



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 15, 2019 – 12:50 PM EDT

PDB ID : 6IMR
Title : Crystal structure of PDE4D complexed with a novel inhibitor
Authors : Zhang, X.L.; Su, H.X.; Xu, Y.C.
Deposited on : 2018-10-23
Resolution : 1.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.0 (224370), CSD as540be (2019)
Xtriage (Phenix) : 1.13
EDS : 2.4
buster-report : 1.1.7 (2018)
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) : 2.4

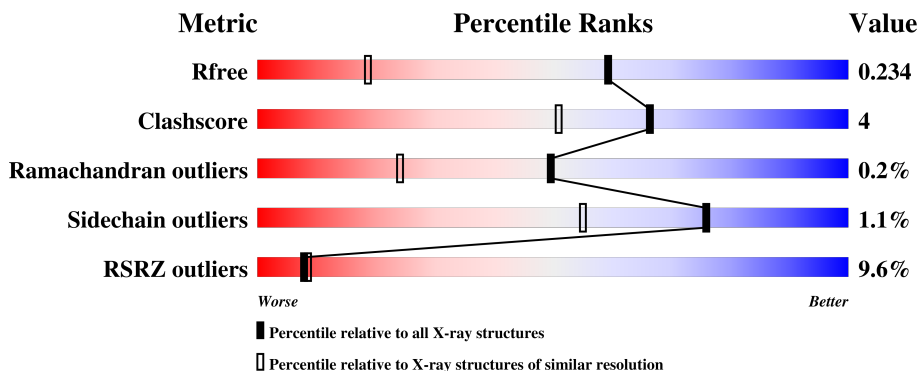
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

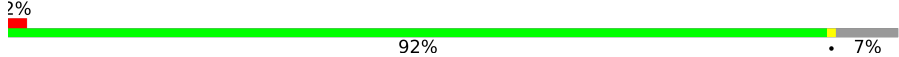
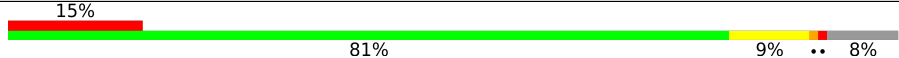
The reported resolution of this entry is 1.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	111664	2534 (1.50-1.50)
Clashscore	122126	2727 (1.50-1.50)
Ramachandran outliers	120053	2661 (1.50-1.50)
Sidechain outliers	120020	2659 (1.50-1.50)
RSRZ outliers	108989	2481 (1.50-1.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	349	
1	B	349	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5684 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 cAMP-specific 3',5'-cyclic phosphodiesterase 4D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	0	0	0
			2604	1651	444	495	14			
1	B	320	Total	C	N	O	S	0	0	0
			2517	1601	428	474	14			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	65	MET	-	expression tag	UNP Q08499
A	66	GLY	-	expression tag	UNP Q08499
A	67	SER	-	expression tag	UNP Q08499
A	68	SER	-	expression tag	UNP Q08499
A	69	HIS	-	expression tag	UNP Q08499
A	70	HIS	-	expression tag	UNP Q08499
A	71	HIS	-	expression tag	UNP Q08499
A	72	HIS	-	expression tag	UNP Q08499
A	73	HIS	-	expression tag	UNP Q08499
A	74	HIS	-	expression tag	UNP Q08499
A	75	SER	-	expression tag	UNP Q08499
A	76	SER	-	expression tag	UNP Q08499
A	77	GLY	-	expression tag	UNP Q08499
A	78	LEU	-	expression tag	UNP Q08499
A	79	VAL	-	expression tag	UNP Q08499
A	80	PRO	-	expression tag	UNP Q08499
A	81	ARG	-	expression tag	UNP Q08499
A	82	GLY	-	expression tag	UNP Q08499
A	83	SER	-	expression tag	UNP Q08499
A	84	HIS	-	expression tag	UNP Q08499
A	85	MET	-	expression tag	UNP Q08499
B	65	MET	-	expression tag	UNP Q08499
B	66	GLY	-	expression tag	UNP Q08499
B	67	SER	-	expression tag	UNP Q08499
B	68	SER	-	expression tag	UNP Q08499

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Chain	Residue	Modelled	Actual	Comment	Reference
B	69	HIS	-	expression tag	UNP Q08499
B	70	HIS	-	expression tag	UNP Q08499
B	71	HIS	-	expression tag	UNP Q08499
B	72	HIS	-	expression tag	UNP Q08499
B	73	HIS	-	expression tag	UNP Q08499
B	74	HIS	-	expression tag	UNP Q08499
B	75	SER	-	expression tag	UNP Q08499
B	76	SER	-	expression tag	UNP Q08499
B	77	GLY	-	expression tag	UNP Q08499
B	78	LEU	-	expression tag	UNP Q08499
B	79	VAL	-	expression tag	UNP Q08499
B	80	PRO	-	expression tag	UNP Q08499
B	81	ARG	-	expression tag	UNP Q08499
B	82	GLY	-	expression tag	UNP Q08499
B	83	SER	-	expression tag	UNP Q08499
B	84	HIS	-	expression tag	UNP Q08499
B	85	MET	-	expression tag	UNP Q08499

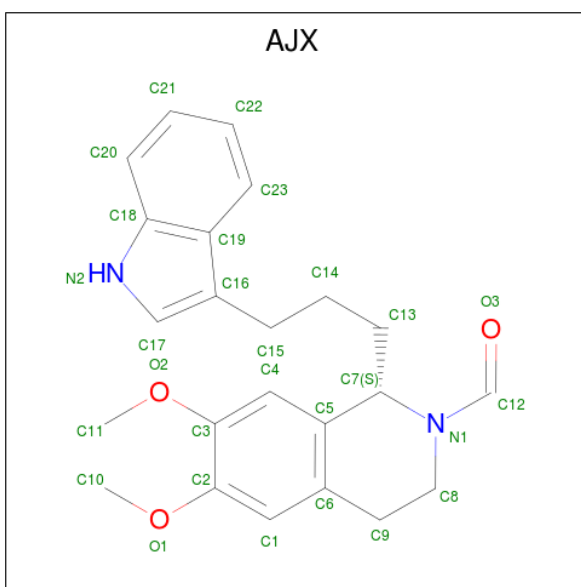
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Zn 1 1	0	0
2	A	1	Total Zn 1 1	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

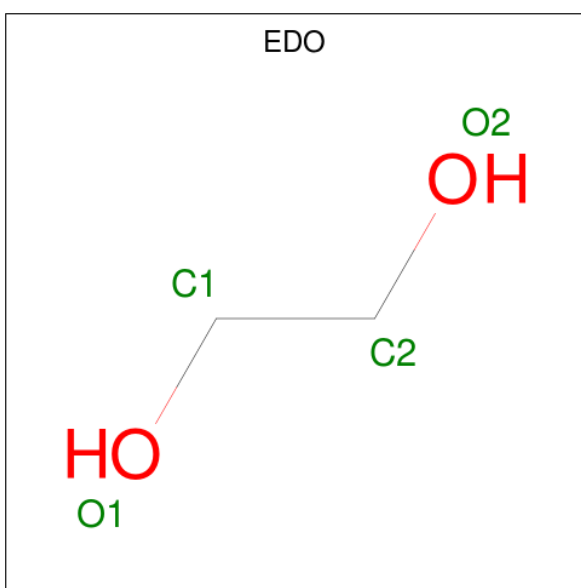
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Mg 1 1	0	0
3	A	1	Total Mg 1 1	0	0

- Molecule 4 is (1S)-1-[3-(1H-indol-3-yl)propyl]-6,7-dimethoxy-3,4-dihydroisoquinoline-2(1H)-carbaldehyde (three-letter code: AJX) (formula: C₂₃H₂₆N₂O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			28	23	2	3		
4	B	1	Total	C	N	O	0	0
			28	23	2	3		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	273	Total O 273 273	0	0
6	B	174	Total O 174 174	0	0

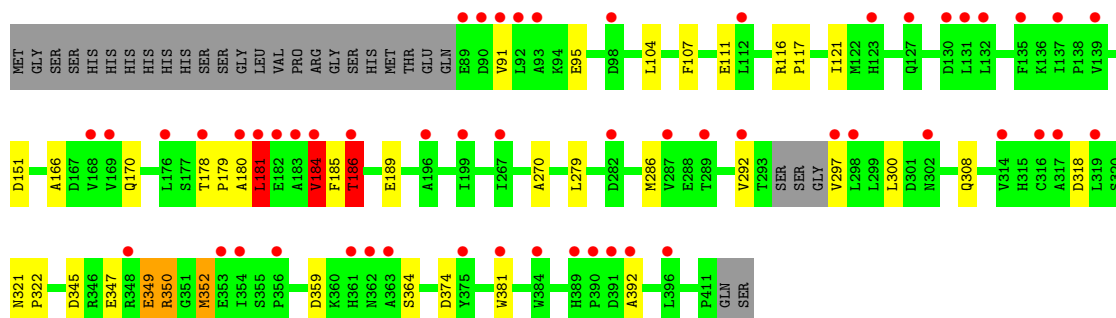
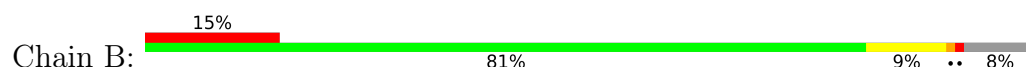
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: cAMP-specific 3',5'-cyclic phosphodiesterase 4D



- Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4D



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.12Å 80.62Å 163.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.90 – 1.50 30.90 – 1.50	Depositor EDS
% Data completeness (in resolution range)	99.5 (30.90-1.50) 99.5 (30.90-1.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.10 (at 1.50Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, R_{free}	0.211 , 0.237 0.212 , 0.234	Depositor DCC
R_{free} test set	6014 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	12.2	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 51.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5684	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.80% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: AJX, ZN, MG, EDO

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	0.37	0/2658	0.53	0/3616
1	B	0.99	7/2570 (0.3%)	0.85	12/3505 (0.3%)
All	All	0.74	7/5228 (0.1%)	0.70	12/7121 (0.2%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	349	GLU	CD-OE2	-6.20	1.18	1.25
1	B	300	LEU	C-O	-5.84	1.12	1.23
1	B	95	GLU	CD-OE2	-5.84	1.19	1.25
1	B	352	MET	C-O	-5.34	1.13	1.23
1	B	381	TRP	CE3-CZ3	-5.33	1.29	1.38
1	B	185	PHE	N-CA	-5.18	1.35	1.46
1	B	364	SER	C-O	-5.03	1.13	1.23

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	181	LEU	CB-CG-CD1	10.04	128.07	111.00
1	B	374	ASP	CB-CG-OD2	-7.15	111.87	118.30
1	B	350	ARG	NE-CZ-NH2	-6.67	116.97	120.30
1	B	184	VAL	N-CA-C	6.36	128.18	111.00
1	B	181	LEU	CB-CG-CD2	-6.14	100.56	111.00
1	B	318	ASP	CB-CG-OD2	-6.00	112.91	118.30
1	B	364	SER	N-CA-CB	5.85	119.27	110.50
1	B	359	ASP	CB-CG-OD2	-5.70	113.17	118.30
1	B	184	VAL	CA-CB-CG2	-5.68	102.38	110.90
1	B	186	THR	OG1-CB-CG2	-5.54	97.27	110.00
1	B	185	PHE	N-CA-C	-5.33	96.60	111.00
1	B	181	LEU	CA-CB-CG	5.30	127.49	115.30

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	2604	0	2540	2	0
1	B	2517	0	2408	39	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	28	0	0	1	0
4	B	28	0	0	1	0
5	A	28	0	42	1	0
5	B	28	0	42	2	0
6	A	273	0	0	0	0
6	B	174	0	0	2	0
All	All	5684	0	5032	44	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 4.

All (44) 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:B:350:ARG:HD2	1:B:352:MET:CE	1.65	1.26
1:B:350:ARG:CD	1:B:352:MET:HE2	1.68	1.21
1:B:350:ARG:CD	1:B:352:MET:CE	2.27	1.09
1:B:350:ARG:HD3	1:B:352:MET:HE2	1.32	1.08
1:B:186:THR:HG22	1:B:189:GLU:CD	1.81	1.00
1:B:186:THR:HG22	1:B:189:GLU:OE1	1.62	0.99
1:B:186:THR:CG2	1:B:189:GLU:CG	2.45	0.95
1:B:186:THR:HG23	1:B:189:GLU:CG	1.99	0.93
1:B:186:THR:CG2	1:B:189:GLU:CD	2.39	0.91
1:B:350:ARG:HD2	1:B:352:MET:HE1	1.50	0.91
1:B:186:THR:CG2	1:B:189:GLU:HG3	2.05	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:186:THR:HG23	1:B:189:GLU:CB	2.09	0.83
1:B:186:THR:HG23	1:B:189:GLU:HG3	1.69	0.73
1:B:186:THR:HG23	1:B:189:GLU:HB2	1.72	0.71
1:B:184:VAL:HG13	1:B:184:VAL:O	1.91	0.70
1:B:186:THR:CG2	1:B:189:GLU:OE1	2.43	0.63
1:B:186:THR:HG21	1:B:189:GLU:HG3	1.82	0.61
1:A:290:LYS:HA	5:A:507:EDO:H11	1.86	0.57
1:B:292:VAL:HA	1:B:297:VAL:O	2.04	0.56
1:B:350:ARG:HB2	1:B:352:MET:CE	2.35	0.56
1:B:350:ARG:CB	1:B:352:MET:HE2	2.35	0.56
1:B:392:ALA:HA	6:B:744:HOH:O	2.09	0.53
1:B:345:ASP:O	1:B:349:GLU:HG3	2.09	0.52
1:B:116:ARG:N	1:B:117:PRO:CD	2.73	0.51
1:B:350:ARG:HB2	1:B:352:MET:HE3	1.92	0.51
1:A:282:ASP:HB3	1:A:308:GLN:NE2	2.26	0.51
1:B:104:LEU:HD22	1:B:170:GLN:HG3	1.93	0.50
1:B:350:ARG:HD3	1:B:352:MET:CE	2.15	0.50
1:B:186:THR:HG23	1:B:189:GLU:H	1.78	0.49
5:B:506:EDO:H22	6:B:625:HOH:O	2.12	0.48
1:B:179:PRO:O	1:B:181:LEU:N	2.46	0.48
1:B:107:PHE:O	1:B:111:GLU:HG3	2.13	0.47
1:B:292:VAL:CA	1:B:297:VAL:O	2.62	0.47
1:B:350:ARG:HD2	1:B:352:MET:HE3	1.82	0.47
1:B:151:ASP:HA	5:B:506:EDO:H21	1.98	0.45
1:B:321:ASN:HB2	1:B:322:PRO:HD3	1.99	0.44
1:B:347:GLU:HA	1:B:352:MET:HE3	1.98	0.44
4:B:503:AJX:C13	4:B:503:AJX:C17	2.96	0.44
4:A:503:AJX:C17	4:A:503:AJX:C13	2.95	0.44
1:B:178:THR:HB	1:B:181:LEU:HD22	1.99	0.44
1:B:350:ARG:CB	1:B:352:MET:CE	2.95	0.44
1:B:121:ILE:HD12	1:B:166:ALA:HB1	2.01	0.42
1:B:270:ALA:HB1	1:B:279:LEU:HD11	2.02	0.41
1:B:179:PRO:C	1:B:181:LEU:H	2.24	0.41

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	323/349 (93%)	317 (98%)	6 (2%)	0	100	100
1	B	316/349 (90%)	307 (97%)	8 (2%)	1 (0%)	43	19
All	All	639/698 (92%)	624 (98%)	14 (2%)	1 (0%)	49	24

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	180	ALA

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	289/318 (91%)	289 (100%)	0	100	100
1	B	270/318 (85%)	264 (98%)	6 (2%)	55	23
All	All	559/636 (88%)	553 (99%)	6 (1%)	76	55

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	91	VAL
1	B	181	LEU
1	B	184	VAL
1	B	186	THR
1	B	286	MET

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Mol	Chain	Res	Type
1	B	308	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	389	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 4 are monoatomic - leaving 16 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
4	AJX	A	503	-	29,31,31	0.65	0	32,43,43	1.25	5 (15%)
5	EDO	A	504	-	3,3,3	0.46	0	2,2,2	0.44	0
5	EDO	A	505	-	3,3,3	0.44	0	2,2,2	0.60	0
5	EDO	A	506	-	3,3,3	0.51	0	2,2,2	0.50	0
5	EDO	A	507	-	3,3,3	0.47	0	2,2,2	0.45	0
5	EDO	A	508	-	3,3,3	0.47	0	2,2,2	0.39	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	A	509	-	3,3,3	0.47	0	2,2,2	0.61	0
5	EDO	A	510	-	3,3,3	0.45	0	2,2,2	0.54	0
4	AJX	B	503	-	29,31,31	0.61	0	32,43,43	1.13	3 (9%)
5	EDO	B	504	-	3,3,3	0.45	0	2,2,2	0.57	0
5	EDO	B	505	-	3,3,3	0.45	0	2,2,2	0.63	0
5	EDO	B	506	-	3,3,3	0.47	0	2,2,2	0.46	0
5	EDO	B	507	-	3,3,3	0.48	0	2,2,2	0.49	0
5	EDO	B	508	-	3,3,3	0.46	0	2,2,2	0.48	0
5	EDO	B	509	-	3,3,3	0.46	0	2,2,2	0.38	0
5	EDO	B	510	-	3,3,3	0.50	0	2,2,2	0.39	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
4	AJX	A	503	-	-	6/12/25/25	0/4/4/4
5	EDO	A	504	-	-	0/1/1/1	-
5	EDO	A	505	-	-	0/1/1/1	-
5	EDO	A	506	-	-	0/1/1/1	-
5	EDO	A	507	-	-	0/1/1/1	-
5	EDO	A	508	-	-	0/1/1/1	-
5	EDO	A	509	-	-	1/1/1/1	-
5	EDO	A	510	-	-	0/1/1/1	-
4	AJX	B	503	-	-	6/12/25/25	0/4/4/4
5	EDO	B	504	-	-	0/1/1/1	-
5	EDO	B	505	-	-	0/1/1/1	-
5	EDO	B	506	-	-	1/1/1/1	-
5	EDO	B	507	-	-	0/1/1/1	-
5	EDO	B	508	-	-	0/1/1/1	-
5	EDO	B	509	-	-	0/1/1/1	-
5	EDO	B	510	-	-	0/1/1/1	-

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	503	AJX	C4-C5-C7	-2.71	115.30	119.39
4	B	503	AJX	C4-C5-C7	-2.41	115.75	119.39
4	A	503	AJX	C6-C5-C7	2.27	124.22	121.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	503	AJX	C21-C20-C18	-2.24	116.80	120.08
4	B	503	AJX	C21-C20-C18	-2.09	117.02	120.08
4	B	503	AJX	C6-C5-C7	2.06	123.98	121.62
4	A	503	AJX	O1-C2-C1	2.06	127.65	124.14
4	A	503	AJX	C9-C6-C5	-2.00	118.06	121.12

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	503	AJX	C14-C13-C7-C5
4	B	503	AJX	C14-C13-C7-N1
4	B	503	AJX	C14-C15-C16-C19
4	A	503	AJX	C14-C13-C7-C5
4	A	503	AJX	C14-C13-C7-N1
4	A	503	AJX	C14-C15-C16-C19
4	A	503	AJX	C13-C14-C15-C16
4	B	503	AJX	C13-C14-C15-C16
4	B	503	AJX	O3-C12-N1-C8
4	A	503	AJX	O3-C12-N1-C8
5	B	506	EDO	O1-C1-C2-O2
4	A	503	AJX	C7-C13-C14-C15
4	B	503	AJX	C7-C13-C14-C15
5	A	509	EDO	O1-C1-C2-O2

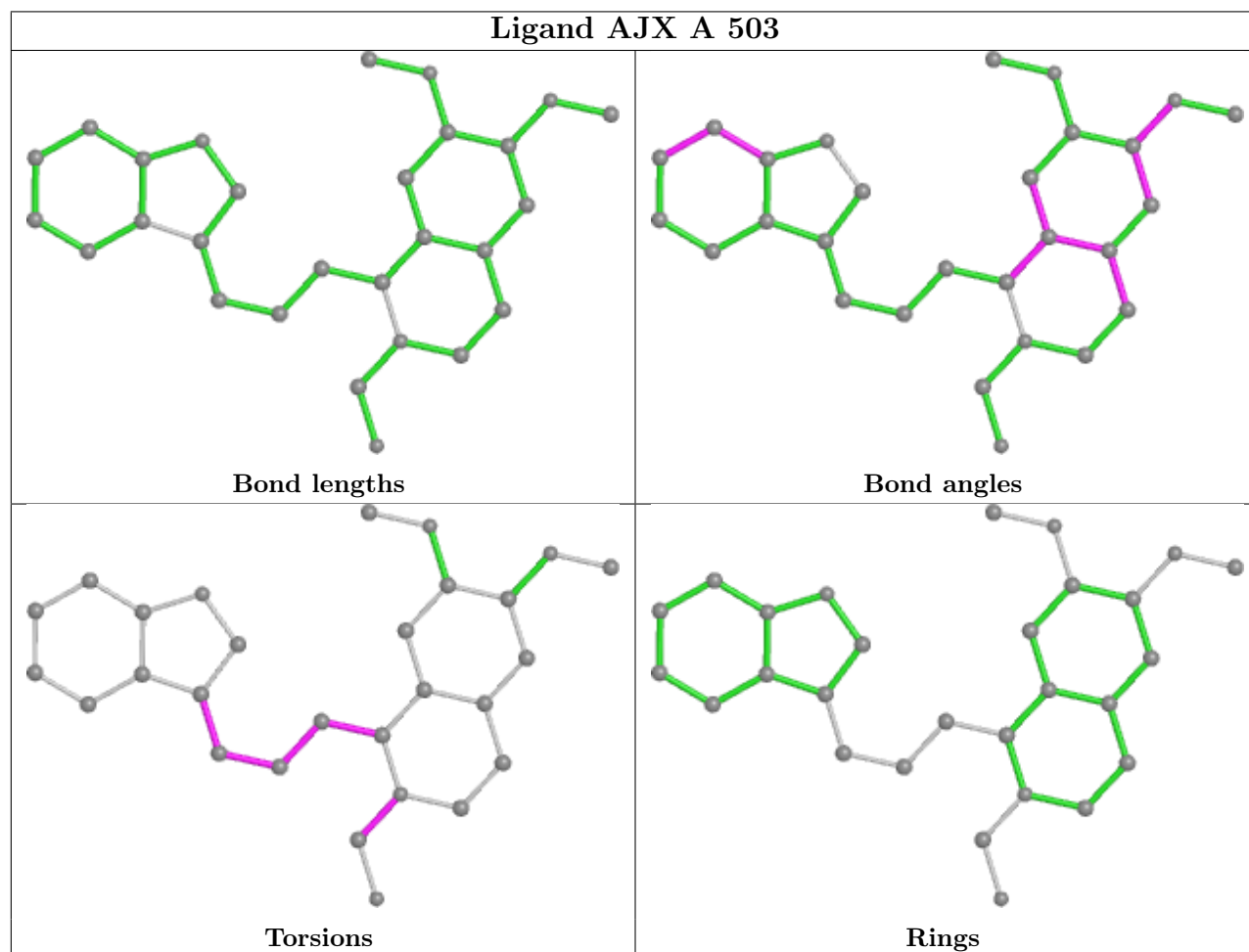
There are no ring outliers.

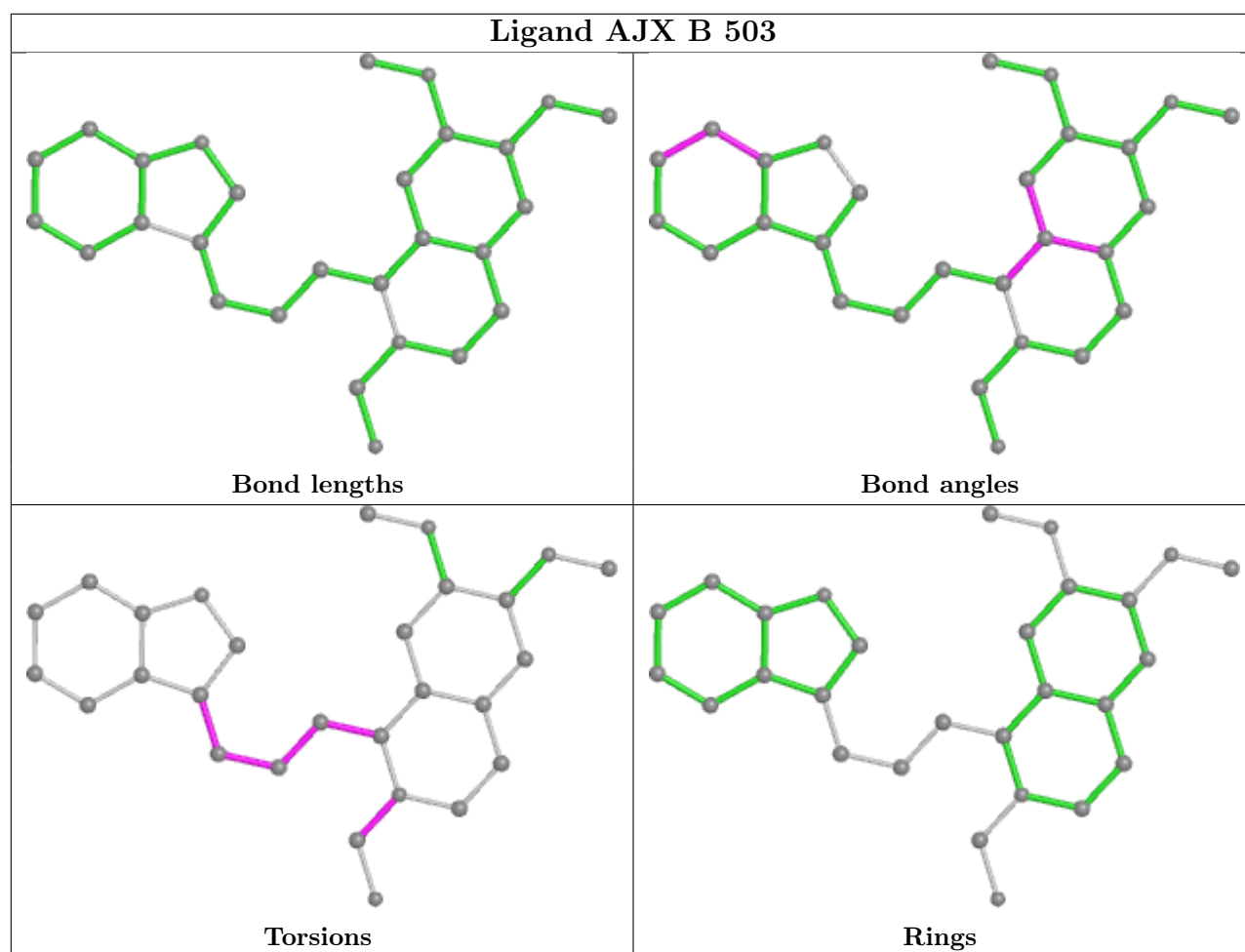
4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	503	AJX	1	0
5	A	507	EDO	1	0
4	B	503	AJX	1	0
5	B	506	EDO	2	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.





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	325/349 (93%)	0.19	8 (2%) 57 63	5, 13, 26, 37	0
1	B	320/349 (91%)	1.05	54 (16%) 1 2	7, 21, 38, 56	0
All	All	645/698 (92%)	0.62	62 (9%) 8 8	5, 17, 34, 56	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	91	VAL	5.7
1	B	390	PRO	5.3
1	B	297	VAL	4.9
1	B	181	LEU	4.6
1	B	183	ALA	4.3
1	B	289	THR	4.3
1	B	314	VAL	4.0
1	B	287	VAL	3.9
1	B	89	GLU	3.9
1	B	186	THR	3.7
1	B	184	VAL	3.7
1	B	363	ALA	3.5
1	B	389	HIS	3.5
1	B	354	ILE	3.4
1	B	375	TYR	3.3
1	B	298	LEU	3.3
1	B	130	ASP	3.1
1	B	392	ALA	3.1
1	B	348	ARG	3.1
1	B	168	VAL	3.1
1	B	93	ALA	3.1
1	B	182	GLU	3.0
1	B	90	ASP	3.0
1	B	137	ILE	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	361	HIS	2.9
1	B	131	LEU	2.9
1	B	391	ASP	2.8
1	B	139	VAL	2.8
1	A	87	GLU	2.8
1	B	319	LEU	2.8
1	B	98	ASP	2.8
1	B	169	VAL	2.8
1	B	317	ALA	2.8
1	A	295	SER	2.7
1	B	178	THR	2.6
1	A	168	VAL	2.6
1	B	302	ASN	2.6
1	A	88	GLN	2.6
1	B	132	LEU	2.6
1	A	362	ASN	2.6
1	B	127	GLN	2.6
1	A	363	ALA	2.5
1	B	199	ILE	2.4
1	B	267	ILE	2.4
1	B	353	GLU	2.4
1	B	92	LEU	2.4
1	B	362	ASN	2.3
1	B	135	PHE	2.3
1	B	381	TRP	2.3
1	B	176	LEU	2.3
1	B	282	ASP	2.3
1	A	296	GLY	2.2
1	A	294	SER	2.2
1	B	292	VAL	2.2
1	B	316	CYS	2.2
1	B	112	LEU	2.2
1	B	384	TRP	2.2
1	B	196	ALA	2.1
1	B	180	ALA	2.1
1	B	123	HIS	2.0
1	B	396	LEU	2.0
1	B	356	PRO	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

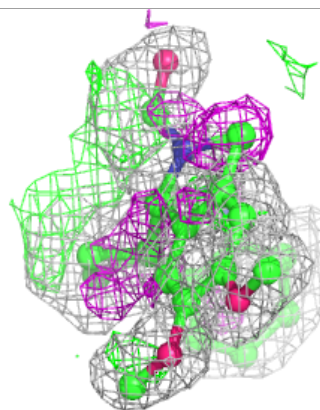
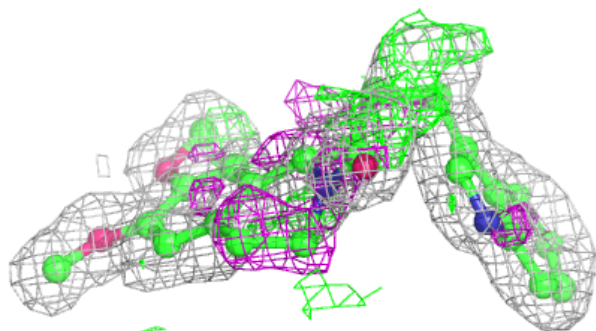
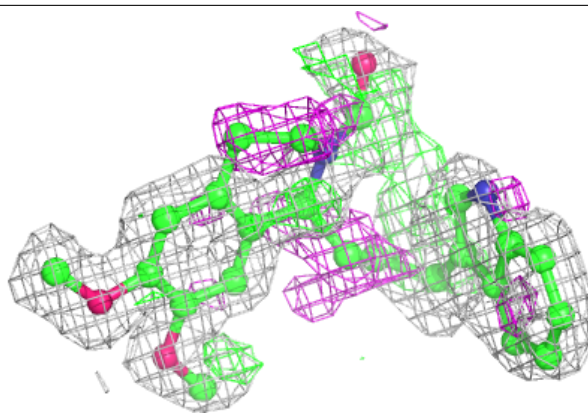
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	EDO	B	508	4/4	0.66	0.19	26,27,30,36	0
5	EDO	A	506	4/4	0.76	0.18	23,26,31,34	0
4	AJX	A	503	28/28	0.79	0.21	9,19,27,33	0
5	EDO	A	507	4/4	0.80	0.15	25,25,25,31	0
4	AJX	B	503	28/28	0.81	0.21	12,22,27,28	0
5	EDO	A	509	4/4	0.82	0.12	20,21,23,24	0
5	EDO	B	505	4/4	0.84	0.18	24,25,26,32	0
5	EDO	B	506	4/4	0.85	0.14	22,25,27,30	0
5	EDO	B	504	4/4	0.88	0.10	20,21,25,25	0
5	EDO	B	507	4/4	0.88	0.18	24,25,28,34	0
5	EDO	B	509	4/4	0.89	0.22	23,25,26,26	0
5	EDO	A	510	4/4	0.91	0.13	17,21,22,26	0
5	EDO	A	505	4/4	0.91	0.10	18,19,20,21	0
5	EDO	B	510	4/4	0.91	0.11	22,22,27,27	0
5	EDO	A	508	4/4	0.94	0.10	19,20,23,31	0
5	EDO	A	504	4/4	0.96	0.06	15,17,17,24	0
2	ZN	B	501	1/1	0.99	0.04	14,14,14,14	0
2	ZN	A	501	1/1	1.00	0.03	11,11,11,11	0
3	MG	B	502	1/1	1.00	0.09	6,6,6,6	0
3	MG	A	502	1/1	1.00	0.05	5,5,5,5	0

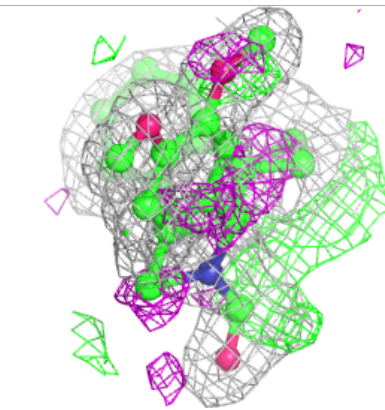
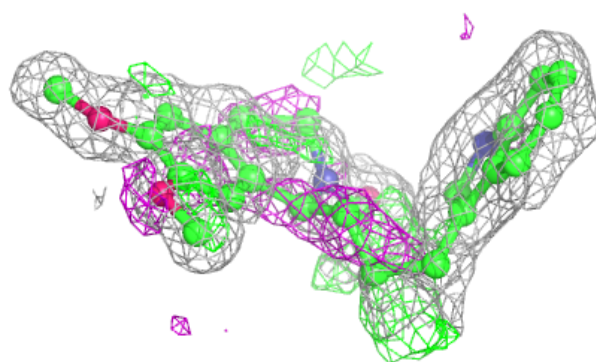
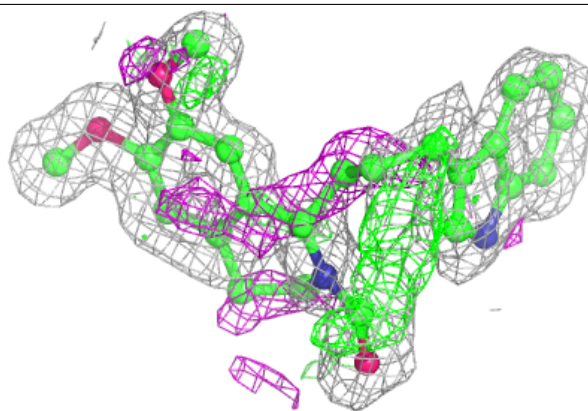
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around AJX A 503:

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

**Electron density around AJX B 503:**

$2mF_o-DF_c$ (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.