

Full wwPDB X-ray Structure Validation Report (i)

Aug 28, 2023 – 03:49 AM EDT

PDB ID : 3LIV

Title : crystal structure of HTLV protease complexed with the inhibitor KNI-10683

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Deposited on : 2010-01-25

Resolution : 2.59 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 (i) 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 (1)) were used in the production of this report:

MolProbity : 4.02b-467

Mogul : 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13

EDS: 2.35

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Refmac : 5.8.0158

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

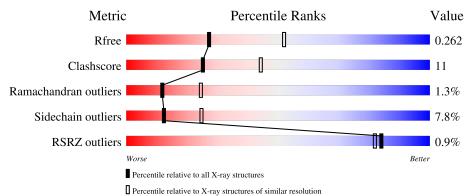
Validation Pipeline (wwPDB-VP) : 2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	Similar resolution		
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$		
R_{free}	130704	3163 (2.60-2.60)		
Clashscore	141614	3518 (2.60-2.60)		
Ramachandran outliers	138981	3455 (2.60-2.60)		
Sidechain outliers	138945	3455 (2.60-2.60)		
RSRZ outliers	127900	3104 (2.60-2.60)		

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 <=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	116	77%	19%	
1	В	116	74%	21%	5%



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 1899 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.

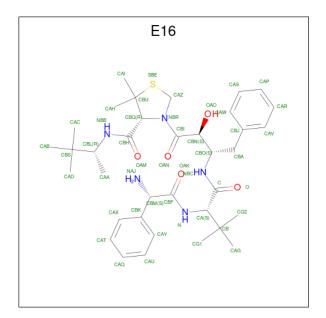
• Molecule 1 is a protein called Protease.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	116	Total 891			O 165	S 3	0	2	0
1	В	116	Total 883			O 164	S 3	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	40	ILE	LEU	engineered mutation	UNP Q82134
В	40	ILE	LEU	engineered mutation	UNP Q82134

• Molecule 2 is (4R)-3-[(2S,3S)-3-[[(2S)-2-[[(2S)-2-azanyl-2-phenyl-ethanoyl]amino]-3,3-dimethylbutanoyl]amino]-2-hydroxy-4-phenyl-but anoyl]-N-[(2R)-3,3-dimethylbutan-2-yl]-5,5-dimethyl-1,3-thiazolidine-4-carboxamide (three-letter code: E16) (formula: $C_{36}H_{53}N_5O_5S$).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
9	B	1	Total	С	N	О	S	0	0
	Ъ	1	47	36	5	5	1	0	0

$\bullet\,$ Molecule 3 is water.

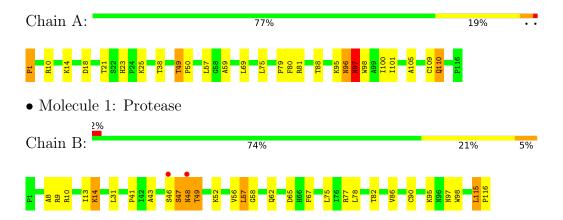
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	39	Total O 39 39	0	0
3	В	39	Total O 39 39	0	0



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: Protease





4 Data and refinement statistics (i)

Property	Value	Source	
Space group	P 63 2 2	Depositor	
Cell constants	77.77Å 77.77Å 162.97Å	Donogitor	
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor	
Resolution (Å)	50.00 - 2.59	Depositor	
rtesolution (A)	42.28 - 2.59	EDS	
% Data completeness	93.8 (50.00-2.59)	Depositor	
(in resolution range)	93.8 (42.28-2.59)	EDS	
R_{merge}	(Not available)	Depositor	
R_{sym}	0.08	Depositor	
$< I/\sigma(I) > 1$	5.48 (at 2.58Å)	Xtriage	
Refinement program	REFMAC 5.4.0057	Depositor	
P. P.	0.201 , 0.262	Depositor	
R, R_{free}	0.193 , 0.262	DCC	
R_{free} test set	268 reflections (2.96%)	wwPDB-VP	
Wilson B-factor (Å ²)	35.7	Xtriage	
Anisotropy	0.345	Xtriage	
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.31 , 31.0	EDS	
L-test for twinning ²	$ < L >=0.51, < L^2>=0.34$	Xtriage	
Estimated twinning fraction	No twinning to report.	Xtriage	
F_o, F_c correlation	0.95	EDS	
Total number of atoms	1899	wwPDB-VP	
Average B, all atoms (Å ²)	40.0	wwPDB-VP	

Xtriage's analysis on translational NCS is as follows: The analyses of the Patterson function reveals a significant off-origin peak that is 34.22 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.0638e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.

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



¹Intensities estimated from amplitudes.

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: E16

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		nd lengths	Bond angles		
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.69	2/915~(0.2%)	0.82	2/1252~(0.2%)	
1	В	0.58	0/901	0.82	1/1233 (0.1%)	
All	All	0.64	2/1816 (0.1%)	0.82	3/2485 (0.1%)	

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(A)
1	A	109	CYS	CB-SG	-6.64	1.71	1.82
1	A	1	PRO	N-CD	5.86	1.56	1.47

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}(^{o})$
1	В	48	ASN	N-CA-C	-5.58	95.95	111.00
1	A	1	PRO	CA-N-CD	-5.07	104.40	111.50
1	A	109	CYS	CB-CA-C	-5.04	100.31	110.40

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	891	0	942	20	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	В	883	0	926	24	0
2	В	47	0	53	1	0
3	A	39	0	0	7	0
3	В	39	0	0	7	0
All	All	1899	0	1921	40	0

The all-atom clash score is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clash score for this structure is 11.

All (40) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${f distance}({ m \AA})$	$-$ overlap (\AA)
1:B:10:ARG:HD3	3:B:120:HOH:O	1.72	0.89
1:B:9:ARG:HD3	3:B:141:HOH:O	1.72	0.86
1:B:46:SER:O	1:B:48:ASN:O	2.00	0.79
1:A:23:HIS:HB3	3:A:131:HOH:O	1.84	0.77
1:B:48:ASN:O	1:B:49:THR:HB	1.96	0.66
1:A:88:THR:HB	3:A:148:HOH:O	1.96	0.65
1:A:80:PHE:HD2	1:A:81:ARG:HG2	1.60	0.65
1:A:1:PRO:HA	3:A:153:HOH:O	1.98	0.64
1:A:1:PRO:N	1:B:115:LEU:O	2.35	0.59
1:A:80:PHE:CD2	1:A:81:ARG:HG2	2.38	0.59
1:A:25:LYS:HE2	3:A:127:HOH:O	2.03	0.57
1:B:14:LYS:HD3	3:B:140:HOH:O	2.04	0.56
1:A:38:THR:HG23	1:A:100:ILE:O	2.07	0.54
1:B:10:ARG:CD	3:B:120:HOH:O	2.40	0.54
1:A:98:TRP:CZ2	2:B:117:E16:HAP	2.43	0.54
1:B:41:PRO:HB3	1:B:97:ASN:HB3	1.90	0.53
1:B:13:ILE:HD13	1:B:78:LEU:HD22	1.94	0.49
1:A:59:ALA:HB3	1:B:58:GLY:HA3	1.94	0.48
1:B:52:LYS:HD3	3:B:145:HOH:O	2.14	0.48
1:A:59:ALA:HA	1:B:98:TRP:HZ3	1.79	0.48
1:B:46:SER:HA	3:B:155:HOH:O	2.14	0.47
1:B:56:VAL:HG21	1:B:67:PHE:HB2	1.97	0.47
1:A:21:THR:CG2	1:A:49:THR:HG21	2.44	0.47
1:A:10:ARG:HD2	3:A:137:HOH:O	2.14	0.47
1:B:41:PRO:CB	1:B:97:ASN:HD22	2.28	0.47
1:B:47:SER:C	1:B:48:ASN:O	2.54	0.45
1:B:48:ASN:O	1:B:49:THR:CB	2.66	0.43
1:A:81:ARG:HG3	3:A:145:HOH:O	2.18	0.42
1:B:8:ALA:HB3	3:B:148:HOH:O	2.20	0.42

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Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:A:14:LYS:O	1:A:79:PRO:HD3	2.19	0.42
3:A:154:HOH:O	1:B:10:ARG:HG3	2.19	0.42
1:B:57:LEU:HG	1:B:62:GLN:NE2	2.35	0.41
1:A:101[A]:ILE:HG23	1:A:105:ALA:HB3	2.02	0.41
1:A:18:ASP:HB3	1:A:75:LEU:HB2	2.03	0.41
1:A:49:THR:HA	1:A:50:PRO:HD3	1.86	0.41
1:A:95:LYS:O	1:A:97:ASN:N	2.49	0.41
1:B:41:PRO:HB3	1:B:97:ASN:HD22	1.86	0.40
1:B:75:LEU:HD22	1:B:86:VAL:HG22	2.02	0.40
1:A:110:GLN:HB2	1:B:116:PRO:HG2	2.02	0.40
1:B:43:ALA:H	1:B:97:ASN:ND2	2.19	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	116/116 (100%)	112 (97%)	2 (2%)	2 (2%)	9 18
1	В	114/116~(98%)	108 (95%)	5 (4%)	1 (1%)	17 35
All	All	230/232~(99%)	220 (96%)	7 (3%)	3 (1%)	12 24

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	96	ASN
1	В	49	THR
1	A	97	ASN



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	104/102 (102%)	98 (94%)	6 (6%)	20	40	
1	В	102/102 (100%)	92 (90%)	10 (10%)	8	15	
All	All	206/204 (101%)	190 (92%)	16 (8%)	12	25	

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	THR
1	A	57	LEU
1	A	69	LEU
1	A	96	ASN
1	A	97	ASN
1	A	110	GLN
1	В	14	LYS
1	В	31	LEU
1	В	47	SER
1	В	57	LEU
1	В	65	ASP
1	В	77	ARG
1	В	82	THR
1	В	90	CYS
1	В	95	LYS
1	В	115	LEU

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

Mol	Chain	Res	Type
1	A	96	ASN
1	A	97	ASN
1	В	23	HIS
1	В	62	GLN
1	В	97	ASN



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 monosaccharides in this entry.

5.6 Ligand geometry (i)

1 ligand is 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 Chai	Chain	Chain	Chain	Chain	Chain	Chain	Chain	Chain Res		B	ond leng	gths	В	ond ang	gles
WIOI	Moi Type Chain	nes	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2						
2	E16	В	117	-	47,49,49	4.78	23 (48%)	64,73,73	1.64	13 (20%)					

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	E16	В	117	-	-	14/56/72/72	0/3/3/3

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	Observed(A)	$Ideal(\AA)$
2	В	117	E16	CBQ-NBR	13.17	1.58	1.46
2	В	117	E16	CAY-CBK	10.53	1.56	1.39
2	В	117	E16	CAX-CBK	10.22	1.55	1.39

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Mol	Chain	Res	Type	Atoms	\mathbf{Z}	Observed(A)	$Ideal(\AA)$
2	В	117	E16	CAU-CAY	7.90	1.55	1.38
2	В	117	E16	CAT-CAX	7.88	1.55	1.38
2	В	117	E16	CAR-CAV	7.17	1.54	1.38
2	В	117	E16	CAS-CAW	6.85	1.53	1.38
2	В	117	E16	CAU-CAQ	6.71	1.55	1.38
2	В	117	E16	CAT-CAQ	6.61	1.55	1.38
2	В	117	E16	CAW-CBJ	6.55	1.53	1.38
2	В	117	E16	CAS-CAP	6.18	1.54	1.38
2	В	117	E16	CAZ-SBE	-6.15	1.68	1.83
2	В	117	E16	CAR-CAP	6.09	1.54	1.38
2	В	117	E16	CAV-CBJ	5.94	1.51	1.38
2	В	117	E16	CBF-N	5.66	1.46	1.34
2	В	117	E16	C-NBC	5.42	1.46	1.34
2	В	117	E16	CBH-NBB	5.16	1.45	1.34
2	В	117	E16	CAZ-NBR	4.80	1.55	1.46
2	В	117	E16	CBI-NBR	4.16	1.44	1.34
2	В	117	E16	CBK-CBM	3.42	1.56	1.52
2	В	117	E16	CBU-SBE	-3.29	1.78	1.85
2	В	117	E16	OAN-CBI	-2.82	1.17	1.22
2	В	117	E16	OAM-CBH	-2.41	1.18	1.23

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
2	В	117	E16	CBN-CBO-NBC	-5.36	99.95	110.03
2	В	117	E16	CBJ-CBA-CBO	-4.50	105.52	113.33
2	В	117	E16	CBM-CBF-N	3.27	118.76	114.78
2	В	117	E16	CAA-CBL-NBB	-3.06	103.17	109.25
2	В	117	E16	CAD-CBS-CAB	-2.97	103.03	108.80
2	В	117	E16	CAI-CBU-CBQ	-2.75	105.47	111.57
2	В	117	E16	CBO-NBC-C	-2.72	118.27	123.07
2	В	117	E16	CA-N-CBF	-2.46	116.97	122.05
2	В	117	E16	CBQ-CBH-NBB	2.19	118.87	114.48
2	В	117	E16	CBS-CBL-NBB	2.17	115.14	110.60
2	В	117	E16	CAI-CBU-CAH	-2.11	107.34	110.78
2	В	117	E16	CAH-CBU-CBQ	2.06	116.15	111.57
2	В	117	E16	CAW-CBJ-CAV	2.04	121.37	118.17

There are no chirality outliers.

All (14) torsion outliers are listed below:



Mol	Chain	Res	Type	Atoms
2	В	117	E16	O-C-CA-CB
2	В	117	E16	NBC-C-CA-CB
2	В	117	E16	CBS-CBL-NBB-CBH
2	В	117	E16	CAA-CBL-CBS-CAD
2	В	117	E16	CAA-CBL-CBS-CAC
2	В	117	E16	CAA-CBL-NBB-CBH
2	В	117	E16	NBB-CBL-CBS-CAD
2	В	117	E16	N-CA-CB-CG1
2	В	117	E16	CBI-CBN-CBO-NBC
2	В	117	E16	N-CA-CB-CAG
2	В	117	E16	C-CA-CB-CG1
2	В	117	E16	N-CA-CB-CG2
2	В	117	E16	C-CA-CB-CAG
2	В	117	E16	C-CA-CB-CG2

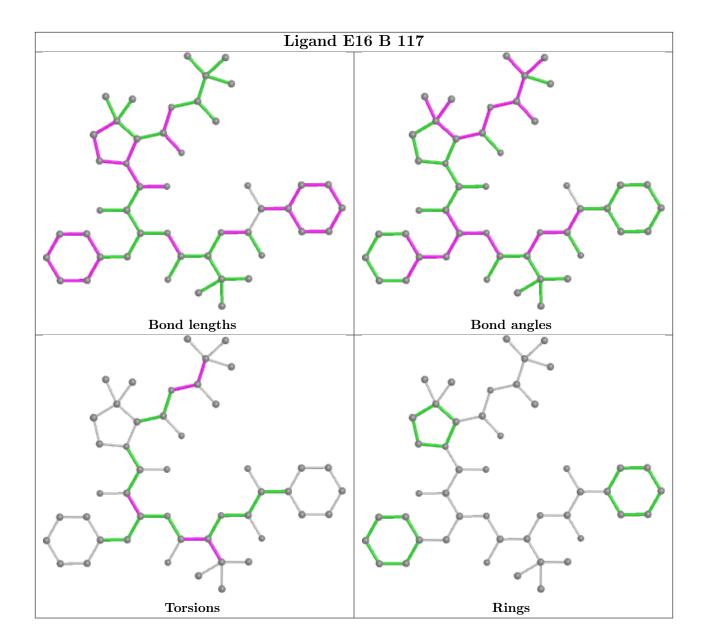
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	В	117	E16	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 then 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^{th} 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	$OWAB(A^2)$	Q < 0.9
1	A	116/116 (100%)	-0.36	0 100 100	19, 36, 66, 85	0
1	В	116/116 (100%)	-0.31	2 (1%) 70 66	18, 36, 69, 85	0
All	All	232/232 (100%)	-0.33	2 (0%) 84 82	18, 36, 66, 85	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	48	ASN	3.9
1	В	46	SER	2.4

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 (i)

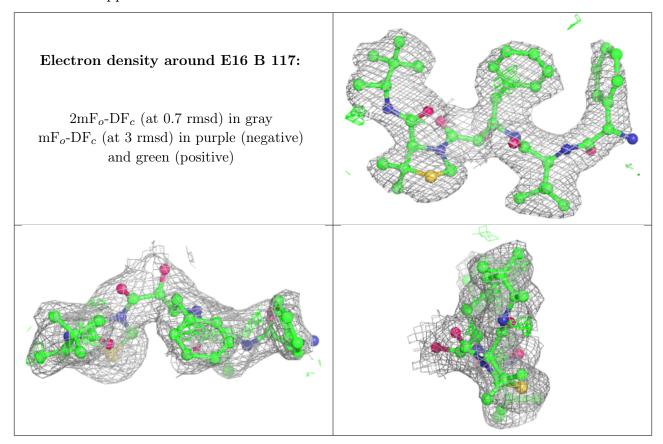
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	E16	В	117	47/47	0.96	0.17	20,26,55,58	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.

