

Full wwPDB X-ray Structure Validation Report (i)

May 23, 2020 – 10:53 pm BST

PDB ID : 5OGR

Title : Structure of cathepsin B1 from Schistosoma mansoni in complex with WRR286

inhibitor

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Deposited on : 2017-07-13

Resolution : 1.55 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 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 : 2.11

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)

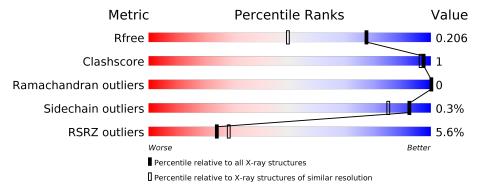
Validation Pipeline (wwPDB-VP) : 2.11

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 1.55 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar resolution} \\ (\#{\rm Entries, resolution range(\AA)}) \end{array}$
R_{free}	130704	1483 (1.56-1.56)
Clashscore	141614	1529 (1.56-1.56)
Ramachandran outliers	138981	1498 (1.56-1.56)
Sidechain outliers	138945	1495 (1.56-1.56)
RSRZ outliers	127900	1465 (1.56-1.56)

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 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	254	96%	•				
1	В	254	96%					
1	С	254	98%	•				



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 7148 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 Cathepsin B-like peptidase (C01 family).

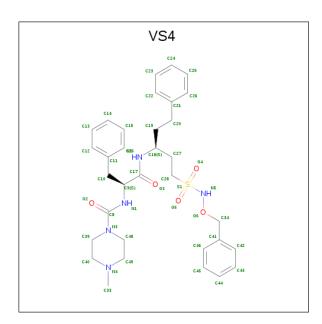
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	Λ	254	Total	С	N	О	S	0	14	0
1	1 A	204	2077	1309	365	386	17	U	14	
1	D	254	Total	С	N	О	S	0	7	0
1		204	2035	1278	356	384	17	U		
1	C	C 254	Total	С	N	О	S	0	9	0
1			2008	1261	352	378	17	0	2	U

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	168	ALA	THR	engineered mutation	UNP Q8MNY2
A	283	ALA	THR	engineered mutation	- 1
В	168	ALA	THR	engineered mutation	UNP Q8MNY2
В	283	ALA	THR	engineered mutation	UNP Q8MNY2
С	168	ALA	THR	engineered mutation	
С	283	ALA	THR	engineered mutation	UNP Q8MNY2

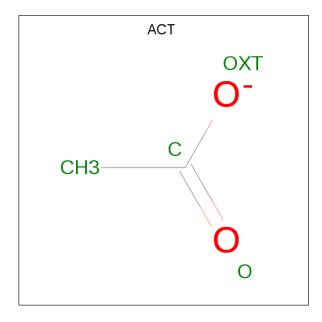
• Molecule 2 is 3-[[N-[4-METHYL-PIPERAZINYL]CARBONYL]-PHENYLALANINYL-A MINO]-5-PHENYL-PENTANE-1-SULFONIC ACID BENZYLOXY-AMIDE (three-letter code: VS4) (formula: $C_{33}H_{43}N_5O_5S$).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf		
9	2 A	1	Total	С	N	О	S	0	0	
		1	44	33	5	5	1	0	0	
9	B	1	Total	С	N	О	S	0	0	
	2 D	1	44	33	5	5	1			
9	С	1	Total	С	N	О	S	0	0	
			44	33	5	5	1	U		

 \bullet Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $\mathrm{C_2H_3O_2}).$



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



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\mathbf{Mol}	Chain	Residues	Atoms	ZeroOcc	AltConf
3	В	1	Total C O 4 2 2	0	0
3	С	1	Total C O 4 2 2	0	0

• Molecule 4 is water.

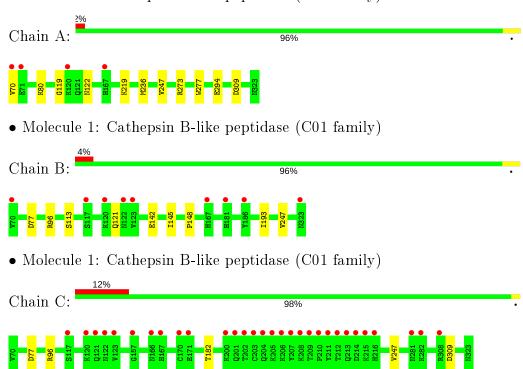
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	356	Total O 362 362	0	13
4	В	313	Total O 318 318	0	11
4	С	204	Total O 204 204	0	1



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: Cathepsin B-like peptidase (C01 family)





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	65.48	D : 4
a, b, c, α , β , γ	90.00° 91.24° 90.00°	Depositor
Resolution (Å)	39.04 - 1.55	Depositor
Resolution (A)	39.04 - 1.55	EDS
% Data completeness	98.9 (39.04-1.55)	Depositor
(in resolution range)	98.9 (39.04-1.55)	EDS
R_{merge}	0.06	Depositor
R_{sum}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.46 (at 1.55Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
D D.	0.162 , 0.203	Depositor
R, R_{free}	0.173 , 0.206	DCC
R_{free} test set	1121 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å ²)	18.6	Xtriage
Anisotropy	0.133	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.34 , 45.7	EDS
L-test for twinning ²	$< L >=0.48, < L^2>=0.31$	Xtriage
	0.020 for -h,-l,-k	
Estimated twinning fraction	0.006 for -h,l,k	Xtriage
	0.026 for h,-k,-l	
F_o, F_c correlation	0.97	EDS
Total number of atoms	7148	wwPDB-VP
Average B, all atoms (\mathring{A}^2)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.24% of the height of the origin peak. No significant pseudotranslation is detected.

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: VS4, ACT

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	Boı	nd lengths	Bond angles		
Mol Chain		RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.79	$1/2171 \ (0.0\%)$	0.83	$2/2921 \ (0.1\%)$	
1	В	0.79	0/2105	0.83	3/2836 (0.1%)	
1	С	0.70	0/2066	0.77	3/2784 (0.1%)	
All	All	0.76	$1/6342 \ (0.0\%)$	0.81	8/8541 (0.1%)	

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(ext{\AA})$
1	A	294	GLU	CD-OE1	5.70	1.31	1.25

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
1	В	96	ARG	NE-CZ-NH1	6.94	123.77	120.30
1	В	77	ASP	CB-CG-OD1	6.71	124.34	118.30
1	С	77	ASP	CB-CG-OD1	5.89	123.60	118.30
1	A	273	ARG	NE-CZ-NH2	-5.70	117.45	120.30
1	С	96	ARG	NE-CZ-NH1	5.67	123.14	120.30
1	С	309	ASP	CB-CG-OD1	5.54	123.29	118.30
1	В	77	ASP	CB-CG-OD2	-5.17	113.64	118.30
1	A	309	ASP	CB-CG-OD1	5.02	122.82	118.30

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	2077	0	2042	5	0
1	В	2035	0	1967	5	0
1	С	2008	0	1932	1	0
2	A	44	0	42	1	0
2	В	44	0	42	2	0
2	С	44	0	42	1	0
3	A	4	0	3	0	0
3	В	4	0	3	0	0
3	С	4	0	3	0	0
4	A	362	0	0	1	0
4	В	318	0	0	0	0
4	С	204	0	0	0	0
All	All	7148	0	6076	10	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 1.

All (10) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{aligned} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{aligned}$	$egin{array}{c} ext{Clash} \ ext{overlap } (ext{Å}) \end{array}$
1:B:247:VAL:HG21	2:B:401:VS4:H341	1.74	0.70
1:C:247:VAL:HG21	2:C:401:VS4:H341	1.76	0.66
1:A:247:VAL:HG21	2:A:401:VS4:H341	1.82	0.61
1:A:236[A]:MET:HG3	1:A:277:TRP:HH2	1.73	0.54
1:A:119:GLY:HA2	1:A:122:ASN:OD1	2.10	0.52
1:B:113[B]:SER:OG	1:B:121:GLN:HB3	2.13	0.48
1:B:193:ILE:CD1	2:B:401:VS4:H251	2.47	0.45
1:B:145:ILE:HG22	1:B:148:PRO:HD2	2.03	0.41
1:A:80:LYS:HE2	1:B:142:GLU:HB3	2.02	0.40
1:A:219:LYS:NZ	4:A:510:HOH:O	2.51	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	$266/254 \ (105\%)$	256 (96%)	10 (4%)	0	100	100	
1	В	$259/254 \ (102\%)$	250 (96%)	9 (4%)	0	100	100	
1	С	$254/254 \ (100\%)$	242 (95%)	12 (5%)	0	100	100	
All	All	779/762 (102%)	748 (96%)	31 (4%)	0	100	100	

There are no Ramachandran outliers to report.

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	$228/214 \ (106\%)$	227 (100%)	1 (0%)	91	82	
1	В	221/214 (103%)	221 (100%)	0	100	100	
1	С	216/214 (101%)	215 (100%)	1 (0%)	88	78	
All	All	665/642 (104%)	663 (100%)	2 (0%)	92	85	

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	${f Res}$	\mathbf{Type}
1	A	70	VAL
1	С	182	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

6 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).

Mal	T	Chain	Dag	Link	В	ond leng	gths	Bond angles		
Mol	Type	Chain	Res	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ACT	A	402	-	1,3,3	2.57	1 (100%)	0,3,3	0.00	=
3	ACT	С	402	-	1,3,3	1.62	0	0,3,3	0.00	-
2	VS4	A	401	1	47,47,47	1.48	6 (12%)	55,62,62	1.75	6 (10%)
3	ACT	В	402	-	1,3,3	1.17	0	0,3,3	0.00	-
2	VS4	С	401	1	47,47,47	1.46	5 (10%)	55,62,62	1.70	11 (20%)
2	VS4	В	401	1	47,47,47	1.54	7 (14%)	55,62,62	1.27	5 (9%)

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	\mathbf{Res}	Link	Chirals	${f Torsions}$	Rings
2	VS4	A	401	1	-	2/34/48/48	0/4/4/4
2	VS4	С	401	1	-	5/34/48/48	0/4/4/4
2	VS4	В	401	1	-	4/34/48/48	0/4/4/4

All (19) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	$Ideal(\AA)$
2	В	401	VS4	O4-S1	5.95	1.51	1.43
2	С	401	VS4	O4-S1	5.10	1.50	1.43
2	A	401	VS4	O4-S1	4.03	1.49	1.43
2	В	401	VS4	C49-N4	3.88	1.54	1.46
2	С	401	VS4	C49-N4	3.50	1.53	1.46
2	A	401	VS4	C27-C28	3.44	1.56	1.52
2	A	401	VS4	C49-N4	3.26	1.53	1.46
2	С	401	VS4	C8-N1	2.75	1.41	1.35
2	В	401	VS4	O6-C34	2.65	1.49	1.43
3	A	402	ACT	СН3-С	2.57	1.52	1.48
2	В	401	VS4	C8-N1	2.56	1.40	1.35
2	A	401	VS4	C8-N3	2.46	1.41	1.36
2	С	401	VS4	O6-N5	-2.44	1.39	1.42
2	A	401	VS4	C8-N1	2.38	1.40	1.35
2	A	401	VS4	O6-C34	2.35	1.49	1.43
2	В	401	VS4	C48-N3	-2.17	1.43	1.47
2	С	401	VS4	C48-N3	-2.12	1.43	1.47
2	В	401	VS4	C8-N3	2.06	1.40	1.36
2	В	401	VS4	S1-N5	2.02	1.68	1.62

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	A	401	VS4	O5-S1-O4	-8.87	106.50	119.35
2	С	401	VS4	C33-N4-C40	5.85	119.41	110.66
2	С	401	VS4	C48-C49-N4	-3.95	106.34	110.80
2	В	401	VS4	O6-C34-C41	3.50	124.59	109.52
2	A	401	VS4	O4-S1-C28	3.46	113.38	107.86
2	С	401	VS4	O5-S1-O4	-3.32	114.53	119.35
2	С	401	VS4	O5-S1-C28	-3.16	102.82	107.86
2	С	401	VS4	C27-C18-N2	-3.12	106.15	110.54
2	A	401	VS4	O6-C34-C41	3.06	122.72	109.52
2	В	401	VS4	O5-S1-O4	-2.88	115.18	119.35
2	С	401	VS4	O6-C34-C41	2.87	121.87	109.52
2	В	401	VS4	N1-C8-N3	-2.82	112.07	117.21
2	A	401	VS4	C24-C25-C26	-2.82	115.90	120.19
2	A	401	VS4	C48-N3-C39	2.65	117.72	112.62
2	С	401	VS4	O2-C8-N3	2.64	125.49	121.78
2	С	401	VS4	C14-C15-C16	-2.58	116.27	120.19
2	В	401	VS4	C48-C49-N4	-2.45	108.04	110.80
2	В	401	VS4	C39-N3-C8	-2.36	113.26	121.94
2	С	401	VS4	C33-N4-C49	-2.20	107.36	110.66
2	С	401	VS4	C39-N3-C8	-2.15	114.01	121.94



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Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
2	A	401	VS4	C23-C22-C21	-2.11	117.39	120.63
2	С	401	VS4	C48-N3-C39	2.01	116.49	112.62

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	VS4	C27-C28-S1-N5
2	С	401	VS4	C27-C28-S1-N5
2	С	401	VS4	C27-C28-S1-O4
2	С	401	VS4	C27-C28-S1-O5
2	В	401	VS4	C27-C28-S1-N5
2	В	401	VS4	C27-C28-S1-O4
2	В	401	VS4	C41-C34-O6-N5
2	A	401	VS4	C27-C28-S1-O5
2	С	401	VS4	C41-C34-O6-N5
2	С	401	VS4	C19-C18-C27-C28
2	В	401	VS4	C19-C18-C27-C28

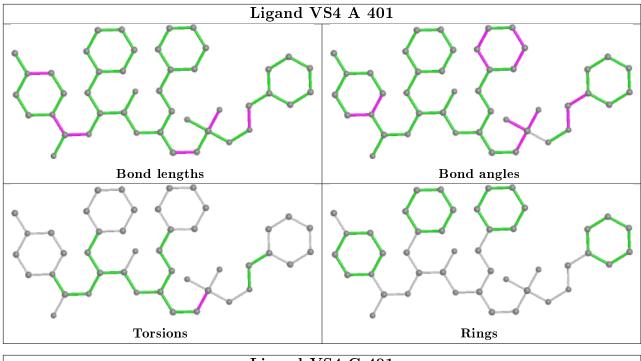
There are no ring outliers.

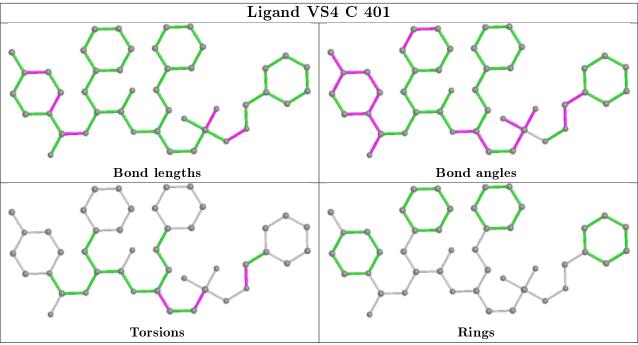
3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	VS4	1	0
2	С	401	VS4	1	0
2	В	401	VS4	2	0

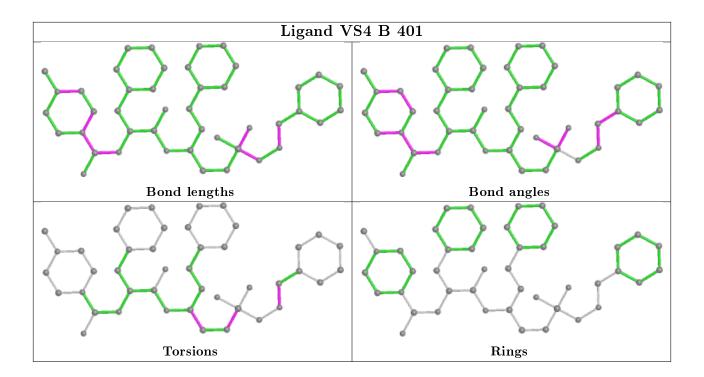
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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 $>$	$\#\mathrm{RSRZ}{>}2$	$\mathbf{OWAB}(\mathrm{\AA}^2)$	Q < 0.9
1	A	$254/254 \ (100\%)$	-0.07	4 (1%) 72 77	13, 19, 36, 52	0
1	В	$254/254 \ (100\%)$	0.03	9 (3%) 44 52	14, 20, 42, 71	0
1	С	254/254 (100%)	0.57	30 (11%) 4 4	15, 29, 69, 82	0
All	All	762/762 (100%)	0.18	43 (5%) 24 28	13, 22, 59, 82	0

All (43) RSRZ outliers are listed below:

Mol	Chain	V 1		RSRZ
1	С	210	PRO	7.5
1	С	123	VAL	7.4
1	С	211	TYR	7.0
1	С	208	LYS	6.6
1	С	212	THR	6.0
1	С	206	LYS	5.9
1	С	209	THR	5.8
1	С	213	GLN	5.6
1	С	202	THR	5.1
1	С	167	HIS	4.7
1	С	203	CYS	4.6
1	С	121	GLN	4.6
1	С	120	LYS	4.4
1	С	122	ASN	3.9
1	A	70	VAL	3.9
1	С	170	CYS	3.8
1	С	207	TYR	3.3
1	В	120	LYS	3.2
1	С	205	LYS	3.1
1	С	281	ASN	3.0
1	В	70	VAL	2.9
1	В	181	HIS	2.8
1	В	323	ASN	2.7



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Mol	Chain	Res	Type	RSRZ	
1	В	167	HIS	2.6	
1	С	201	GLN	2.6	
1	С	216	HIS	2.5	
1	С	117	SER	2.5	
1	A	71	GLU	2.5	
1	С	214	ASP	2.4	
1	В	123	VAL	2.4	
1	В	186	TYR	2.3	
1	С	204	GLN	2.3	
1	A	167	HIS	2.3	
1	С	157	GLY	2.3	
1	С	282	LYS	2.2	
1	С	166	ASN	2.2	
1	С	200	LYS	2.2	
1	С	171	GLU	2.2	
1	С	308	ARG	2.2	
1	A	120	LYS	2.1	
1	В	122	ASN	2.1	
1	В	117	SER	2.0	
1	С	215	LYS	2.0	

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

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

6.3 Carbohydrates (i)

There are no carbohydrates 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.

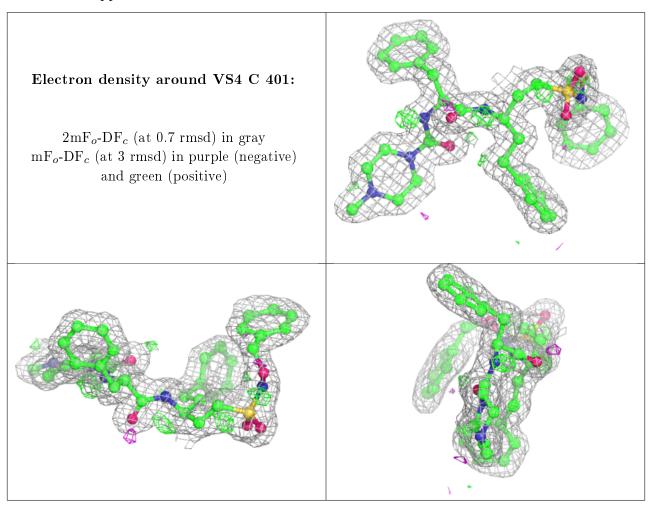
Mol	Type	Chain	Res	Atoms	RSCC	RSR	${f B-factors}({f \AA}^2)$	Q<0.9
3	ACT	С	402	4/4	0.91	0.09	20,24,24,26	0
2	VS4	С	401	44/44	0.95	0.11	15,21,34,35	0
2	VS4	A	401	44/44	0.96	0.11	13,15,31,35	0



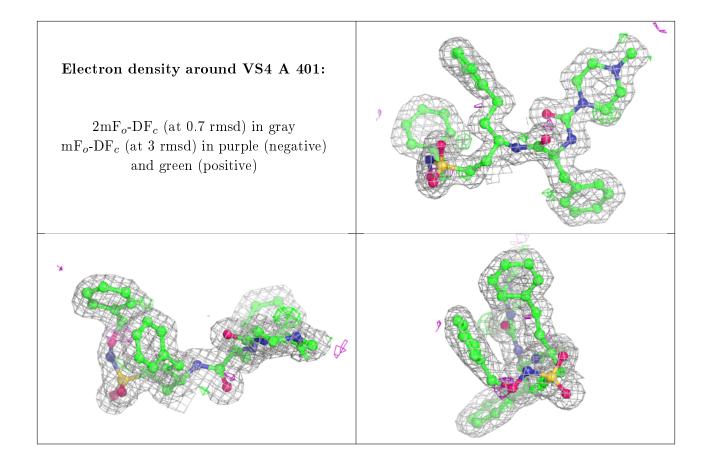
 $Continued\ from\ previous\ page...$

Mol	Type	Chain	Res	Atoms	RSCC	RSR	${f B\text{-factors}}({f \AA}^2)$	Q<0.9
3	ACT	В	402	4/4	0.96	0.05	20,25,25,25	0
3	ACT	A	402	4/4	0.96	0.10	19,25,26,26	0
2	VS4	В	401	44/44	0.96	0.09	11,13,32,33	0

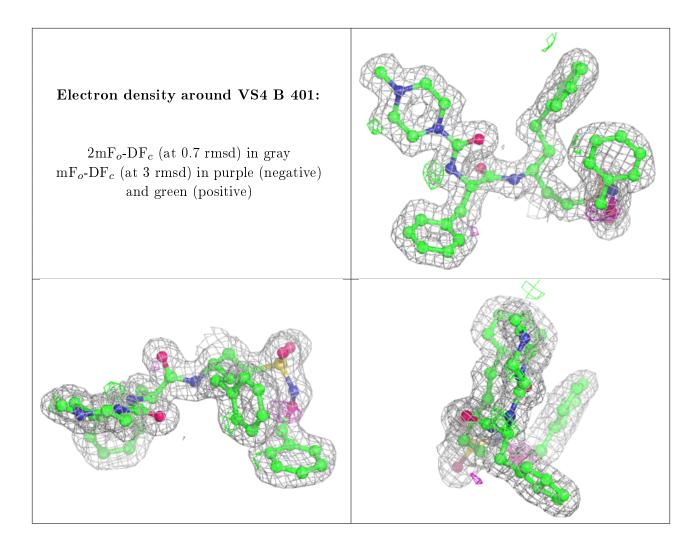
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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











6.5 Other polymers (i)

There are no such residues in this entry.

