



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 24, 2025 – 12:14 PM EDT

PDB ID : 9BXD  
Title : CRYSTAL STRUCTURE OF HIV-1 LM/HS CLADE A/E CRF01 GP120  
CORE IN COMPLEX WITH HZ-IV-188  
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Deposited on : 2024-05-22  
Resolution : 1.85 Å(reported)

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

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

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The types of validation reports are described at

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

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.41.4

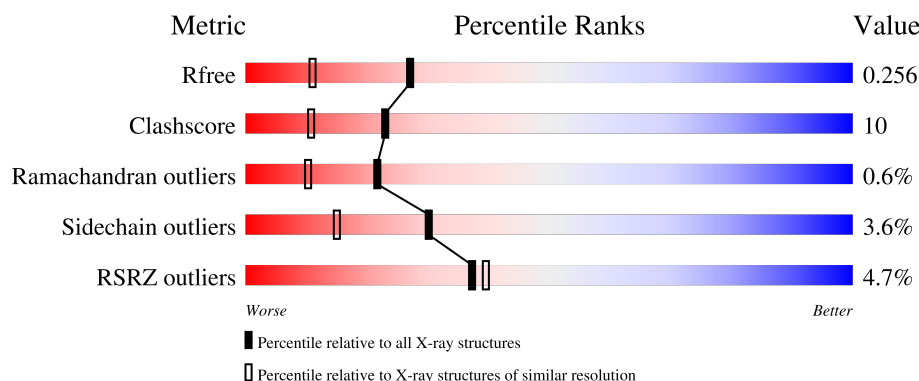
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


The reported resolution of this entry is 1.85 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))
$R_{free}$	164625	3097 (1.86-1.86)
Clashscore	180529	3359 (1.86-1.86)
Ramachandran outliers	177936	3335 (1.86-1.86)
Sidechain outliers	177891	3335 (1.86-1.86)
RSRZ outliers	164620	3097 (1.86-1.86)

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  $\geq 3$ , 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	355	

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3024 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 HIV-1 LM/HS clade A/E CRF01 gp120 core.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	339	Total	C	N	O	S	0	0	0
			2659	1667	460	509	23			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	42	VAL	-	expression tag	UNP A0A0M3KKW9
A	43	PRO	-	expression tag	UNP A0A0M3KKW9
A	61	TYR	HIS	engineered mutation	UNP A0A0M3KKW9
A	105	HIS	GLN	engineered mutation	UNP A0A0M3KKW9
A	108	ILE	VAL	engineered mutation	UNP A0A0M3KKW9
A	375	SER	HIS	engineered mutation	UNP A0A0M3KKW9
A	474	ASP	ASN	engineered mutation	UNP A0A0M3KKW9
A	475	MET	ILE	engineered mutation	UNP A0A0M3KKW9
A	476	ARG	LYS	engineered mutation	UNP A0A0M3KKW9

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



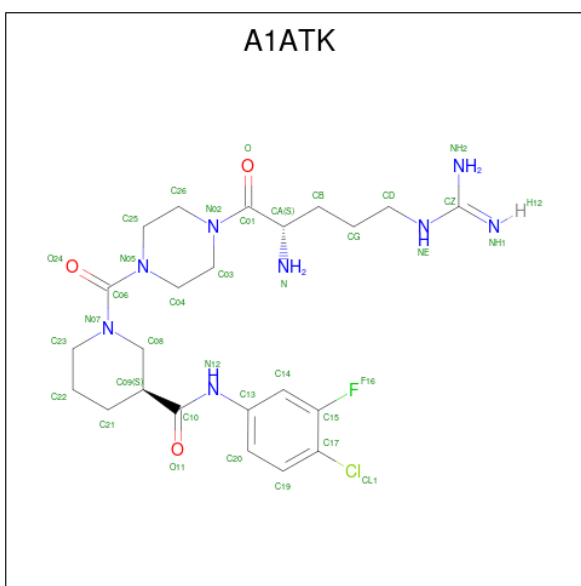
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C<sub>8</sub>H<sub>18</sub>N<sub>2</sub>O<sub>4</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

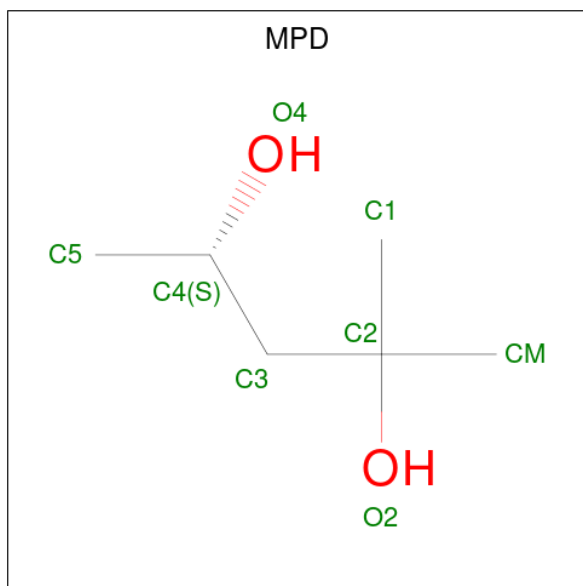
- Molecule 4 is (3S)-1-(4-L-arginylpiperazine-1-carbonyl)-N-(4-chloro-3-fluorophenyl)piperidine-3-carboxamide (three-letter code: A1ATK) (formula:  $C_{23}H_{34}ClFN_8O_3$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	Cl	F	N	O	0	0
			36	23	1	1	8	3		

- Molecule 5 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula:

C<sub>6</sub>H<sub>14</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			8	6	2		

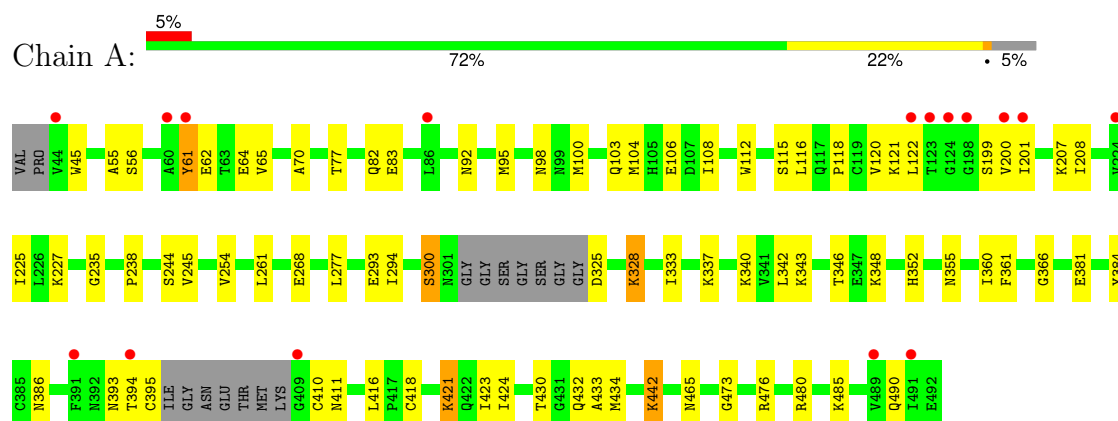
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	180	Total	O	0	0
			180	180		

### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: HIV-1 LM/HS clade A/E CRF01 gp120 core



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.01Å 65.84Å 90.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.43 – 1.85 27.43 – 1.85	Depositor EDS
% Data completeness (in resolution range)	92.3 (27.43-1.85) 92.3 (27.43-1.85)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.14 (at 1.85Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, $R_{free}$	0.230 , 0.259 0.230 , 0.256	Depositor DCC
$R_{free}$ test set	1584 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	36.0	Xtriage
Anisotropy	0.313	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 23.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3024	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.90% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: MPD, EPE, NAG, A1ATK

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	0.33	0/2715	0.57	0/3684

There are no bond length outliers.

There are no bond angle outliers.

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	2659	0	2582	55	0
2	A	126	0	117	2	0
3	A	15	0	17	0	0
4	A	36	0	0	1	0
5	A	8	0	14	2	0
6	A	180	0	0	5	0
All	All	3024	0	2730	57	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 10.

All (57) 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:300:SER:H	1:A:442:LYS:HZ3	1.30	0.79
1:A:65:VAL:HB	1:A:115:SER:HB3	1.66	0.78
1:A:62:GLU:HG3	1:A:64:GLU:H	1.54	0.71
1:A:325:ASP:N	6:A:603:HOH:O	2.24	0.69
1:A:118:PRO:HG2	1:A:121:LYS:HG2	1.76	0.67
1:A:366:GLY:N	6:A:604:HOH:O	2.26	0.67
1:A:65:VAL:HG11	1:A:208:ILE:HD12	1.80	0.64
1:A:207:LYS:NZ	1:A:381:GLU:OE2	2.28	0.63
1:A:337:LYS:NZ	6:A:606:HOH:O	2.31	0.63
1:A:82:GLN:NE2	1:A:83:GLU:O	2.32	0.62
1:A:343:LYS:O	1:A:346:THR:OG1	2.20	0.59
1:A:346:THR:HG22	1:A:361:PHE:HE2	1.68	0.58
1:A:92:ASN:HD22	1:A:238:PRO:HA	1.67	0.58
1:A:340:LYS:HD2	1:A:343:LYS:NZ	2.18	0.58
2:A:504:NAG:H61	5:A:512:MPD:HM2	1.85	0.57
1:A:98:ASN:OD1	1:A:100:MET:HG3	2.05	0.56
1:A:122:LEU:HD23	1:A:200:VAL:HG12	1.87	0.56
1:A:227:LYS:HE3	1:A:485:LYS:HE3	1.88	0.55
1:A:116:LEU:HD13	1:A:208:ILE:HD11	1.90	0.52
1:A:328:LYS:HE2	1:A:418:CYS:O	2.09	0.52
1:A:268:GLU:HB2	6:A:629:HOH:O	2.09	0.52
1:A:361:PHE:O	1:A:393:ASN:ND2	2.40	0.51
1:A:118:PRO:HB3	1:A:433:ALA:HB1	1.92	0.50
1:A:360:ILE:HD12	1:A:465:ASN:HB3	1.92	0.50
1:A:340:LYS:HD2	1:A:343:LYS:HZ1	1.76	0.50
1:A:55:ALA:HB1	1:A:77:THR:HA	1.94	0.49
1:A:103:GLN:HA	1:A:106:GLU:HG2	1.96	0.47
1:A:476:ARG:O	1:A:480:ARG:HB2	2.14	0.47
1:A:348:LYS:HD3	1:A:348:LYS:HA	1.52	0.47
1:A:386:ASN:O	1:A:416:LEU:HD22	2.14	0.47
1:A:56:SER:OG	1:A:70:ALA:HB1	2.14	0.47
1:A:95:MET:SD	1:A:235:GLY:HA3	2.55	0.47
1:A:104:MET:O	1:A:108:ILE:HG12	2.15	0.46
1:A:423:ILE:HG12	1:A:434:MET:HG3	1.98	0.46
1:A:112:TRP:CZ3	1:A:116:LEU:HD23	2.51	0.46
1:A:112:TRP:CE3	1:A:116:LEU:HD23	2.52	0.45
1:A:293:GLU:HB2	6:A:606:HOH:O	2.17	0.45
1:A:384:TYR:CE1	1:A:421:LYS:HB2	2.52	0.45
1:A:120:VAL:HA	1:A:201:ILE:O	2.16	0.44
1:A:95:MET:HE2	1:A:480:ARG:HG2	2.00	0.44
1:A:410:CYS:SG	1:A:410:CYS:O	2.75	0.44
1:A:343:LYS:HZ2	1:A:343:LYS:HB3	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:442:LYS:HB2	1:A:442:LYS:HE2	1.47	0.44
2:A:504:NAG:H61	5:A:512:MPD:CM	2.48	0.43
1:A:294:ILE:HD12	1:A:333:ILE:HD11	1.99	0.43
1:A:480:ARG:HH11	1:A:480:ARG:HG3	1.82	0.43
1:A:225:ILE:HB	1:A:245:VAL:HG23	2.00	0.43
1:A:95:MET:CE	1:A:480:ARG:HG2	2.49	0.43
1:A:95:MET:HG2	1:A:235:GLY:O	2.19	0.42
1:A:254:VAL:HG11	1:A:261:LEU:HB2	2.01	0.42
1:A:424:ILE:O	1:A:432:GLN:HA	2.20	0.41
1:A:65:VAL:CB	1:A:115:SER:HB3	2.43	0.40
1:A:277:LEU:CD1	1:A:352:HIS:HB3	2.51	0.40
1:A:342:LEU:HD12	1:A:342:LEU:HA	1.88	0.40
1:A:293:GLU:H	1:A:337:LYS:NZ	2.19	0.40
1:A:473:GLY:HA2	4:A:511:A1ATK:O11	2.21	0.40
1:A:45:TRP:HA	1:A:490:GLN:O	2.21	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	333/355 (94%)	319 (96%)	12 (4%)	2 (1%)	22	10

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	411	ASN
1	A	61	TYR

### 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	303/313 (97%)	292 (96%)	11 (4%)	30 15

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	TYR
1	A	199	SER
1	A	244	SER
1	A	300	SER
1	A	328	LYS
1	A	355	ASN
1	A	394	THR
1	A	395	CYS
1	A	421	LYS
1	A	430	THR
1	A	442	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	ASN
1	A	352	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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

12 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
2	NAG	A	508	1	14,14,15	0.40	0	17,19,21	0.42	0
2	NAG	A	505	1	14,14,15	0.29	0	17,19,21	0.57	0
2	NAG	A	501	1	14,14,15	0.18	0	17,19,21	0.45	0
2	NAG	A	506	1	14,14,15	0.23	0	17,19,21	0.61	1 (5%)
2	NAG	A	509	1	14,14,15	0.40	0	17,19,21	0.56	0
2	NAG	A	504	1	14,14,15	0.25	0	17,19,21	0.60	0
3	EPE	A	510	-	15,15,15	0.69	1 (6%)	19,20,20	1.71	4 (21%)
2	NAG	A	502	1	14,14,15	0.75	1 (7%)	17,19,21	0.57	0
5	MPD	A	512	-	7,7,7	0.39	0	9,10,10	0.53	0
2	NAG	A	503	1	14,14,15	0.23	0	17,19,21	0.59	0
2	NAG	A	507	1	14,14,15	0.43	0	17,19,21	0.52	0
4	A1ATK	A	511	-	38,38,38	3.31	9 (23%)	50,52,52	2.29	16 (32%)

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	508	1	-	2/6/23/26	0/1/1/1
2	NAG	A	505	1	-	0/6/23/26	0/1/1/1
2	NAG	A	501	1	-	2/6/23/26	0/1/1/1
2	NAG	A	506	1	-	0/6/23/26	0/1/1/1
2	NAG	A	509	1	-	2/6/23/26	0/1/1/1
2	NAG	A	504	1	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EPE	A	510	-	-	2/9/19/19	0/1/1/1
2	NAG	A	502	1	-	0/6/23/26	0/1/1/1
5	MPD	A	512	-	-	2/5/5/5	-
2	NAG	A	503	1	-	0/6/23/26	0/1/1/1
2	NAG	A	507	1	-	2/6/23/26	0/1/1/1
4	A1ATK	A	511	-	-	3/31/51/51	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	511	A1ATK	C01-N02	10.41	1.50	1.35
4	A	511	A1ATK	CZ-NE	9.23	1.51	1.33
4	A	511	A1ATK	C06-N05	7.78	1.52	1.35
4	A	511	A1ATK	C06-N07	7.17	1.51	1.35
4	A	511	A1ATK	C10-N12	6.37	1.50	1.35
4	A	511	A1ATK	C08-N07	3.68	1.51	1.46
4	A	511	A1ATK	C09-C10	2.93	1.56	1.51
4	A	511	A1ATK	O11-C10	-2.88	1.17	1.23
3	A	510	EPE	C10-S	2.22	1.80	1.77
2	A	502	NAG	O5-C1	-2.11	1.40	1.43
4	A	511	A1ATK	C22-C21	-2.06	1.48	1.53

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	511	A1ATK	CA-C01-N02	6.81	129.10	118.85
4	A	511	A1ATK	O-C01-N02	-6.16	114.25	121.61
4	A	511	A1ATK	CB-CA-C01	5.59	121.74	110.75
3	A	510	EPE	C5-N4-C3	4.81	119.21	108.84
4	A	511	A1ATK	O24-C06-N07	-4.09	115.32	124.05
4	A	511	A1ATK	O24-C06-N05	-3.86	115.80	124.05
4	A	511	A1ATK	C04-C03-N02	3.59	117.56	110.42
4	A	511	A1ATK	C03-C04-N05	3.50	117.37	110.42
4	A	511	A1ATK	C21-C09-C10	-3.15	104.16	110.79
3	A	510	EPE	C7-N4-C3	2.99	119.20	111.24
4	A	511	A1ATK	C08-C09-C10	2.79	115.31	110.07
4	A	511	A1ATK	CB-CA-N	-2.75	102.96	110.12
4	A	511	A1ATK	C09-C08-N07	-2.53	105.13	110.14
4	A	511	A1ATK	C04-N05-C06	2.19	130.42	121.53
3	A	510	EPE	O3S-S-C10	2.17	110.25	106.00
3	A	510	EPE	C5-C6-N1	-2.14	106.32	110.65

*Continued on next page...*

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	511	A1ATK	C01-CA-N	-2.12	101.47	109.20
4	A	511	A1ATK	C08-N07-C06	2.11	130.07	121.53
2	A	506	NAG	C1-O5-C5	2.05	114.94	112.19
4	A	511	A1ATK	C03-N02-C01	2.05	130.97	123.30
4	A	511	A1ATK	C19-C17-C15	2.04	120.79	119.04

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	509	NAG	O5-C5-C6-O6
2	A	509	NAG	C4-C5-C6-O6
4	A	511	A1ATK	NE-CD-CG-CB
2	A	501	NAG	O5-C5-C6-O6
2	A	501	NAG	C4-C5-C6-O6
5	A	512	MPD	C2-C3-C4-C5
3	A	510	EPE	C10-C9-N1-C2
3	A	510	EPE	C10-C9-N1-C6
2	A	508	NAG	O5-C5-C6-O6
2	A	508	NAG	C4-C5-C6-O6
4	A	511	A1ATK	CG-CD-NE-CZ
2	A	504	NAG	C4-C5-C6-O6
2	A	507	NAG	O5-C5-C6-O6
5	A	512	MPD	C2-C3-C4-O4
2	A	507	NAG	C4-C5-C6-O6
4	A	511	A1ATK	N-CA-CB-CG

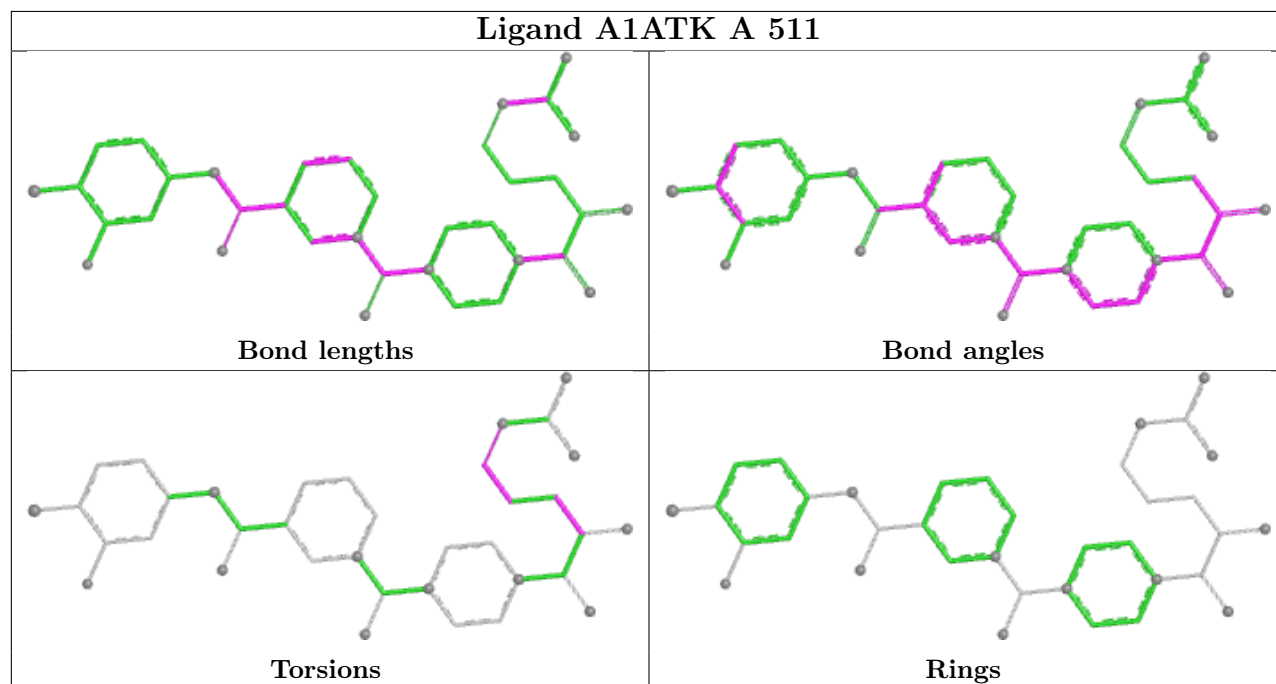
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	504	NAG	2	0
5	A	512	MPD	2	0
4	A	511	A1ATK	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.



## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	339/355 (95%)	0.56	16 (4%) 37 39	24, 39, 60, 81	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	200	VAL	3.6
1	A	122	LEU	3.5
1	A	123	THR	3.5
1	A	391	PHE	3.2
1	A	224	VAL	3.0
1	A	409	GLY	2.9
1	A	124	GLY	2.8
1	A	394	THR	2.6
1	A	60	ALA	2.5
1	A	198	GLY	2.4
1	A	44	VAL	2.3
1	A	201	ILE	2.2
1	A	61	TYR	2.2
1	A	491	ILE	2.2
1	A	489	VAL	2.1
1	A	86	LEU	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 monosaccharides in this entry.

## 6.4 Ligands

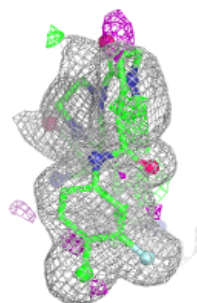
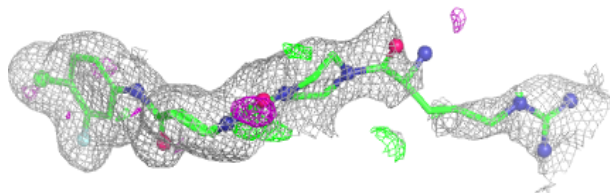
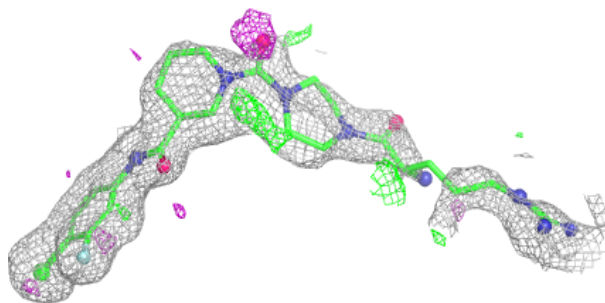
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	NAG	A	502	14/15	0.69	0.15	63,71,77,77	0
2	NAG	A	509	14/15	0.73	0.10	51,60,67,74	0
5	MPD	A	512	8/8	0.74	0.15	60,64,74,77	0
2	NAG	A	506	14/15	0.76	0.11	48,52,60,68	0
2	NAG	A	504	14/15	0.80	0.10	49,57,66,69	0
2	NAG	A	501	14/15	0.81	0.12	43,46,53,68	0
4	A1ATK	A	511	36/36	0.86	0.14	24,48,67,69	0
2	NAG	A	508	14/15	0.87	0.11	39,46,53,55	0
2	NAG	A	507	14/15	0.87	0.10	49,54,57,59	0
2	NAG	A	505	14/15	0.88	0.09	45,48,53,57	0
2	NAG	A	503	14/15	0.94	0.07	24,32,41,45	0
3	EPE	A	510	15/15	0.94	0.09	30,36,38,41	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 A1ATK A 511:**

$2mF_o-DF_c$  (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.