

Full wwPDB X-ray Structure Validation Report (i)

Oct 31, 2023 - 03:12 PM JST

PDB ID	:	8HCE
Title	:	Crystal structure of mTREX1-CMP complex
Authors	:	Hsiao, Y.Y.; Huang, K.W.; Wu, C.Y.
Deposited on		
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 (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:

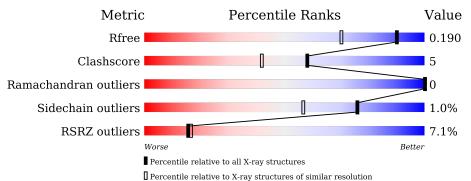
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7(2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

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	$\begin{array}{l} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	2936 (1.50-1.50)
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)
RSRZ outliers	127900	2884 (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 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	А	250	81%	6% ·	11%
1	В	250	<u>6%</u> 81%	9%	10%



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2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 7686 atoms, of which 3522 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
1	Δ	222	Total	С	Η	Ν	0	S	0	0	0
			3453	1091	1730	304	319	9	0		
1	Р	226	Total	С	Η	Ν	0	S	0	1	0
	1 B	220	3507	1108	1755	307	328	9	0	L	0

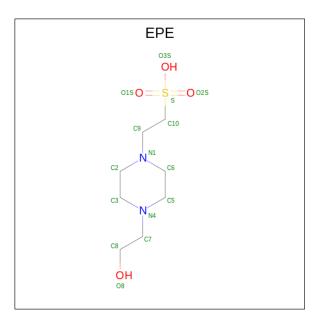
• Molecule 1 is a protein called Three-prime repair exonuclease 1.

A243LEU-expression tagA244GLU-expression tagA245HIS-expression tagA246HIS-expression tag	UNP Q91XB0 UNP Q91XB0
A 245 HIS - expression tag	UNP Q91XB0
	•
A 246 HIS overossion tag	-
A 240 IIIS - expression tag	UNP Q91XB0
A 247 HIS - expression tag	UNP Q91XB0
A 248 HIS - expression tag	UNP Q91XB0
A 249 HIS - expression tag	UNP Q91XB0
A 250 HIS - expression tag	UNP Q91XB0
B 243 LEU - expression tag	UNP Q91XB0
B 244 GLU - expression tag	UNP Q91XB0
B 245 HIS - expression tag	UNP Q91XB0
B 246 HIS - expression tag	UNP Q91XB0
B 247 HIS - expression tag	UNP Q91XB0
B 248 HIS - expression tag	UNP Q91XB0
B 249 HIS - expression tag	UNP Q91XB0
B 250 HIS - expression tag	UNP Q91XB0

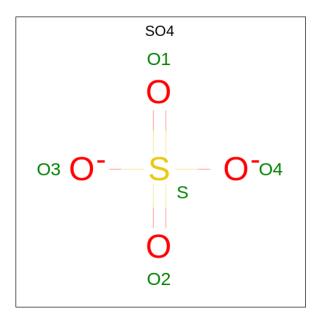
There are 16 discrepancies between the modelled and reference sequences:

• Molecule 2 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: $C_8H_{18}N_2O_4S$).





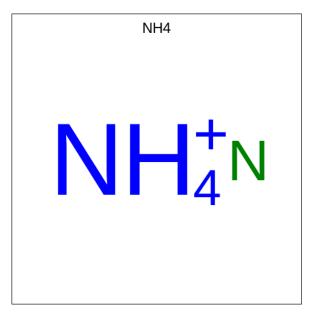
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf		
2	А	1	Total 32		H 17	N 2	0 4	S 1	0	0



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
3	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0

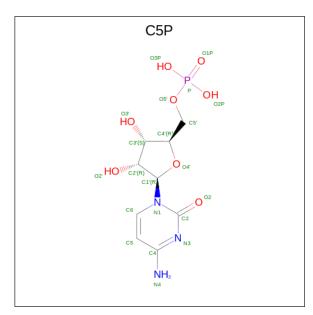
• Molecule 4 is AMMONIUM ION (three-letter code: NH4) (formula: H_4N).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	1	$\begin{array}{rrrr} \text{Total} & \text{H} & \text{N} \\ 5 & 4 & 1 \end{array}$	0	0
4	А	1	$\begin{array}{rrrr} \text{Total} & \text{H} & \text{N} \\ 5 & 4 & 1 \end{array}$	0	0

• Molecule 5 is CYTIDINE-5'-MONOPHOSPHATE (three-letter code: C5P) (formula: $C_9H_{14}N_3O_8P$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf		
5	В	1	Total 33	-	Н 12		-	Р 1	0	0



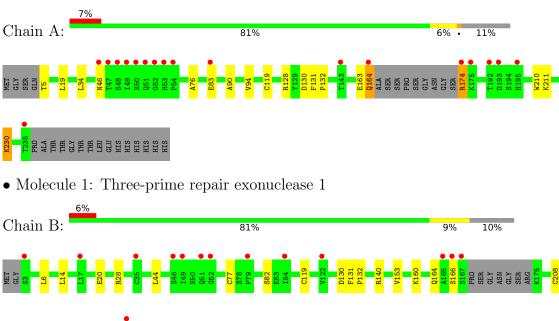
• Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	312	Total O 312 312	0	0
6	В	329	Total O 329 329	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: Three-prime repair exonuclease 1



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	64.04Å 85.66Å 100.03Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	26.08 - 1.50	Depositor
	26.08 - 1.50	EDS
% Data completeness	96.2(26.08-1.50)	Depositor
(in resolution range)	96.3(26.08-1.50)	EDS
R_{merge}	0.04	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$2.57 (at 1.50 \text{\AA})$	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.157 , 0.190	Depositor
It, Itfree	0.157 , 0.190	DCC
R_{free} test set	6626 reflections $(7.78%)$	wwPDB-VP
Wilson B-factor ($Å^2$)	18.1	Xtriage
Anisotropy	0.174	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.41 , 67.1	EDS
L-test for twinning ²	$ < L >=0.48, < L^2>=0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7686	wwPDB-VP
Average B, all atoms $(Å^2)$	31.0	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 25.54 % 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 3.1114e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.

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.



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: NH4, C5P, EPE, SO4 $\,$

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 lengths		Bond	angles
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5		
1	А	0.51	2/1765~(0.1%)	0.68	0/2407		
1	В	0.49	1/1794~(0.1%)	0.66	0/2447		
All	All	0.50	3/3559~(0.1%)	0.67	0/4854		

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	В	119	CYS	CB-SG	-6.99	1.70	1.82
1	А	164	GLN	CB-CG	-6.20	1.35	1.52
1	А	119	CYS	CB-SG	-5.21	1.73	1.81

There are no bond angle outliers.

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	А	1723	1730	1729	19	1
1	В	1752	1755	1754	15	0
2	А	15	17	17	0	0
3	А	10	0	0	1	0
4	А	2	8	0	0	0
5	В	21	12	12	0	0



	0	Non-H	1 0	H(added)	Clashes	Symm-Clashes
6	А	312	0	0	8	0
6	В	329	0	0	7	1
All	All	4164	3522	3512	34	1

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 5.

All (34) 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	distance (\AA)	overlap (Å)
1:A:164:GLN:HA	1:A:164:GLN:OE1	1.85	0.74
1:B:6:LEU:O	6:B:401:HOH:O	2.07	0.70
1:A:230:LYS:NZ	6:A:404:HOH:O	2.25	0.69
1:A:46:ASN:OD1	6:A:401:HOH:O	2.12	0.67
1:A:164:GLN:OE1	1:A:164:GLN:CA	2.44	0.65
1:B:20:GLU:OE1	6:B:402:HOH:O	2.15	0.64
1:A:19:LEU:CD2	1:A:34:LEU:HD13	2.30	0.62
1:A:5:THR:N	6:A:406:HOH:O	2.33	0.62
1:B:208:CYS:HB3	1:B:215:LEU:HD22	1.82	0.61
1:B:28:ARG:NH1	1:B:233:TYR:HB2	2.17	0.60
1:A:174:ARG:O	6:A:402:HOH:O	2.17	0.59
1:A:19:LEU:HD21	1:A:34:LEU:HD13	1.85	0.58
1:A:163:GLU:O	1:A:164:GLN:HG2	2.07	0.54
1:B:213:GLN:NE2	6:B:403:HOH:O	2.20	0.53
1:A:76:ALA:HB1	3:A:302:SO4:O3	2.09	0.53
1:A:46:ASN:ND2	6:A:403:HOH:O	2.23	0.53
1:B:140:ARG:HB3	1:B:140:ARG:CZ	2.40	0.51
1:A:83:GLU:HG3	6:A:438:HOH:O	2.10	0.51
1:B:160:LYS:HB3	6:B:422:HOH:O	2.11	0.49
1:A:83:GLU:CD	6:A:415:HOH:O	2.51	0.48
1:B:164:GLN:HG3	6:B:422:HOH:O	2.14	0.48
1:A:210:TRP:CZ2	1:A:211:LYS:HE3	2.50	0.47
1:A:131:PHE:HB2	1:A:132:PRO:HD3	1.96	0.47
1:A:163:GLU:C	1:A:164:GLN:HG2	2.36	0.46
1:B:131:PHE:HB2	1:B:132:PRO:HD3	1.99	0.45
1:B:14:LEU:HD22	1:B:44:LEU:CD1	2.46	0.45
1:A:230:LYS:NZ	6:A:424:HOH:O	2.48	0.45
1:B:211:LYS:HA	6:B:403:HOH:O	2.17	0.43
1:B:166:SER:O	1:B:166:SER:OG	2.37	0.43
1:B:77:CYS:HB2	1:B:82:SER:OG	2.18	0.43
1:A:90:ALA:O	1:A:94:VAL:HG13	2.19	0.42



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:153:VAL:CG1	1:B:223:ALA:HA	2.49	0.42
1:B:217:GLN:HG2	6:B:657:HOH:O	2.20	0.41
1:A:230:LYS:HE2	1:A:230:LYS:HB2	1.82	0.41

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:ARG:NH1	6:B:415:HOH:O[2_655]	2.03	0.17

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	Perce	ntiles
1	А	218/250~(87%)	217 (100%)	1 (0%)	0	100	100
1	В	223/250~(89%)	222 (100%)	1 (0%)	0	100	100
All	All	441/500 (88%)	439 (100%)	2~(0%)	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 side chain 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	А	192/214~(90%)	189~(98%)	3~(2%)	62 36	
				Continued of	n nert nage	



Conti	naea jion	i previous puye				
Mol	Chain	Analysed	Rotameric	Outliers	Perce	entiles
1	В	196/214~(92%)	195 (100%)	1 (0%)	88	78
All	All	388/428~(91%)	384 (99%)	4 (1%)	76	57

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

Mol	Chain	Res	Type
1	А	130	ASP
1	А	174	ARG
1	А	230	LYS
1	В	130	ASP

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

Mol	Chain	Res	Type
1	А	98	GLN
1	В	98	GLN

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)

Of 6 ligands modelled in this entry, 2 are modelled with single atom - leaving 4 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



Mol	Tune Chain		Res	Link	B	Bond lengths			Bond angles		
10101	Type	Chain	nes	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
2	EPE	А	301	-	15,15,15	0.82	1 (6%)	18,20,20	1.76	5 (27%)	
5	C5P	В	301	-	22,22,22	3.44	11 (50%)	33,33,33	0.95	2 (6%)	
3	SO4	А	303	-	4,4,4	0.16	0	6,6,6	0.09	0	
3	SO4	А	302	-	4,4,4	0.14	0	$6,\!6,\!6$	0.42	0	

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).

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	EPE	А	301	-	-	3/9/19/19	0/1/1/1
5	C5P	В	301	-	-	2/10/26/26	0/2/2/2

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\operatorname{Observed}(\operatorname{\AA})$	$\mathrm{Ideal}(\mathrm{\AA})$
5	В	301	C5P	C2'-C3'	-8.82	1.29	1.53
5	В	301	C5P	C2-N3	5.74	1.48	1.36
5	В	301	C5P	O4'-C4'	-5.49	1.32	1.45
5	В	301	C5P	C6-C5	5.27	1.47	1.35
5	В	301	C5P	C4-N3	4.21	1.43	1.34
5	В	301	C5P	O4'-C1'	3.91	1.51	1.42
5	В	301	C5P	C3'-C4'	3.81	1.62	1.53
5	В	301	C5P	C4-N4	3.06	1.41	1.33
5	В	301	C5P	C6-N1	2.92	1.45	1.38
5	В	301	C5P	C2-N1	2.88	1.46	1.40
2	А	301	EPE	C10-S	2.58	1.81	1.77
5	В	301	C5P	C5-C4	2.13	1.47	1.42

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	А	301	EPE	C5-N4-C3	4.51	118.97	108.83
2	А	301	EPE	C7-N4-C5	3.31	119.69	111.23
5	В	301	C5P	C4'-O4'-C1'	-2.87	103.13	109.47
2	А	301	EPE	C7-N4-C3	2.53	117.70	111.23



Mol	Chain	Res	Type	Atoms		$Observed(^{o})$	$Ideal(^{o})$
5	В	301	C5P	O4'-C1'-N1	-2.44	102.78	108.36
2	А	301	EPE	C6-N1-C2	2.30	114.00	108.83
2	А	301	EPE	C3-C2-N1	2.02	114.78	110.64

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	А	301	EPE	C9-C10-S-O1S
2	А	301	EPE	C9-C10-S-O2S
2	А	301	EPE	C9-C10-S-O3S
5	В	301	C5P	C5'-O5'-P-O3P
5	В	301	C5P	C5'-O5'-P-O1P

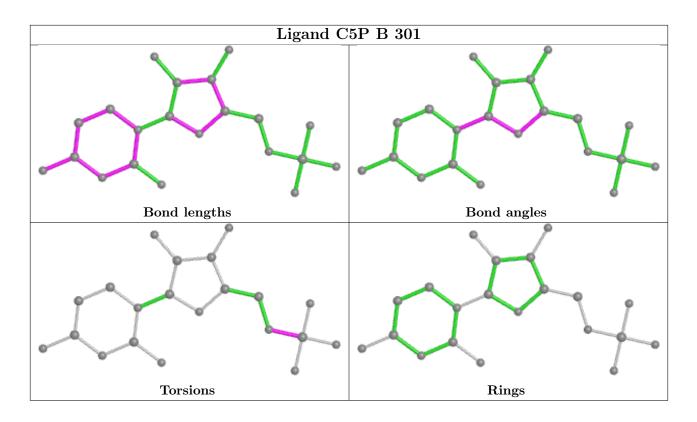
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	А	302	SO4	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	$\langle RSRZ \rangle$	#RSRZ>2		$OWAB(Å^2)$	Q<0.9	
1	А	222/250~(88%)	-0.05	18 (8%)	12	12	12, 20, 56, 105	0
1	В	226/250~(90%)	-0.06	14 (6%)	20	22	12, 21, 52, 104	0
All	All	448/500~(89%)	-0.05	32 (7%)	16	17	12, 21, 55, 105	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	49	ILE	8.5
1	В	167	SER	8.3
1	А	51	GLN	7.1
1	А	235	THR	6.5
1	А	192	THR	6.4
1	А	174	ARG	6.3
1	В	49	ILE	5.9
1	В	235	THR	5.3
1	А	48	SER	5.0
1	А	52	GLY	4.4
1	А	50	SER	4.1
1	А	195	HIS	3.8
1	В	51	GLN	3.6
1	В	166	SER	3.6
1	А	164	GLN	3.5
1	А	53	HIS	3.3
1	В	17	LEU	3.2
1	А	46	ASN	2.9
1	В	165	ALA	2.8
1	А	83	GLU	2.6
1	А	47	THR	2.5
1	А	143	THR	2.4
1	А	54	PRO	2.3
1	В	48	SER	2.3



Mol	Chain	Res	Type	RSRZ
1	В	3	SER	2.2
1	В	79	PRO	2.2
1	В	52	GLY	2.2
1	А	193	ASP	2.2
1	В	122	VAL	2.1
1	В	84	ILE	2.1
1	В	35	CYS	2.1
1	А	175	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 monosaccharides in this entry.

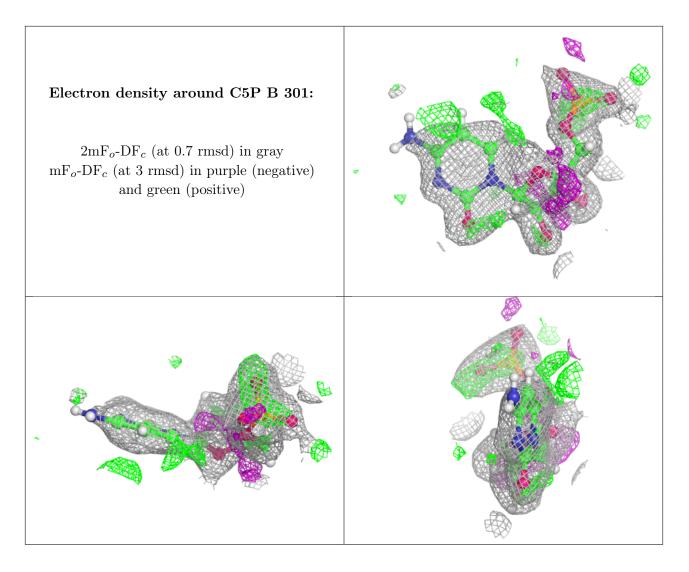
6.4 Ligands (i)

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	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q < 0.9
5	C5P	В	301	21/21	0.78	0.15	26,48,96,111	0
3	SO4	А	303	5/5	0.82	0.24	48,60,70,93	0
3	SO4	А	302	5/5	0.87	0.17	$26,\!67,\!102,\!105$	0
2	EPE	А	301	15/15	0.95	0.20	28,45,71,75	0
4	NH4	А	305	1/1	0.96	0.11	31,37,37,37	0
4	NH4	А	304	1/1	0.98	0.13	$15,\!17,\!17,\!17$	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.

