND2	2.03	0.57
1:A:456:TRP:CE2	1:A:480:PRO:HA	2.40	0.56
1:A:489:PHE:CE1	1:A:610:MET:HE3	2.40	0.56
1:B:181:ASN:HD22	1:B:186:GLU:HG2	1.70	0.56
1:B:572:TRP:N	1:B:600:HIS:HD2	1.91	0.56
1:B:672:TYR:CG	1:B:734:ILE:HG21	2.41	0.56
1:B:103:TRP:H	1:B:357:ASN:HD21	1.54	0.56
1:B:728:SER:N	1:B:729:PRO:CD	2.69	0.56
1:B:179:SER:HB3	1:B:185:TYR:CE2	2.41	0.56
1:A:598:GLU:OE2	1:A:600:HIS:HE1	1.88	0.55
7:A:1059:HOH:O	1:B:262:LYS:HD2	2.05	0.55
1:B:728:SER:N	1:B:729:PRO:HD3	2.20	0.55
1:B:707:TYR:HA	1:B:770:THR:HG21	1.89	0.55
1:B:235:TYR:OH	1:B:249:GLN:NE2	2.39	0.55
1:B:786:ASN:HB2	1:B:793:GLN:NE2	2.23	0.54
1:B:385:GLY:HA2	1:B:387:TRP:CZ3	2.43	0.54
1:A:658:ILE:HD11	1:A:662:GLU:OE1	2.08	0.53
1:A:575:ARG:HH22	1:A:592:GLN:HE22	1.57	0.53
1:A:203:LEU:HD11	1:A:301:ILE:CG2	2.39	0.53
1:A:313:THR:O	1:A:316:ASP:HB2	2.08	0.53
1:A:234[B]:THR:HG21	1:B:268:PRO:HG3	1.89	0.53
1:B:194:ARG:HG3	1:B:194:ARG:NH1	2.22	0.53
1:B:114:GLN:HE22	1:B:414:ARG:HH22	1.57	0.52
1:B:194:ARG:CG	1:B:194:ARG:NH1	2.70	0.52
1:B:786:ASN:HD21	1:B:788:GLU:HB2	1.75	0.52
1:B:386:PHE:CB	7:B:1141:HOH:O	2.58	0.52
1:A:545:PHE:CZ	1:A:549:ARG:HD2	2.45	0.51
1:A:635:ASP:O	1:A:639:HIS:HD2	1.93	0.51
1:B:766:ASN:O	1:B:770:THR:CG2	2.58	0.51
1:B:659:ASP:HB3	1:B:665:LYS:HE2	1.91	0.51
1:B:114:GLN:HE21	1:B:414:ARG:HH12	1.54	0.51
1:A:126:THR:HG22	7:A:1186:HOH:O	2.09	0.51
1:A:451:MET:HE3	1:A:540:LEU:HD22	1.92	0.51
1:B:217:HIS:CD2	1:B:219:LEU:H	2.28	0.51
1:B:95:ASN:HD22	1:B:95:ASN:C	2.18	0.50
1:B:149:LEU:HD11	1:B:295:GLY:HA2	1.93	0.50
1:B:456:TRP:CE2	1:B:480:PRO:HA	2.47	0.49
1:B:728:SER:H	1:B:729:PRO:CD	2.25	0.49
1:A:570:TYR:HB3	1:A:603:LEU:HG	1.94	0.49
1:A:128:SER:O	1:A:143:ALA:HA	2.13	0.49
1:A:515:LEU:O	1:A:519:LEU:HB2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:423:TRP:H	1:A:484:ASN:HD22	1.58	0.49
1:A:456:TRP:CD2	1:A:480:PRO:HA	2.48	0.49
1:B:226:GLY:H	1:B:227:GLN:NE2	2.11	0.48
1:B:149:LEU:CD1	1:B:295:GLY:HA2	2.44	0.48
1:B:95:ASN:HD22	1:B:97:GLY:H	1.62	0.48
1:A:733:ASN:HB3	1:A:801:TRP:CG	2.48	0.47
1:B:145:ILE:CD1	1:B:301:ILE:HD12	2.44	0.47
1:B:385:GLY:HA2	1:B:387:TRP:CH2	2.49	0.47
1:B:762:ARG:HD2	4:B:907:SO4:O3	2.14	0.47
1:B:672:TYR:CE1	1:B:711:SER:HA	2.50	0.47
1:B:538:TYR:HB3	1:B:539:PRO:HD3	1.96	0.47
1:B:217:HIS:CE1	1:B:267:ASP:O	2.62	0.47
1:B:256:ASP:O	1:B:259:VAL:HG22	2.15	0.47
1:A:786:ASN:ND2	1:A:788:GLU:H	2.13	0.46
1:A:720:GLY:HA2	1:A:724:ASN:OD1	2.15	0.46
1:A:672:TYR:CZ	1:A:711:SER:HA	2.50	0.46
1:A:426:GLY:HA3	1:A:487:THR:OG1	2.15	0.46
1:A:748:GLN:HE21	1:A:748:GLN:HB3	1.62	0.46
1:A:234[A]:THR:CG2	1:A:284:GLY:HA2	2.46	0.46
1:B:125:ILE:HA	1:B:146:LYS:O	2.15	0.46
1:B:591:PRO:HD2	1:B:598:GLU:OE2	2.15	0.46
1:A:234[A]:THR:HG22	1:A:285:ASN:H	1.80	0.45
1:A:786:ASN:HD22	1:A:788:GLU:H	1.63	0.45
1:A:441:LEU:HD22	1:A:533:TYR:CE1	2.51	0.45
1:B:707:TYR:CA	1:B:770:THR:HG21	2.46	0.45
1:B:67:LYS:HG2	1:B:168:GLU:CD	2.42	0.45
1:B:497:GLU:OE1	1:B:497:GLU:HA	2.16	0.45
1:B:64:ARG:HD3	1:B:408:PHE:CE1	2.52	0.45
1:A:672:TYR:CE1	1:A:711:SER:HA	2.52	0.45
1:B:128:SER:O	1:B:143:ALA:HA	2.17	0.45
1:B:226:GLY:H	1:B:227:GLN:HE22	1.64	0.45
1:A:594:PRO:HA	1:A:598:GLU:OE1	2.17	0.45
1:A:724:ASN:HA	1:A:727:ARG:HB2	1.99	0.45
1:A:451:MET:HE1	1:A:540:LEU:CA	2.47	0.44
1:A:728:SER:N	1:A:729:PRO:CD	2.80	0.44
1:A:110:ARG:HD2	1:A:110:ARG:HA	1.78	0.44
1:A:601:VAL:HG21	1:A:653:TYR:HB3	1.99	0.44
1:B:244:LYS:HB3	1:B:245:PRO:HD3	1.99	0.44
1:B:727:ARG:O	1:B:728:SER:CB	2.63	0.44
1:A:727:ARG:O	1:A:728:SER:CB	2.66	0.44
1:B:456:TRP:CD2	1:B:480:PRO:HA	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:672:TYR:CE2	1:B:711:SER:HB3	2.52	0.44
1:A:372:ASP:HA	1:A:399:GLU:HA	2.00	0.44
1:A:49:LEU:HA	7:A:1093:HOH:O	2.18	0.43
1:B:485:PRO:HB3	1:B:606:TRP:CE2	2.53	0.43
1:B:604:MET:HE3	1:B:641:LEU:HG	1.99	0.43
1:A:355:PHE:HB2	1:A:806:VAL:HG21	2.01	0.43
1:B:591:PRO:HG2	1:B:598:GLU:HG2	2.00	0.43
1:A:234[B]:THR:HA	1:A:285:ASN:OD1	2.19	0.43
1:A:672:TYR:CG	1:A:734:ILE:HG21	2.54	0.43
1:B:575:ARG:HE	1:B:575:ARG:HB2	1.65	0.43
1:A:646:TRP:HE1	1:A:648:GLU:HG2	1.84	0.43
1:A:575:ARG:HH22	1:A:592:GLN:NE2	2.16	0.43
1:A:110:ARG:NH2	1:A:323:GLN:HG3	2.33	0.42
1:B:798:PHE:C	1:B:800:GLY:HA3	2.44	0.42
1:B:75:TRP:CZ2	1:B:89:ARG:HG3	2.55	0.42
1:B:203:LEU:HD11	1:B:301:ILE:CG2	2.50	0.42
1:A:598:GLU:OE2	1:A:600:HIS:CE1	2.70	0.42
1:A:203:LEU:HD11	1:A:301:ILE:HG23	2.01	0.42
1:A:481:HIS:H	1:A:481:HIS:CD2	2.37	0.42
1:B:145:ILE:HD13	1:B:301:ILE:CD1	2.50	0.42
1:B:588:TYR:O	1:B:590:ARG:HG2	2.20	0.42
1:A:723:GLU:O	1:A:794:ARG:NH1	2.52	0.41
1:B:592:GLN:HA	1:B:593:PRO:HA	1.92	0.41
1:A:308:ALA:O	1:A:309:GLY:O	2.38	0.41
1:A:526:ASN:HB2	1:A:529:VAL:CG2	2.50	0.41
1:B:423:TRP:H	1:B:484:ASN:ND2	2.19	0.41
1:A:195:SER:O	1:A:199:GLY:N	2.49	0.41
1:A:659:ASP:CB	1:A:661:PHE:H	2.29	0.41
1:B:105:GLU:HA	1:B:356:SER:OG	2.20	0.41
1:A:719:TYR:CE1	1:A:728:SER:HB2	2.55	0.41
1:B:604:MET:HE2	1:B:645:HIS:HD2	1.86	0.41
1:A:603:LEU:HD12	1:A:603:LEU:O	2.20	0.41
1:B:386:PHE:HA	7:B:1141:HOH:O	2.20	0.41
1:B:662:GLU:OE2	1:B:662:GLU:HA	2.21	0.41
1:A:719:TYR:CE1	1:A:728:SER:CB	3.04	0.41
1:B:49:LEU:O	1:B:63:PRO:HA	2.21	0.41
1:A:499:LEU:HD12	1:A:499:LEU:HA	1.90	0.40
1:A:428:HIS:O	1:A:431:PRO:HD2	2.21	0.40
1:B:365:PHE:O	1:B:406:GLU:HA	2.21	0.40
1:A:114:GLN:HB3	1:A:127:THR:OG1	2.21	0.40
1:A:234[A]:THR:HA	1:A:285:ASN:OD1	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:181:ASN:HD22	1:B:186:GLU:CG	2.32	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	760/819 (93%)	736 (97%)	22 (3%)	2 (0%)	36	46
1	B	760/819 (93%)	744 (98%)	15 (2%)	1 (0%)	48	60
All	All	1520/1638 (93%)	1480 (97%)	37 (2%)	3 (0%)	43	55

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	728	SER
1	B	728	SER
1	A	309	GLY

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	627/707 (89%)	605 (96%)	22 (4%)	32	48
1	B	632/707 (89%)	600 (95%)	32 (5%)	21	32

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	1259/1414 (89%)	1205 (96%)	54 (4%)	26	39

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

Mol	Chain	Res	Type
1	A	44	ASN
1	A	45	ASN
1	A	49	LEU
1	A	212	ILE
1	A	218	ASP
1	A	280	LYS
1	A	322	LYS
1	A	343	GLN
1	A	348	LYS
1	A	417	PHE
1	A	438	ASP
1	A	526	ASN
1	A	550	LYS
1	A	558	SER
1	A	573	ARG
1	A	611	VAL
1	A	616	SER
1	A	635	ASP
1	A	657	THR
1	A	658	ILE
1	A	786	ASN
1	A	795	THR
1	B	33	GLU
1	B	34	SER
1	B	44	ASN
1	B	45	ASN
1	B	95	ASN
1	B	98	MET
1	B	105	GLU
1	B	126	THR
1	B	145	ILE
1	B	173	GLU
1	B	194	ARG
1	B	221	ARG
1	B	227	GLN
1	B	244	LYS
1	B	274	VAL

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Mol	Chain	Res	Type
1	B	311	GLU
1	B	323	GLN
1	B	369	SER
1	B	417	PHE
1	B	438	ASP
1	B	449	ASN
1	B	474	GLU
1	B	573	ARG
1	B	632	LYS
1	B	657	THR
1	B	662	GLU
1	B	676	PHE
1	B	692	LYS
1	B	770	THR
1	B	777	GLU
1	B	788	GLU
1	B	805	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	120	GLN
1	A	157	GLN
1	A	167	GLN
1	A	200	ASN
1	A	242	GLN
1	A	249	GLN
1	A	323	GLN
1	A	368	HIS
1	A	397	HIS
1	A	398	GLN
1	A	461	GLN
1	A	478	GLN
1	A	484	ASN
1	A	592	GLN
1	A	600	HIS
1	A	639	HIS
1	A	748	GLN
1	A	786	ASN
1	A	793	GLN
1	B	95	ASN
1	B	114	GLN

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Mol	Chain	Res	Type
1	B	120	GLN
1	B	137	HIS
1	B	167	GLN
1	B	181	ASN
1	B	217	HIS
1	B	227	GLN
1	B	249	GLN
1	B	250	GLN
1	B	323	GLN
1	B	357	ASN
1	B	368	HIS
1	B	398	GLN
1	B	461	GLN
1	B	481	HIS
1	B	484	ASN
1	B	600	HIS
1	B	664	HIS
1	B	748	GLN
1	B	786	ASN
1	B	793	GLN
1	B	812	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

18 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	A	903	-	5,5,5	0.11	0	5,5,5	0.41	0
5	NAG	B	902	1	14,14,15	0.73	0	17,19,21	2.53	5 (29%)
4	SO4	A	909	-	4,4,4	0.31	0	6,6,6	0.11	0
4	SO4	B	907	-	4,4,4	0.32	0	6,6,6	0.14	0
4	SO4	B	908	-	4,4,4	0.26	0	6,6,6	0.10	0
3	GOL	B	904	-	5,5,5	0.47	0	5,5,5	1.00	0
4	SO4	A	905	-	4,4,4	0.30	0	6,6,6	0.10	0
4	SO4	A	904	-	4,4,4	0.35	0	6,6,6	0.06	0
2	UQH	B	901[A]	-	36,36,36	3.09	10 (27%)	48,49,49	4.44	5 (10%)
4	SO4	B	906	-	4,4,4	0.32	0	6,6,6	0.06	0
2	UQH	B	901[B]	-	36,36,36	3.09	10 (27%)	48,49,49	4.44	5 (10%)
4	SO4	A	908	-	4,4,4	0.29	0	6,6,6	0.14	0
3	GOL	A	902	-	5,5,5	0.17	0	5,5,5	0.66	0
4	SO4	A	906	-	4,4,4	0.39	0	6,6,6	0.35	0
6	BTB	B	903	-	13,13,13	1.39	3 (23%)	7,16,16	0.92	0
4	SO4	A	907	-	4,4,4	0.28	0	6,6,6	0.15	0
2	UQH	A	901	-	36,36,36	3.09	10 (27%)	48,49,49	4.44	5 (10%)
4	SO4	B	905	-	4,4,4	0.39	0	6,6,6	0.27	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	903	-	-	2/4/4/4	-
5	NAG	B	902	1	-	2/6/23/26	0/1/1/1
3	GOL	B	904	-	-	0/4/4/4	-
2	UQH	B	901[A]	-	-	8/20/42/42	0/4/4/4
2	UQH	B	901[B]	-	-	2/20/42/42	0/4/4/4
3	GOL	A	902	-	-	4/4/4/4	-
6	BTB	B	903	-	-	7/21/21/21	-
2	UQH	A	901	-	-	7/20/42/42	0/4/4/4

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	UQH	C5-N1	-15.22	1.20	1.47
2	B	901[A]	UQH	C5-N1	-15.21	1.20	1.47
2	B	901[B]	UQH	C5-N1	-15.19	1.20	1.47
2	B	901[B]	UQH	C25-C1	-4.57	1.45	1.52
2	A	901	UQH	C25-C1	-4.53	1.45	1.52
2	B	901[A]	UQH	C25-C1	-4.46	1.45	1.52
2	B	901[B]	UQH	C18-C19	3.67	1.54	1.49
2	B	901[A]	UQH	C18-C19	3.66	1.54	1.49
2	A	901	UQH	C18-C19	3.63	1.54	1.49
6	B	903	BTB	C2-N	3.42	1.55	1.48
2	B	901[B]	UQH	C15-C14	-3.33	1.39	1.49
2	B	901[A]	UQH	C15-C14	-3.33	1.39	1.49
2	A	901	UQH	C16-C14	-3.33	1.39	1.49
2	B	901[A]	UQH	C16-C14	-3.32	1.39	1.49
2	B	901[B]	UQH	C16-C14	-3.32	1.39	1.49
2	A	901	UQH	C15-C14	-3.31	1.39	1.49
2	B	901[A]	UQH	C11-N2	3.25	1.47	1.38
2	A	901	UQH	C11-N2	3.24	1.47	1.38
2	B	901[B]	UQH	C11-N2	3.21	1.47	1.38
2	B	901[B]	UQH	C16-C15	-2.65	1.39	1.48
2	B	901[B]	UQH	C24-C25	2.64	1.55	1.52
2	B	901[A]	UQH	C24-C25	2.64	1.55	1.52
2	B	901[A]	UQH	C16-C15	-2.64	1.39	1.48
2	A	901	UQH	C16-C15	-2.63	1.39	1.48
2	A	901	UQH	C24-C25	2.60	1.55	1.52
2	B	901[B]	UQH	O1-C1	2.59	1.49	1.43
2	B	901[A]	UQH	O1-C1	2.56	1.49	1.43
2	A	901	UQH	O1-C1	2.55	1.49	1.43
2	B	901[A]	UQH	C21-C20	2.17	1.42	1.38
2	A	901	UQH	C21-C20	2.14	1.42	1.38
2	B	901[B]	UQH	C21-C20	2.13	1.42	1.38
6	B	903	BTB	C5-N	2.12	1.51	1.48
6	B	903	BTB	C7-N	2.03	1.50	1.48

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	901	UQH	C15-C14-C13	21.03	150.02	121.28
2	B	901[B]	UQH	C15-C14-C13	21.02	150.00	121.28
2	B	901[B]	UQH	C16-C14-C13	21.02	150.00	121.28
2	B	901[A]	UQH	C16-C14-C13	21.02	150.00	121.28
2	B	901[A]	UQH	C15-C14-C13	21.01	149.99	121.28
2	A	901	UQH	C16-C14-C13	20.99	149.96	121.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	902	NAG	C2-N2-C7	6.23	131.25	122.90
5	B	902	NAG	C1-O5-C5	5.52	119.59	112.19
2	A	901	UQH	C22-N3-N4	4.01	121.99	118.91
2	B	901[B]	UQH	C22-N3-N4	3.96	121.95	118.91
2	B	901[A]	UQH	C22-N3-N4	3.90	121.90	118.91
5	B	902	NAG	C3-C4-C5	-2.91	104.96	110.23
5	B	902	NAG	C8-C7-N2	2.86	120.86	116.12
2	B	901[B]	UQH	C4-C3-C2	-2.68	108.90	112.93
2	B	901[A]	UQH	C4-C3-C2	-2.67	108.91	112.93
5	B	902	NAG	O7-C7-N2	-2.66	117.29	121.98
2	A	901	UQH	C4-C3-C2	-2.65	108.95	112.93
2	A	901	UQH	C18-C19-N4	2.49	119.59	115.90
2	B	901[B]	UQH	C18-C19-N4	2.46	119.55	115.90
2	B	901[A]	UQH	C18-C19-N4	2.42	119.49	115.90

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	901	UQH	C17-C13-C14-C15
2	A	901	UQH	C12-C13-C14-C16
2	B	901[A]	UQH	C17-C13-C14-C15
2	B	901[A]	UQH	C12-C13-C14-C16
3	A	902	GOL	O1-C1-C2-C3
3	A	902	GOL	C1-C2-C3-O3
3	A	902	GOL	O2-C2-C3-O3
3	A	903	GOL	C1-C2-C3-O3
6	B	903	BTB	N-C2-C3-O3
6	B	903	BTB	C1-C2-C4-O4
6	B	903	BTB	C3-C2-C4-O4
6	B	903	BTB	N-C2-C4-O4
2	A	901	UQH	C23-C11-N2-C10
2	A	901	UQH	C12-C11-N2-C10
5	B	902	NAG	C8-C7-N2-C2
5	B	902	NAG	O7-C7-N2-C2
3	A	903	GOL	O2-C2-C3-O3
2	A	901	UQH	N2-C10-C9-C8
3	A	902	GOL	O1-C1-C2-O2
2	B	901[B]	UQH	N2-C10-C9-C8
6	B	903	BTB	N-C5-C6-O6
6	B	903	BTB	N-C7-C8-O8
2	A	901	UQH	C17-C13-C14-C16

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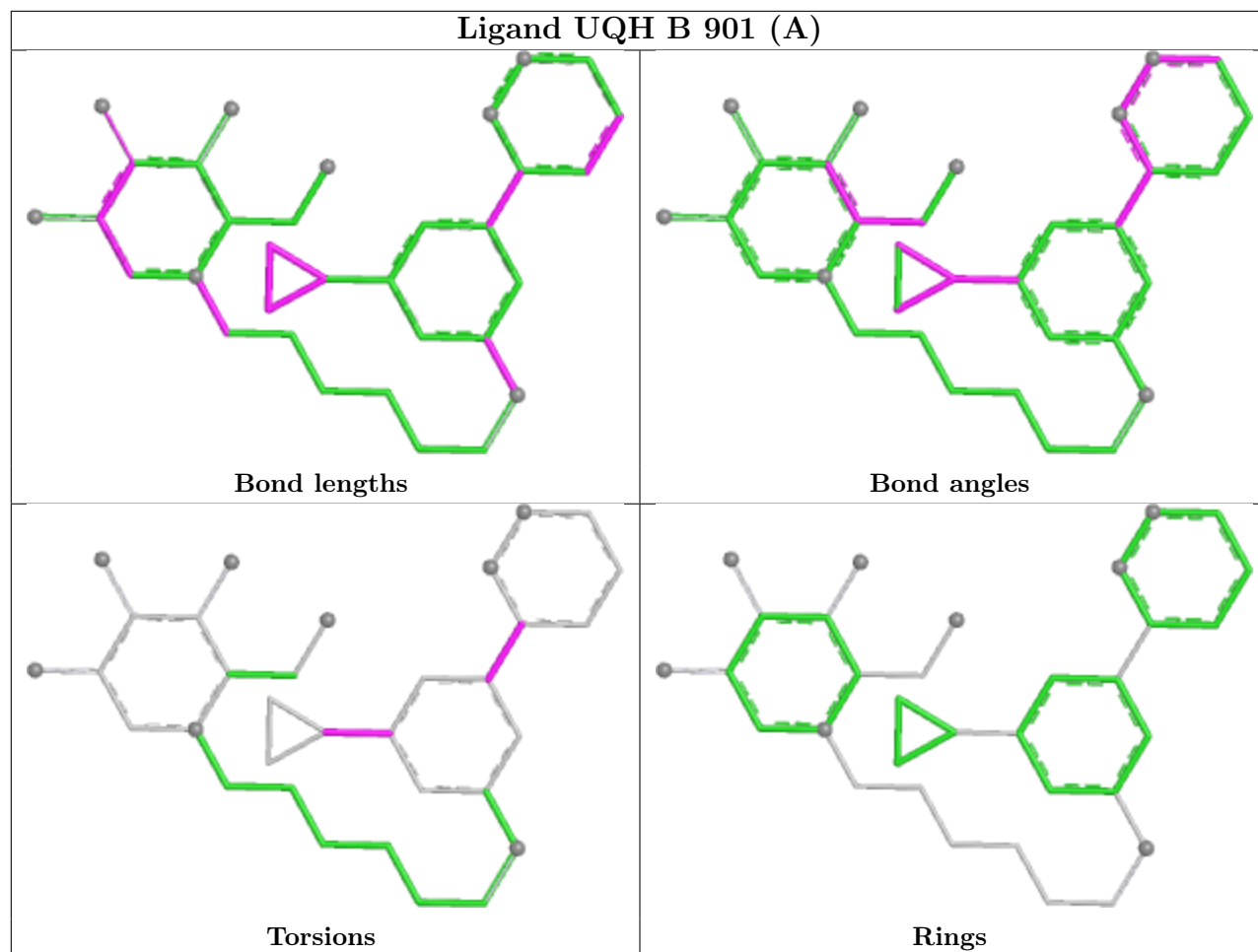
Mol	Chain	Res	Type	Atoms
2	A	901	UQH	C12-C13-C14-C15
2	B	901[A]	UQH	C17-C13-C14-C16
2	B	901[A]	UQH	C12-C13-C14-C15
6	B	903	BTB	C4-C2-C3-O3
2	B	901[A]	UQH	C23-C18-C19-N4
2	B	901[A]	UQH	C17-C18-C19-N4
2	B	901[A]	UQH	C17-C18-C19-C20
2	B	901[A]	UQH	C23-C18-C19-C20
2	B	901[B]	UQH	C7-C8-C9-C10

There are no ring outliers.

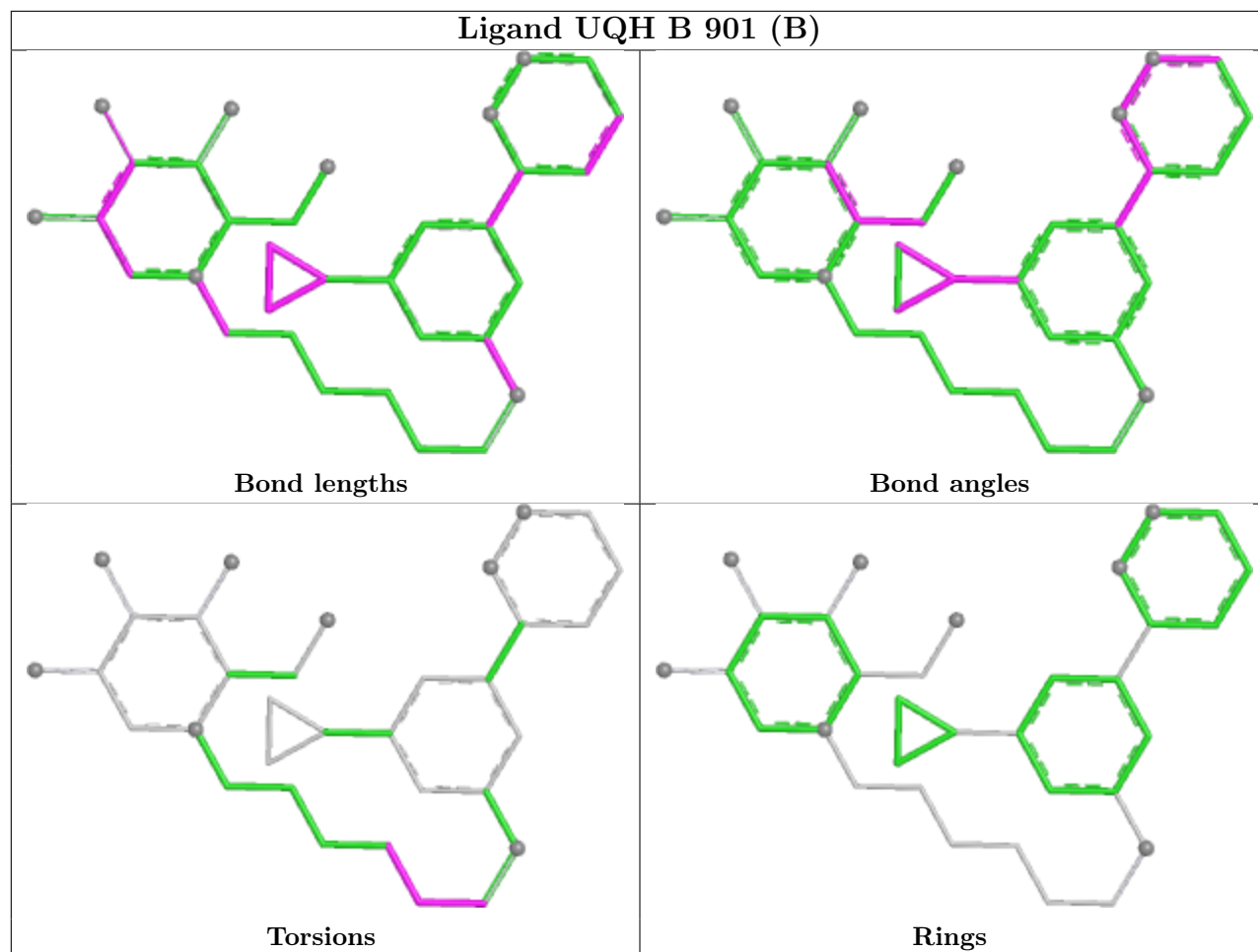
3 monomers are involved in 3 short contacts:

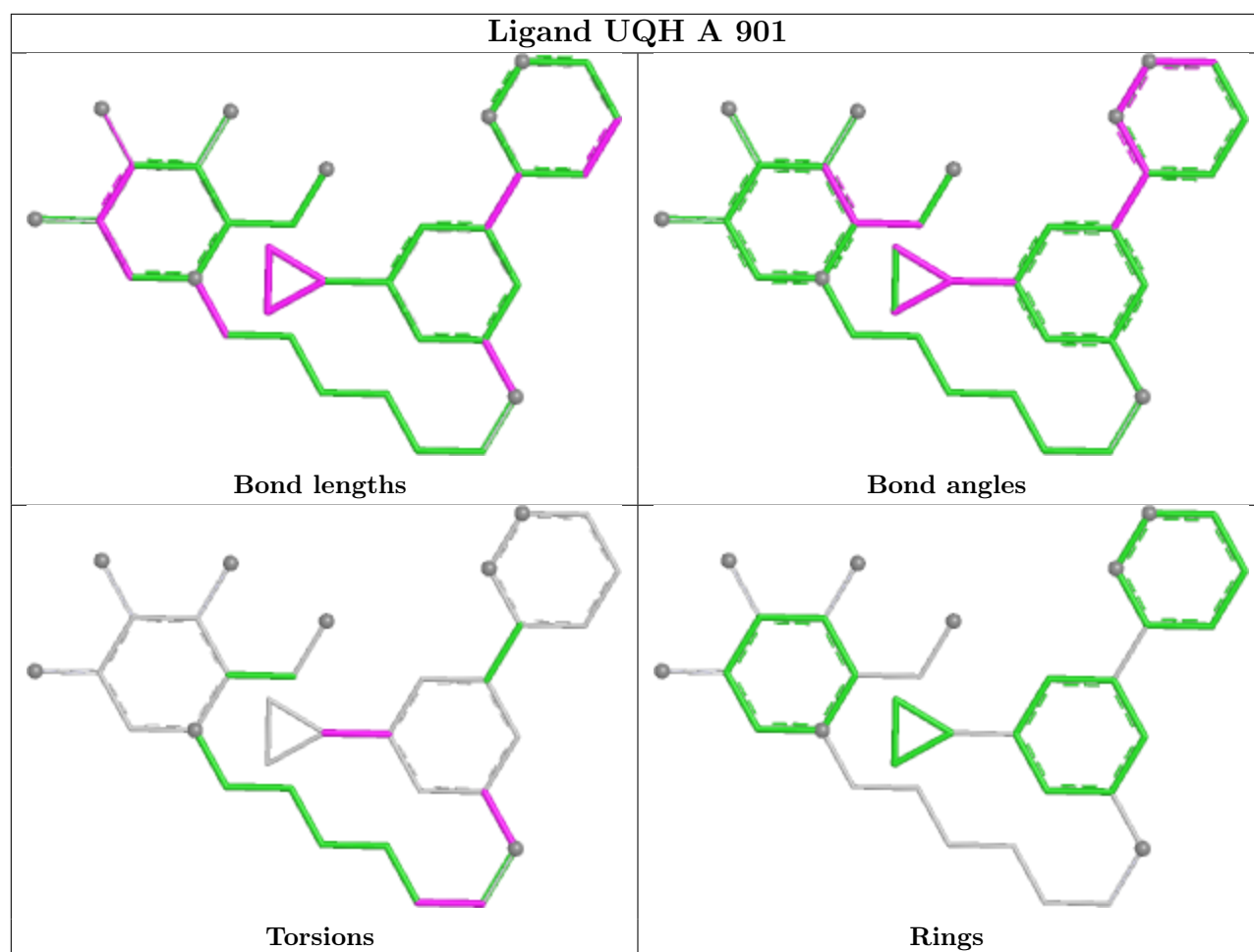
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	907	SO4	1	0
4	A	908	SO4	1	0
6	B	903	BTB	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



Ligand UQH B 901 (B)





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	763/819 (93%)	-0.22	12 (1%) 70 72	18, 35, 60, 103	1 (0%)
1	B	764/819 (93%)	-0.27	13 (1%) 69 70	21, 34, 56, 86	0
All	All	1527/1638 (93%)	-0.24	25 (1%) 70 72	18, 35, 59, 103	1 (0%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	515	LEU	5.6
1	B	518	THR	4.3
1	B	812	HIS	4.2
1	A	560	ASP	3.9
1	B	556	ILE	3.6
1	A	516	ASP	3.5
1	B	33	GLU	3.4
1	B	501	LYS	3.1
1	A	658	ILE	3.1
1	A	517	GLU	3.0
1	A	518	THR	3.0
1	A	812	HIS	2.9
1	B	562	GLU	2.8
1	A	501	LYS	2.6
1	B	592	GLN	2.5
1	B	170	GLU	2.5
1	A	660	GLU	2.4
1	B	663	GLU	2.4
1	B	649	LYS	2.3
1	B	627	VAL	2.3
1	B	560	ASP	2.2
1	A	661	PHE	2.2
1	A	473	LYS	2.1
1	A	42	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	348	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](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

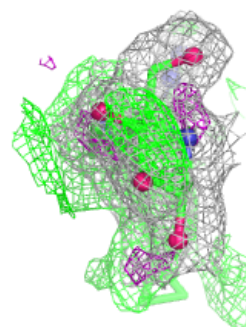
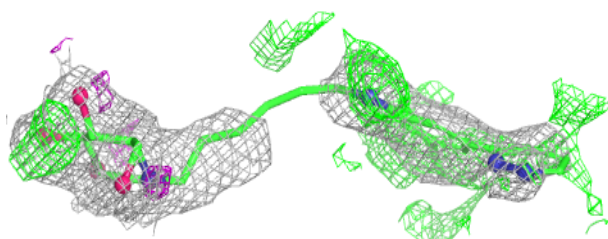
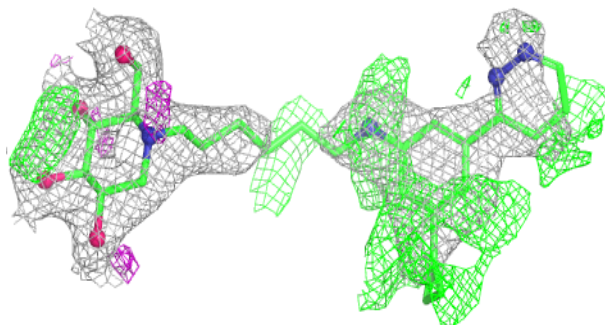
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	SO4	A	909	5/5	0.74	0.12	89,92,95,100	0
4	SO4	A	908	5/5	0.80	0.13	73,77,89,91	0
4	SO4	A	905	5/5	0.82	0.11	96,99,102,104	0
4	SO4	A	904	5/5	0.84	0.11	90,91,94,96	0
5	NAG	B	902	14/15	0.84	0.13	52,65,72,74	0
2	UQH	B	901[B]	33/33	0.85	0.25	22,33,47,49	33
2	UQH	B	901[A]	33/33	0.85	0.25	21,34,56,58	33
6	BTB	B	903	14/14	0.86	0.14	51,63,68,70	0
3	GOL	A	903	6/6	0.87	0.16	50,55,67,69	0
4	SO4	B	906	5/5	0.89	0.10	70,70,80,82	0
2	UQH	A	901	33/33	0.91	0.18	21,52,90,90	0
4	SO4	B	908	5/5	0.92	0.12	51,70,71,73	0
3	GOL	A	902	6/6	0.93	0.11	49,57,61,64	0
3	GOL	B	904	6/6	0.93	0.14	24,41,46,48	0
4	SO4	B	907	5/5	0.95	0.10	53,56,62,68	0
4	SO4	A	907	5/5	0.97	0.19	47,51,61,70	0
4	SO4	A	906	5/5	0.97	0.06	37,38,39,44	0
4	SO4	B	905	5/5	0.99	0.06	29,29,39,42	0

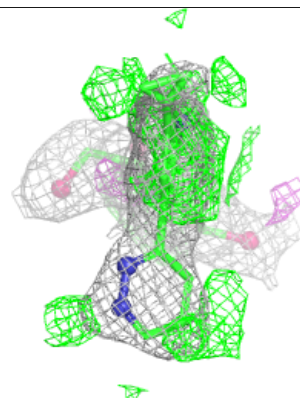
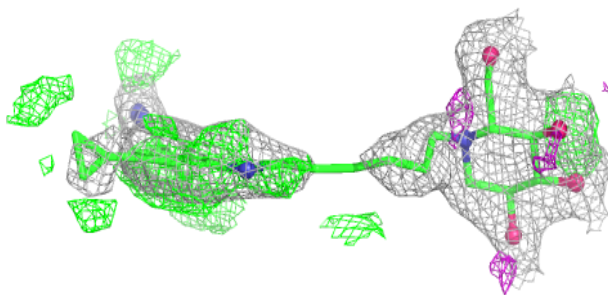
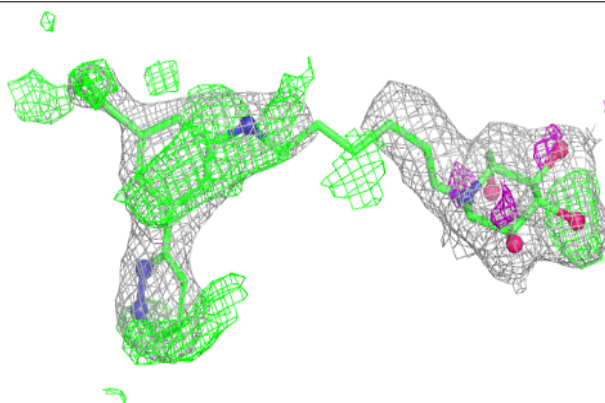
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

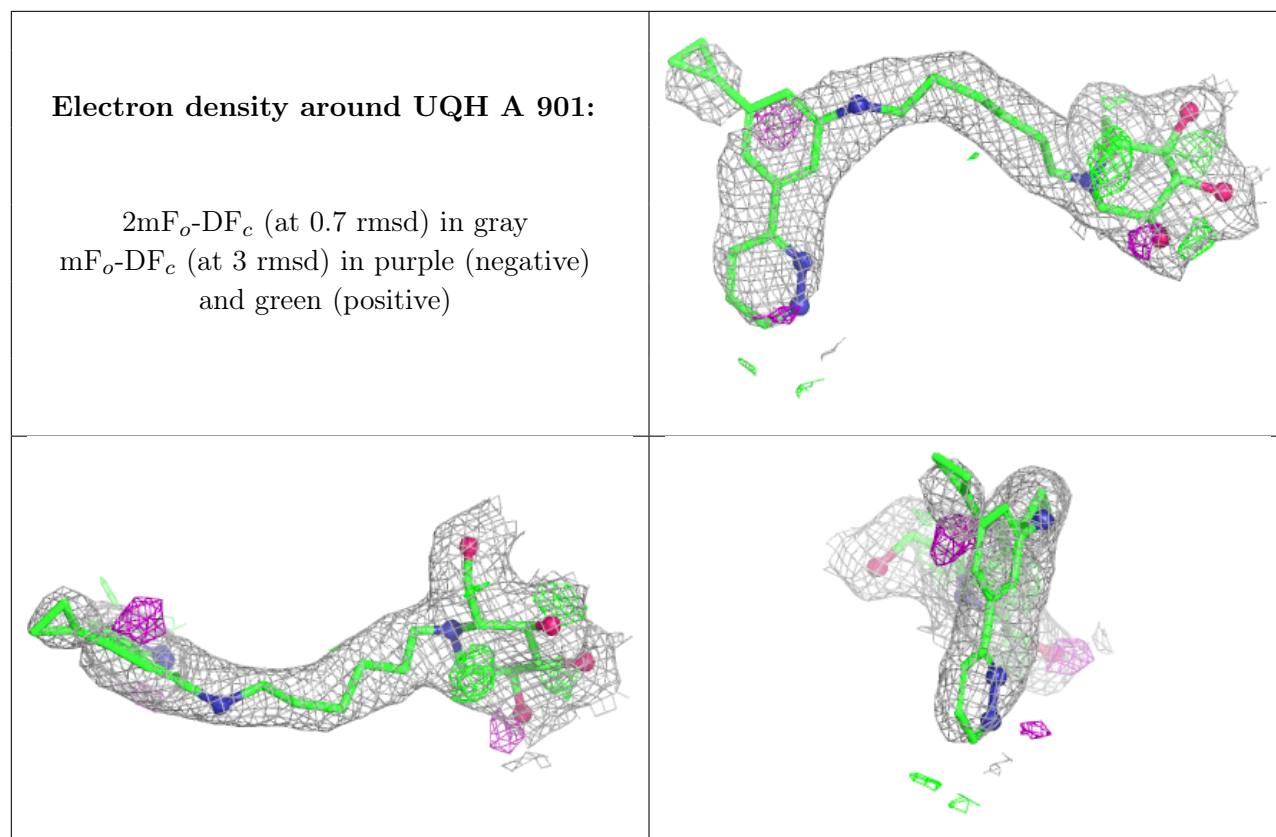
Electron density around UQH B 901 (B):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around UQH B 901 (A):**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.