



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 11:46 AM UTC

PDB ID : 7G30 / pdb_00007g30

Title : Crystal Structure of rat Autotaxin in complex with 3-fluoro-4-[[1-[3-[2-[(5-methyltetrazol-2-yl)methyl]-4-(trifluoromethyl)phenyl]propanoyl]piperidin-4-yl]methylsulfanyl]benzenesulfonamide, i.e. SMILES c1(c(F)cc(S(=O)(=O)N)cc1)SC[C@H]1CCN(CC1)C(=O)CCc1ccc(cc1CN1N=NC(=N1)C)C(F)(F)F with IC50=0.0105744 microM

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Deposited on : 2023-06-05

Resolution : 1.88 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)

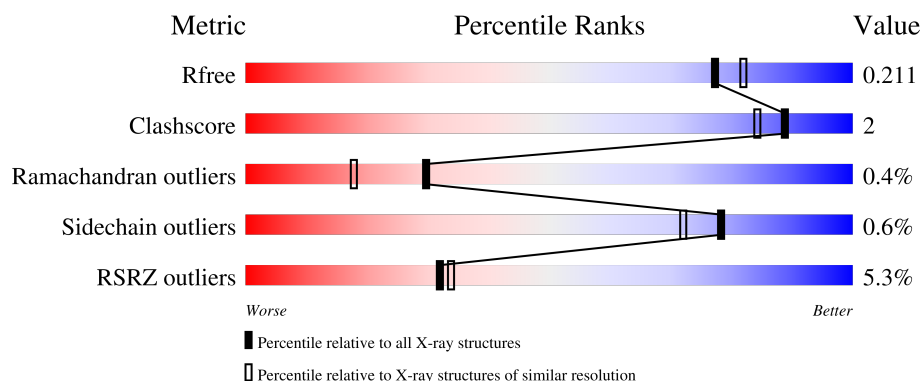
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.88 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	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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	846	5% 90% 5% 5%
2	B	8	100%

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 7120 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.

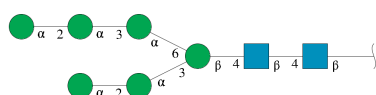
- Molecule 1 is a protein called Isoform 2 of Ectonucleotide pyrophosphatase/phosphodiesterase family member 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	804	Total	C	N	O	S	0	15	0
			6575	4175	1130	1218	52			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	53	ALA	ASN	engineered mutation	UNP Q64610
A	410	ALA	ASN	engineered mutation	UNP Q64610
A	591	THR	ARG	engineered mutation	UNP Q64610
A	863	GLY	-	expression tag	UNP Q64610
A	864	GLY	-	expression tag	UNP Q64610
A	865	ARG	-	expression tag	UNP Q64610
A	866	HIS	-	expression tag	UNP Q64610
A	867	HIS	-	expression tag	UNP Q64610
A	868	HIS	-	expression tag	UNP Q64610
A	869	HIS	-	expression tag	UNP Q64610
A	870	HIS	-	expression tag	UNP Q64610
A	871	HIS	-	expression tag	UNP Q64610
A	872	HIS	-	expression tag	UNP Q64610
A	873	HIS	-	expression tag	UNP Q64610

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	8	Total	C	N	O	0	0	0
			94	52	2	40			

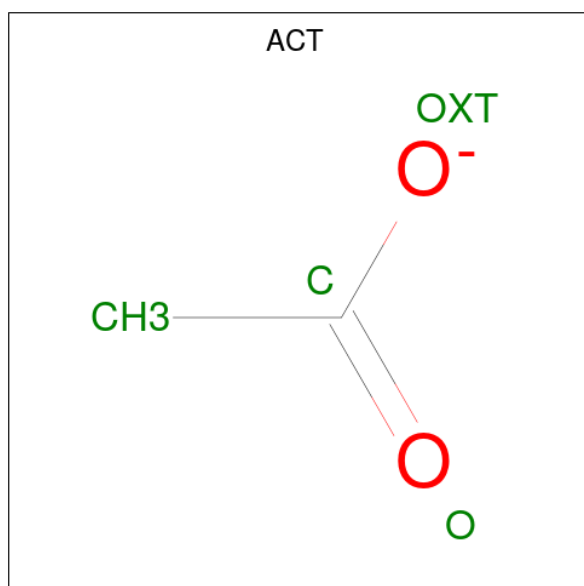
- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 4 is POTASSIUM ION (CCD ID: K) (formula: K).

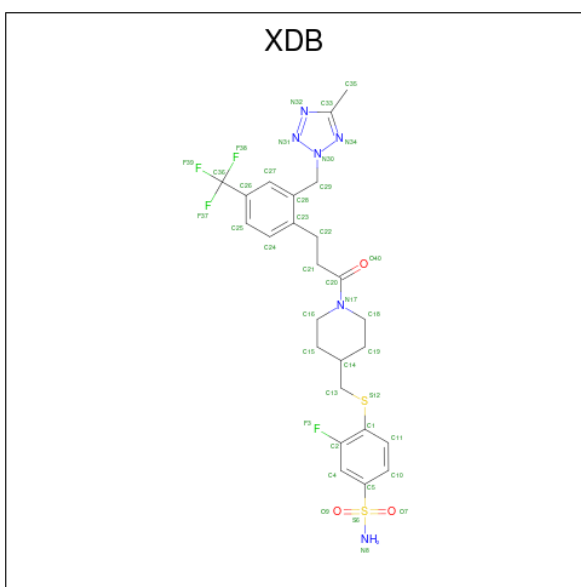
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	K	0	0
			1	1		

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



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

- Molecule 6 is 3-fluoro-4-([1-(3-{2-[(5-methyl-2H-tetrazol-2-yl)methyl]-4-(trifluoromethyl)phenyl}propanoyl)piperidin-4-yl)methyl}sulfanyl)benzene-1-sulfonamide (CCD ID: XDB) (formula: C₂₅H₂₈F₄N₆O₃S₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	F	N	O	S	
			40	25	4	6	3	2	0

- Molecule 7 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 8 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total	Na		
			1	1	0	0

- Molecule 9 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	402	Total	O		
			402	402	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	84.09Å 92.16Å 120.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.08 – 1.88 46.08 – 1.88	Depositor EDS
% Data completeness (in resolution range)	95.2 (46.08-1.88) 95.2 (46.08-1.88)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.86 (at 1.88Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.190 , 0.228 (Not available) , 0.211	Depositor DCC
R_{free} test set	3858 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	28.1	Xtriage
Anisotropy	0.204	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 25.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7120	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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, ZN, CL, MAN, BMA, CA, K, ACT, XDB, NA

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	1.00	1/6797 (0.0%)	0.99	9/9212 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	539	PHE	CA-C	-5.21	1.46	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	539	PHE	N-CA-C	-7.28	97.03	108.90
1	A	577	LYS	N-CA-C	5.99	118.59	111.71
1	A	279	ILE	CA-C-N	-5.88	114.21	120.03
1	A	279	ILE	C-N-CA	-5.88	114.21	120.03
1	A	827	ARG	CG-CD-NE	-5.72	99.41	112.00
1	A	540	ARG	CB-CA-C	5.36	116.22	110.65
1	A	749	GLU	CB-CA-C	-5.22	102.12	110.79
1	A	147	ASP	CA-C-N	-5.03	113.47	122.07
1	A	147	ASP	C-N-CA	-5.03	113.47	122.07

There are no chirality outliers.

There are no planarity outliers.

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	6575	0	6362	25	0
2	B	94	0	79	0	0
3	A	1	0	0	0	0
4	A	1	0	0	0	0
5	A	4	0	3	0	0
6	A	40	0	0	0	0
7	A	1	0	0	0	0
8	A	1	0	0	0	0
9	A	1	0	0	0	0
10	A	402	0	0	3	0
All	All	7120	0	6444	25	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 2.

All (25) 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:148:CYS:HB2	1:A:194[B]:CYS:SG	2.36	0.64
1:A:682:PRO:HB3	1:A:716:GLN:HB3	1.89	0.54
1:A:148:CYS:CB	1:A:194[B]:CYS:SG	2.96	0.53
1:A:793:PHE:CE2	1:A:819:MET:HE3	2.44	0.52
1:A:269:ARG:HB3	1:A:300:PRO:HA	1.94	0.50
1:A:163:PRO:HB3	1:A:350:CYS:O	2.13	0.49
1:A:148:CYS:CB	1:A:194[B]:CYS:HG	2.26	0.49
1:A:539:PHE:O	1:A:541:PRO:HD3	2.13	0.49
1:A:584:ARG:O	1:A:588:LYS:HG2	2.14	0.48
1:A:148:CYS:HB2	1:A:194[B]:CYS:HG	1.77	0.48
1:A:166:ILE:HD12	1:A:351:VAL:HG11	1.96	0.47
1:A:222:PRO:HA	1:A:225:HIS:CE1	2.50	0.47
1:A:201:MET:HE1	1:A:356:VAL:CG2	2.45	0.46
1:A:535:ARG:HG2	1:A:536:THR:HG23	1.97	0.45
1:A:148:CYS:SG	1:A:194[B]:CYS:SG	3.08	0.45
1:A:371:PHE:CE1	1:A:457:PRO:HA	2.52	0.45
1:A:539:PHE:HA	10:A:1037:HOH:O	2.18	0.44
1:A:192:ARG:NH1	10:A:1008:HOH:O	2.46	0.42
1:A:450[A]:ARG:NH1	10:A:1001:HOH:O	2.22	0.42
1:A:178:MET:HE3	1:A:192:ARG:HD2	2.02	0.42
1:A:540:ARG:HD3	1:A:540:ARG:HA	1.88	0.41
1:A:128:ASP:HB3	1:A:290:GLN:OE1	2.21	0.41
1:A:458:LEU:HA	1:A:461:TYR:CE2	2.57	0.40
1:A:721[A]:LYS:HD2	1:A:721[A]:LYS:HA	1.73	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:858[B]:TYR:CD1	1:A:858[B]:TYR:N	2.89	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	813/846 (96%)	785 (97%)	25 (3%)	3 (0%)	30 18

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	378	ASN
1	A	477	ASP
1	A	386	PRO

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	742/765 (97%)	738 (100%)	4 (0%)	81 76

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	183	LYS
1	A	389	LEU
1	A	414	LYS
1	A	458	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	436	ASN
1	A	471	GLN
1	A	484	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 ⓘ

8 monosaccharides 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	B	1	2,1	14,14,15	0.74	1 (7%)	17,19,21	1.17	1 (5%)
2	NAG	B	2	2	14,14,15	0.66	0	17,19,21	2.49	4 (23%)
2	BMA	B	3	2	11,11,12	1.04	1 (9%)	15,15,17	1.88	1 (6%)
2	MAN	B	4	2	11,11,12	1.18	0	15,15,17	1.51	2 (13%)
2	MAN	B	5	2	11,11,12	0.99	0	15,15,17	1.49	3 (20%)
2	MAN	B	6	2	11,11,12	0.79	0	15,15,17	1.10	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MAN	B	7	2	11,11,12	1.34	1 (9%)	15,15,17	1.37	3 (20%)
2	MAN	B	8	2	11,11,12	0.92	0	15,15,17	1.25	2 (13%)

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	B	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1
2	BMA	B	3	2	-	0/2/19/22	0/1/1/1
2	MAN	B	4	2	-	2/2/19/22	0/1/1/1
2	MAN	B	5	2	-	0/2/19/22	0/1/1/1
2	MAN	B	6	2	-	1/2/19/22	0/1/1/1
2	MAN	B	7	2	-	2/2/19/22	0/1/1/1
2	MAN	B	8	2	-	0/2/19/22	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	7	MAN	C1-C2	2.49	1.58	1.52
2	B	1	NAG	C1-C2	-2.11	1.49	1.52
2	B	3	BMA	O3-C3	2.07	1.48	1.43

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	NAG	C1-O5-C5	7.96	122.85	112.19
2	B	3	BMA	C1-O5-C5	5.77	119.92	112.19
2	B	2	NAG	C2-N2-C7	4.49	128.92	122.90
2	B	4	MAN	C1-O5-C5	3.49	116.87	112.19
2	B	7	MAN	C1-C2-C3	2.89	113.85	109.64
2	B	8	MAN	O3-C3-C2	2.82	115.81	110.05
2	B	5	MAN	C3-C4-C5	2.73	115.18	110.23
2	B	6	MAN	O3-C3-C2	-2.32	105.32	110.05
2	B	2	NAG	O7-C7-N2	2.27	125.99	121.98
2	B	2	NAG	O7-C7-C8	-2.27	118.02	122.05
2	B	5	MAN	O5-C1-C2	-2.26	105.40	110.79
2	B	7	MAN	O2-C2-C3	-2.24	105.52	110.15
2	B	7	MAN	O5-C5-C6	2.23	112.01	107.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	4	MAN	O3-C3-C2	2.16	114.47	110.05
2	B	8	MAN	O5-C5-C6	2.16	111.86	107.66
2	B	1	NAG	C2-N2-C7	-2.03	120.18	122.90
2	B	5	MAN	C1-O5-C5	2.00	114.87	112.19

There are no chirality outliers.

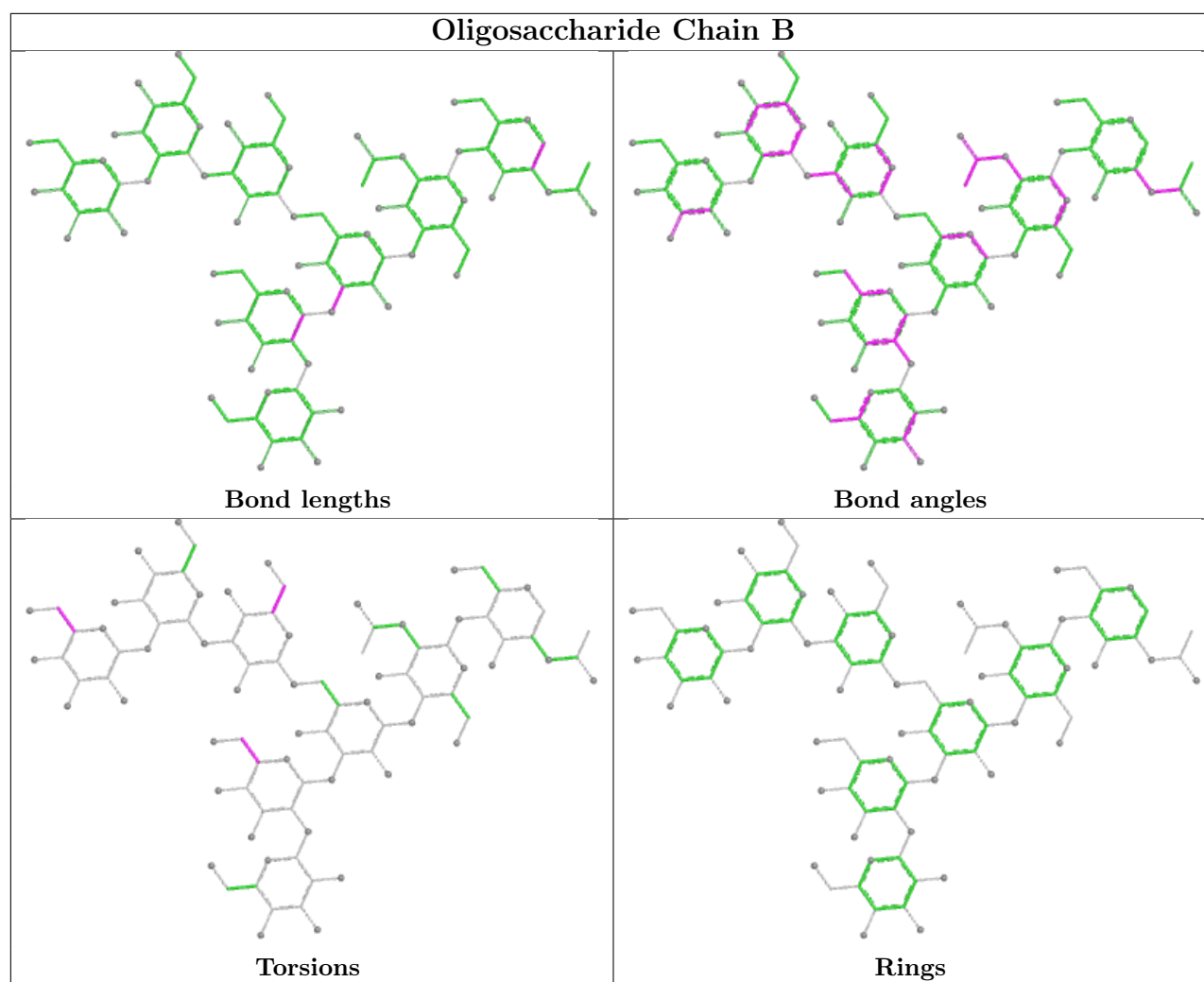
All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	4	MAN	C4-C5-C6-O6
2	B	4	MAN	O5-C5-C6-O6
2	B	7	MAN	O5-C5-C6-O6
2	B	7	MAN	C4-C5-C6-O6
2	B	6	MAN	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 5 are monoatomic - leaving 2 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
6	XDB	A	904	-	42,43,43	2.38	9 (21%)	56,63,63	1.76	15 (26%)
5	ACT	A	903	-	3,3,3	0.88	0	3,3,3	0.89	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
6	XDB	A	904	-	-	4/30/40/40	0/4/4/4

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	904	XDB	C5-S6	-8.76	1.63	1.77
6	A	904	XDB	C1-S12	-6.45	1.67	1.77
6	A	904	XDB	N30-N34	6.00	1.42	1.33
6	A	904	XDB	C20-N17	4.45	1.43	1.35
6	A	904	XDB	N30-N31	4.24	1.38	1.32
6	A	904	XDB	C25-C26	-2.53	1.35	1.39
6	A	904	XDB	C22-C23	-2.44	1.47	1.51
6	A	904	XDB	C29-N30	-2.31	1.42	1.46
6	A	904	XDB	C19-C14	2.24	1.58	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	904	XDB	O9-S6-C5	5.02	113.03	107.35
6	A	904	XDB	C5-S6-N8	-4.33	102.38	108.40
6	A	904	XDB	C15-C14-C13	-3.11	105.31	111.53
6	A	904	XDB	C2-C1-S12	-3.06	115.25	121.78
6	A	904	XDB	C13-S12-C1	3.04	108.89	102.42
6	A	904	XDB	C29-C28-C27	-2.85	115.32	120.37
6	A	904	XDB	C19-C14-C13	2.85	117.23	111.53
6	A	904	XDB	C35-C33-N34	2.42	128.47	122.18
6	A	904	XDB	C15-C16-N17	-2.38	105.99	110.66
6	A	904	XDB	F38-C36-F37	-2.29	97.49	105.77
6	A	904	XDB	C26-C27-C28	-2.26	119.68	122.50
6	A	904	XDB	O7-S6-C5	-2.25	104.81	107.35
6	A	904	XDB	C16-N17-C20	-2.15	115.75	122.83
6	A	904	XDB	C28-C29-N30	-2.07	109.60	111.96
6	A	904	XDB	C22-C21-C20	-2.00	110.06	112.33

There are no chirality outliers.

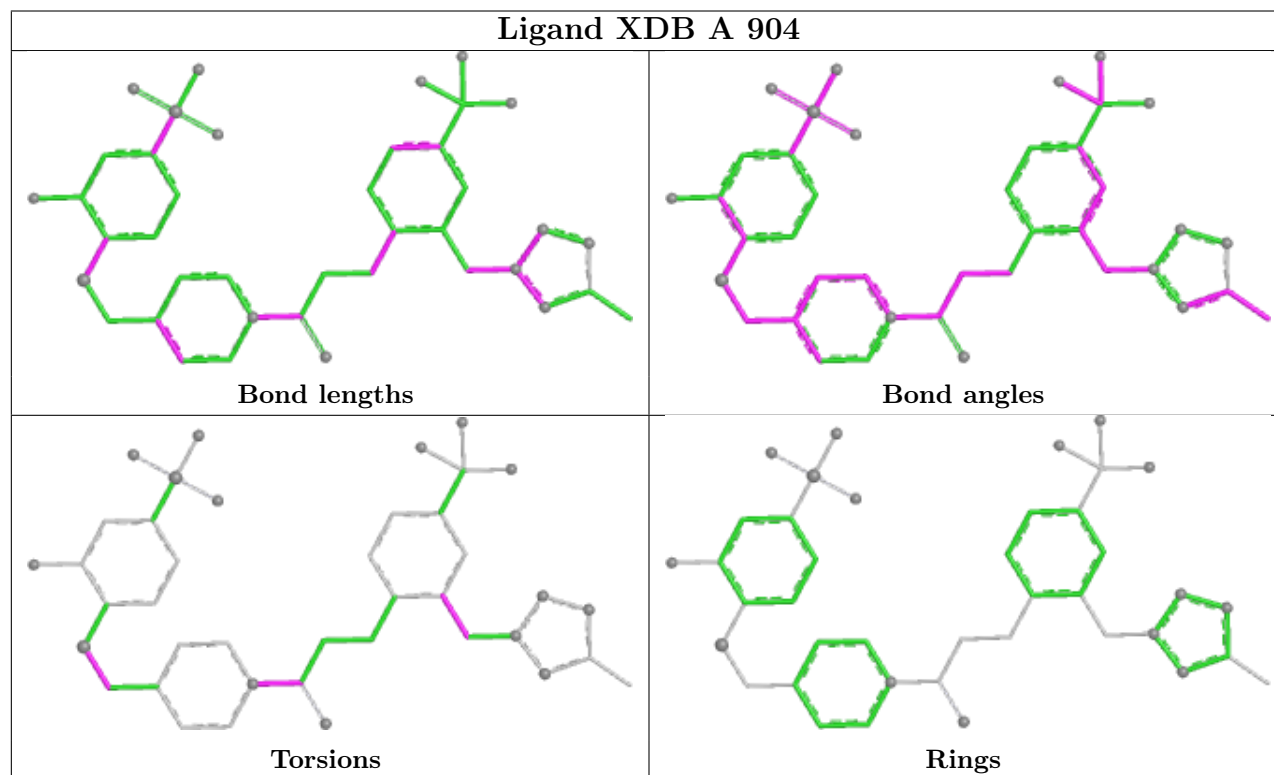
All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	904	XDB	C14-C13-S12-C1
6	A	904	XDB	C21-C20-N17-C18
6	A	904	XDB	O40-C20-N17-C18
6	A	904	XDB	C23-C28-C29-N30

There are no ring outliers.

No monomer is involved in short contacts.

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.



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	804/846 (95%)	0.17	43 (5%) 32 34	15, 31, 64, 80	15 (1%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	468	CYS	6.3
1	A	469	PHE	6.2
1	A	578	LEU	5.7
1	A	572	VAL	5.5
1	A	51	TRP	5.1
1	A	397	ILE	4.6
1	A	70	PRO	4.3
1	A	237	PHE	4.1
1	A	68	VAL	3.9
1	A	581	LEU	3.9
1	A	71	PRO	3.5
1	A	463	LYS	3.1
1	A	573	GLU	2.9
1	A	378	ASN	2.9
1	A	366	CYS	2.8
1	A	236	VAL	2.8
1	A	470	PHE	2.7
1	A	538	THR	2.7
1	A	83	SER	2.7
1	A	577	LYS	2.7
1	A	233	TYR	2.6
1	A	243	LEU	2.6
1	A	242	HIS	2.6
1	A	537	ASN	2.5
1	A	461	TYR	2.5
1	A	379	VAL	2.5
1	A	576	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	69	GLY	2.4
1	A	579	GLU	2.4
1	A	239	ALA	2.4
1	A	402	TYR	2.3
1	A	240	SER	2.3
1	A	346	ARG	2.3
1	A	566	CYS	2.3
1	A	462	LYS	2.2
1	A	55	SER	2.2
1	A	398	ASN	2.2
1	A	54	THR	2.2
1	A	53	ALA	2.2
1	A	542	THR	2.1
1	A	417	ASP	2.1
1	A	249	PHE	2.0
1	A	52	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

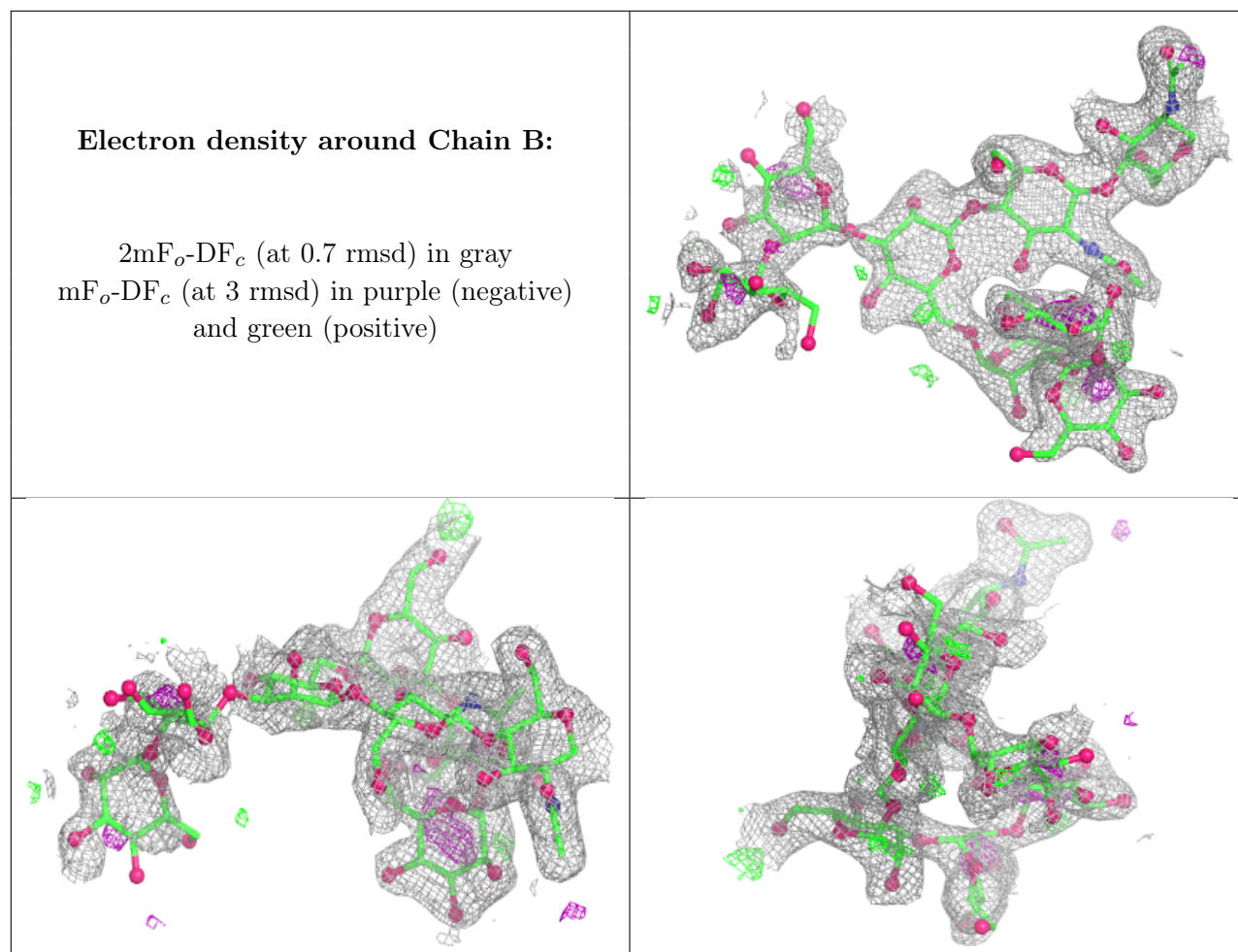
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

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
2	MAN	B	7	11/12	0.40	0.16	80,80,80,80	0
2	BMA	B	3	11/12	0.60	0.15	62,67,76,80	0
2	MAN	B	8	11/12	0.61	0.17	71,80,80,80	0
2	MAN	B	4	11/12	0.71	0.13	59,69,73,73	0
2	MAN	B	5	11/12	0.77	0.14	61,70,75,78	0
2	MAN	B	6	11/12	0.86	0.14	39,55,63,64	0
2	NAG	B	2	14/15	0.91	0.09	33,40,49,56	0
2	NAG	B	1	14/15	0.96	0.07	21,24,26,31	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



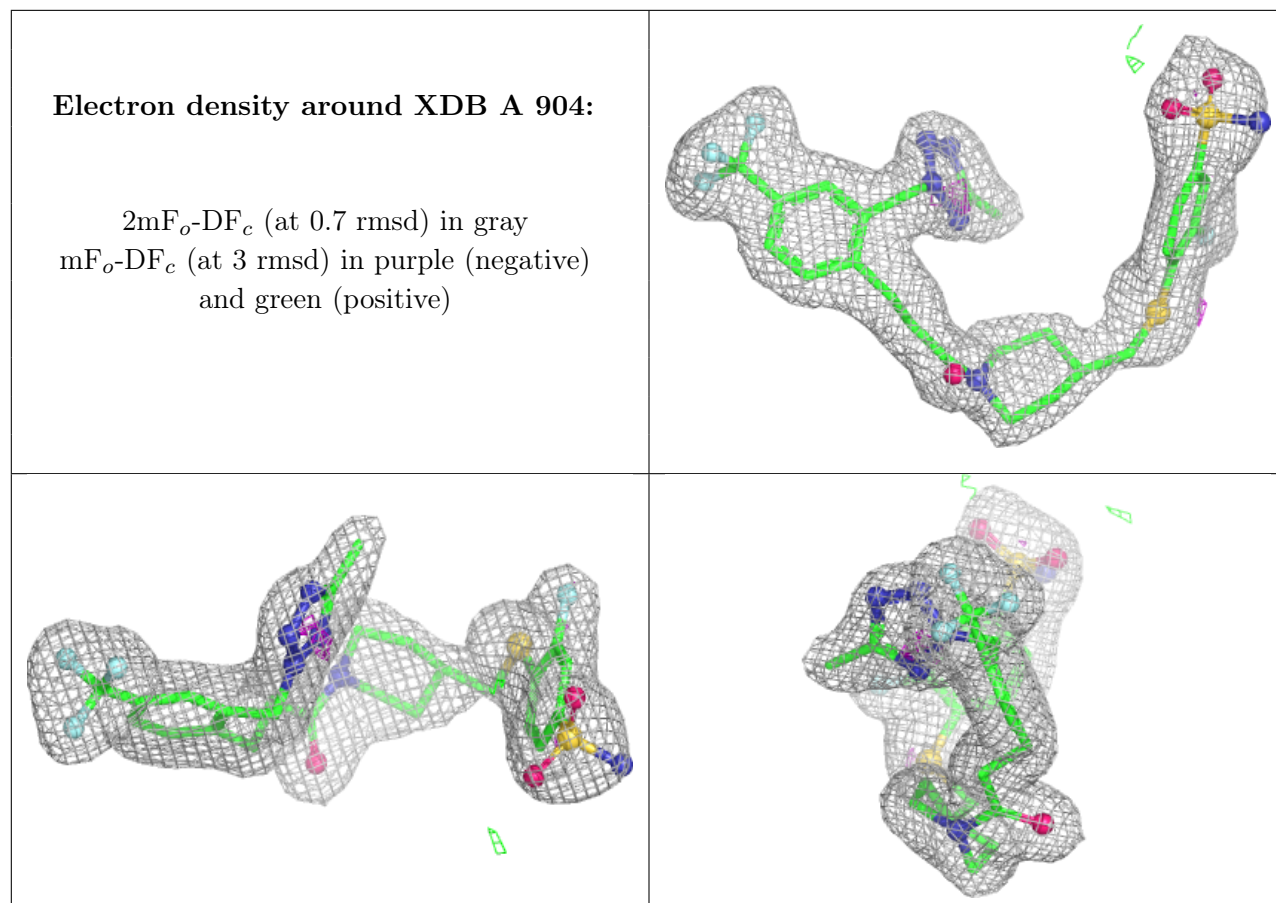
6.4 Ligands [i](#)

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
3	CL	A	901	1/1	0.88	0.13	54,54,54,54	0
5	ACT	A	903	4/4	0.90	0.13	41,41,44,46	0
6	XDB	A	904	40/40	0.95	0.08	27,37,59,60	0
8	NA	A	906	1/1	0.95	0.07	43,43,43,43	0
4	K	A	902	1/1	0.98	0.04	34,34,34,34	0
9	CA	A	907	1/1	0.99	0.03	23,23,23,23	0
7	ZN	A	905	1/1	1.00	0.01	26,26,26,26	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.