



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 07:10 pm BST

PDB ID : 6OIQ
Title : Crystal structure of MYST acetyltransferase domain in complex with inhibitor 63
Authors : Hermans, S.J.; Chung, M.C.; Parker, M.W.; Thomas, T.; Baell, J.B.
Deposited on : 2019-04-09
Resolution : 1.75 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 : 1.8.5 (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

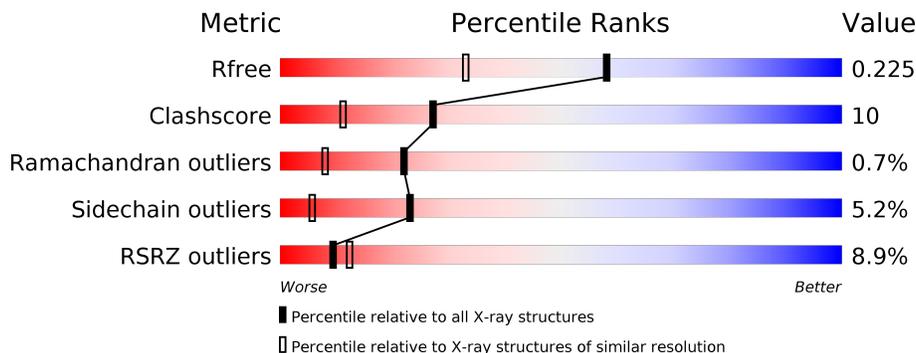
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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}	130704	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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 $\leq 5\%$. 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	273	<div style="display: flex; align-items: center;"> <div style="width: 9%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 81%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 15%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center; margin-top: 5px;">9% 81% 15% •</p>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2514 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 Histone acetyltransferase KAT8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	272	2316	1514	380	410	12	0	9	0

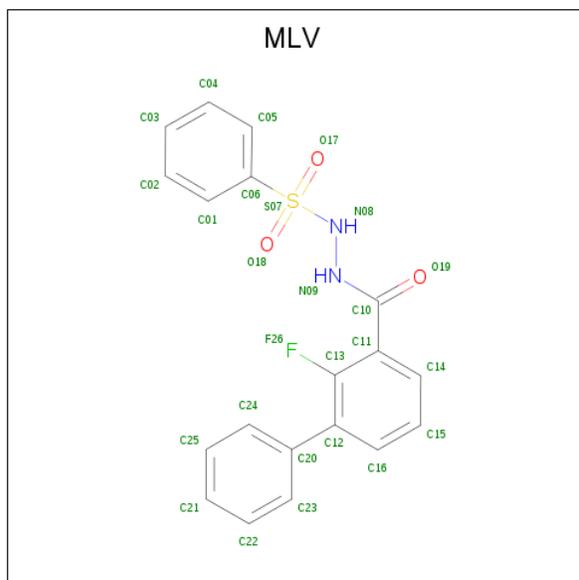
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	579	HIS	TYR	conflict	UNP Q9H7Z6
A	645	SER	ALA	engineered mutation	UNP Q9H7Z6
A	648	MET	LEU	engineered mutation	UNP Q9H7Z6
A	649	ILE	THR	engineered mutation	UNP Q9H7Z6
A	660	ARG	LYS	engineered mutation	UNP Q9H7Z6
A	702	ASN	ILE	conflict	UNP Q9H7Z6

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

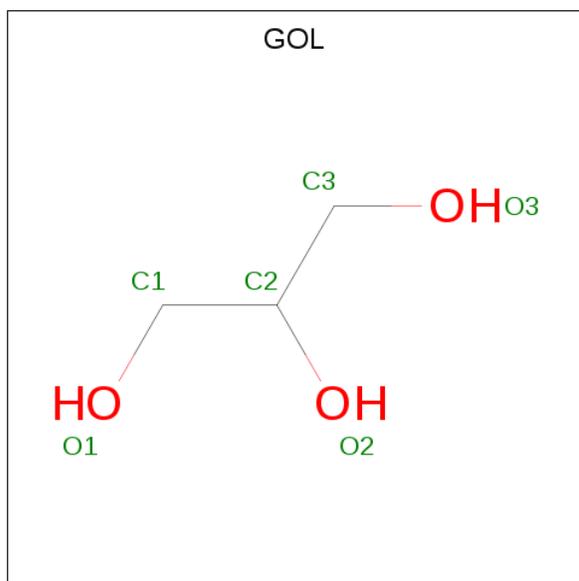
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is 2-fluoro-N'-(phenylsulfonyl)[1,1'-biphenyl]-3-carbohydrazide (three-letter code: MLV) (formula: C₁₉H₁₅FN₂O₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	F	N	O			S
3	A	1	26	19	1	2	3	1	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	C O		
4	A	1	6	3 3	0	0

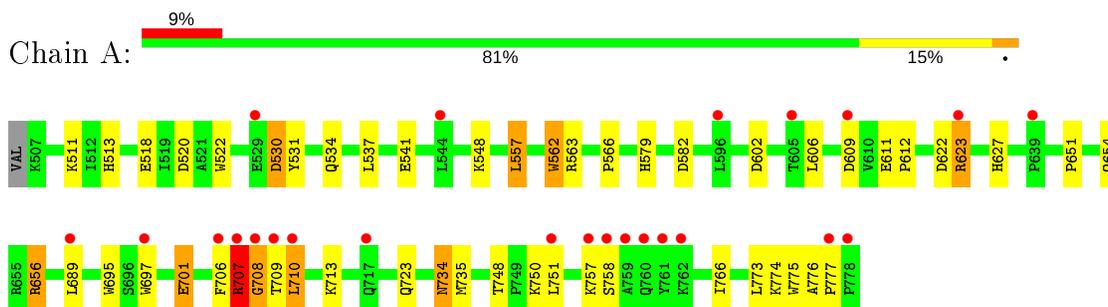
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	165	Total 165	O 165	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: Histone acetyltransferase KAT8



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	46.60Å 56.77Å 122.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	122.28 – 1.75 27.65 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.7 (122.28-1.75) 99.8 (27.65-1.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.17 (at 1.75Å)	Xtrriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.184 , 0.227 0.183 , 0.225	Depositor DCC
R_{free} test set	1695 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	16.7	Xtrriage
Anisotropy	0.210	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 46.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2514	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.51% 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: GOL, ZN, MLV, ALY

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	1.22	6/2403 (0.2%)	1.09	6/3255 (0.2%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	695	TRP	CD2-CE2	7.20	1.50	1.41
1	A	522	TRP	CD2-CE2	5.96	1.48	1.41
1	A	531	TYR	CE1-CZ	5.84	1.46	1.38
1	A	562[A]	TRP	CD2-CE2	5.64	1.48	1.41
1	A	562[B]	TRP	CD2-CE2	5.64	1.48	1.41
1	A	775	TRP	CD2-CE2	5.00	1.47	1.41

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	557	LEU	CB-CG-CD1	9.39	126.96	111.00
1	A	563	ARG	NE-CZ-NH2	-6.55	117.03	120.30
1	A	582	ASP	CB-CG-OD1	5.93	123.64	118.30
1	A	766	ILE	CB-CA-C	-5.36	100.88	111.60
1	A	602	ASP	CB-CG-OD1	5.26	123.04	118.30
1	A	537	LEU	CB-CG-CD2	-5.25	102.08	111.00

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	2316	0	2299	47	0
2	A	1	0	0	0	0
3	A	26	0	0	0	0
4	A	6	0	8	0	0
5	A	165	0	0	14	3
All	All	2514	0	2307	47	3

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 (47) 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:697[A]:TRP:HE3	5:A:1052:HOH:O	1.08	1.31
1:A:697[B]:TRP:CZ2	1:A:701[B]:GLU:OE1	1.98	1.15
1:A:707:ARG:H	1:A:708:GLY:HA3	0.95	1.09
1:A:707:ARG:N	1:A:708:GLY:HA3	1.69	1.04
1:A:757:LYS:HG3	5:A:939:HOH:O	1.57	1.01
1:A:697[B]:TRP:CE2	1:A:701[B]:GLU:OE1	2.22	0.92
1:A:623:ARG:HH11	1:A:623:ARG:HG2	1.35	0.89
1:A:520:ASP:O	5:A:901:HOH:O	1.91	0.89
1:A:707:ARG:H	1:A:708:GLY:CA	1.84	0.89
1:A:689:LEU:CD1	5:A:980:HOH:O	2.22	0.85
1:A:748:THR:HG23	1:A:751:LEU:H	1.41	0.85
1:A:697[B]:TRP:CE2	1:A:701[B]:GLU:CD	2.52	0.83
1:A:697[B]:TRP:CH2	1:A:701[B]:GLU:OE1	2.36	0.79
1:A:707:ARG:N	1:A:708:GLY:CA	2.44	0.77
1:A:697[B]:TRP:CZ2	1:A:701[B]:GLU:CD	2.61	0.73
1:A:541:GLU:OE1	1:A:627:HIS:HD2	1.72	0.72
1:A:622:ASP:OD1	1:A:627:HIS:HE1	1.73	0.69
1:A:689:LEU:HD13	5:A:980:HOH:O	1.88	0.69
1:A:697[B]:TRP:CZ3	5:A:903:HOH:O	2.46	0.67
1:A:707:ARG:HG3	1:A:708:GLY:HA2	1.77	0.67
1:A:611:GLU:HG2	1:A:612:PRO:HD3	1.81	0.63
1:A:706:PHE:O	1:A:707:ARG:HB3	2.00	0.62
1:A:774:LYS:NZ	5:A:902:HOH:O	2.22	0.61
1:A:689:LEU:HD11	5:A:980:HOH:O	1.91	0.60
1:A:513:HIS:HD2	1:A:518:GLU:OE2	1.83	0.60
1:A:530[B]:ASP:OD2	1:A:548:LYS:NZ	2.33	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:706:PHE:HB2	1:A:710:LEU:HD22	1.88	0.54
1:A:697[A]:TRP:CE3	5:A:1052:HOH:O	2.00	0.52
1:A:697[B]:TRP:CE2	1:A:701[B]:GLU:OE2	2.63	0.51
1:A:606:LEU:HD21	1:A:609:ASP:HB3	1.91	0.51
1:A:697[B]:TRP:CZ2	1:A:701[B]:GLU:OE2	2.65	0.50
1:A:623:ARG:NH1	1:A:623:ARG:HG2	2.11	0.49
1:A:562[A]:TRP:HZ3	1:A:566:PRO:HD3	1.78	0.49
1:A:707:ARG:NE	1:A:707:ARG:O	2.45	0.49
1:A:734:ASN:C	5:A:905:HOH:O	2.50	0.49
1:A:707:ARG:HG3	1:A:708:GLY:CA	2.44	0.48
1:A:654:GLN:O	1:A:656:ARG:HD2	2.15	0.47
1:A:534:GLN:HB2	5:A:995:HOH:O	2.16	0.45
1:A:606:LEU:HD13	1:A:606:LEU:C	2.37	0.45
1:A:697[B]:TRP:NE1	1:A:701[B]:GLU:OE2	2.52	0.43
1:A:520:ASP:HB3	5:A:998:HOH:O	2.19	0.43
1:A:511:LYS:NZ	1:A:520:ASP:OD1	2.46	0.43
1:A:735:MET:N	5:A:905:HOH:O	2.50	0.43
1:A:579:HIS:HE1	5:A:906:HOH:O	2.02	0.43
1:A:713:LYS:HG2	1:A:723:GLN:NE2	2.33	0.43
1:A:776:ALA:HB1	1:A:777:PRO:HD2	2.01	0.42
1:A:611:GLU:CD	1:A:611:GLU:H	2.22	0.42

All (3) 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)
5:A:904:HOH:O	5:A:968:HOH:O[4_545]	0.67	1.53
5:A:931:HOH:O	5:A:938:HOH:O[3_555]	1.30	0.90
5:A:998:HOH:O	5:A:1028:HOH:O[4_545]	2.14	0.06

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	278/273 (102%)	269 (97%)	7 (2%)	2 (1%)	22 8

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	707	ARG
1	A	708	GLY

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	259/252 (103%)	243 (94%)	16 (6%)	18 4

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

Mol	Chain	Res	Type
1	A	530[A]	ASP
1	A	530[B]	ASP
1	A	557	LEU
1	A	623	ARG
1	A	651	PRO
1	A	656	ARG
1	A	701[A]	GLU
1	A	701[B]	GLU
1	A	707	ARG
1	A	709	THR
1	A	710	LEU
1	A	734	ASN
1	A	750	LYS
1	A	758[A]	SER
1	A	758[B]	SER
1	A	773	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	513	HIS
1	A	516	ASN
1	A	586	HIS
1	A	624	GLN
1	A	627	HIS
1	A	702	ASN
1	A	723	GLN
1	A	734	ASN
1	A	760	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](#)

1 non-standard protein/DNA/RNA residue 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	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	ALY	A	604	1	10,11,12	0.88	0	7,12,14	2.26	2 (28%)

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
1	ALY	A	604	1	-	1/9/10/12	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	604	ALY	CH3-CH-NZ	-4.32	108.44	116.09
1	A	604	ALY	CD-CE-NZ	2.18	118.44	112.21

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	604	ALY	CG-CD-CE-NZ

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	GOL	A	803	-	5,5,5	0.27	0	5,5,5	1.02	0
3	MLV	A	802	-	28,28,28	3.44	9 (32%)	38,39,39	2.62	16 (42%)

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	GOL	A	803	-	-	3/4/4/4	-
3	MLV	A	802	-	-	1/20/20/20	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	802	MLV	O18-S07	8.51	1.53	1.43
3	A	802	MLV	N08-N09	7.85	1.46	1.40
3	A	802	MLV	S07-N08	7.22	1.75	1.65
3	A	802	MLV	F26-C13	7.13	1.46	1.35
3	A	802	MLV	C06-S07	4.53	1.83	1.76
3	A	802	MLV	C12-C20	-3.89	1.42	1.49
3	A	802	MLV	C11-C10	-3.87	1.42	1.50
3	A	802	MLV	C14-C11	3.22	1.45	1.39
3	A	802	MLV	C16-C12	2.90	1.44	1.40

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	802	MLV	O18-S07-O17	-5.49	112.80	119.55
3	A	802	MLV	C04-C05-C06	-5.31	113.44	118.95
3	A	802	MLV	O18-S07-N08	5.22	115.55	105.57
3	A	802	MLV	C05-C06-C01	5.09	127.55	120.44
3	A	802	MLV	O18-S07-C06	-5.04	101.76	107.97
3	A	802	MLV	C05-C06-S07	-4.94	114.40	119.77
3	A	802	MLV	C16-C15-C14	4.77	127.02	120.25
3	A	802	MLV	C15-C14-C11	-3.34	113.51	119.81
3	A	802	MLV	C03-C02-C01	-2.92	115.75	120.19
3	A	802	MLV	C13-C11-C10	-2.50	120.48	125.33
3	A	802	MLV	C14-C11-C13	2.45	121.01	117.18
3	A	802	MLV	O17-S07-N08	2.45	110.25	105.57
3	A	802	MLV	C04-C03-C02	2.30	124.21	119.93
3	A	802	MLV	C15-C16-C12	-2.24	116.55	120.33
3	A	802	MLV	C22-C23-C20	2.09	123.19	120.56
3	A	802	MLV	F26-C13-C11	-2.02	116.80	121.10

There are no chirality outliers.

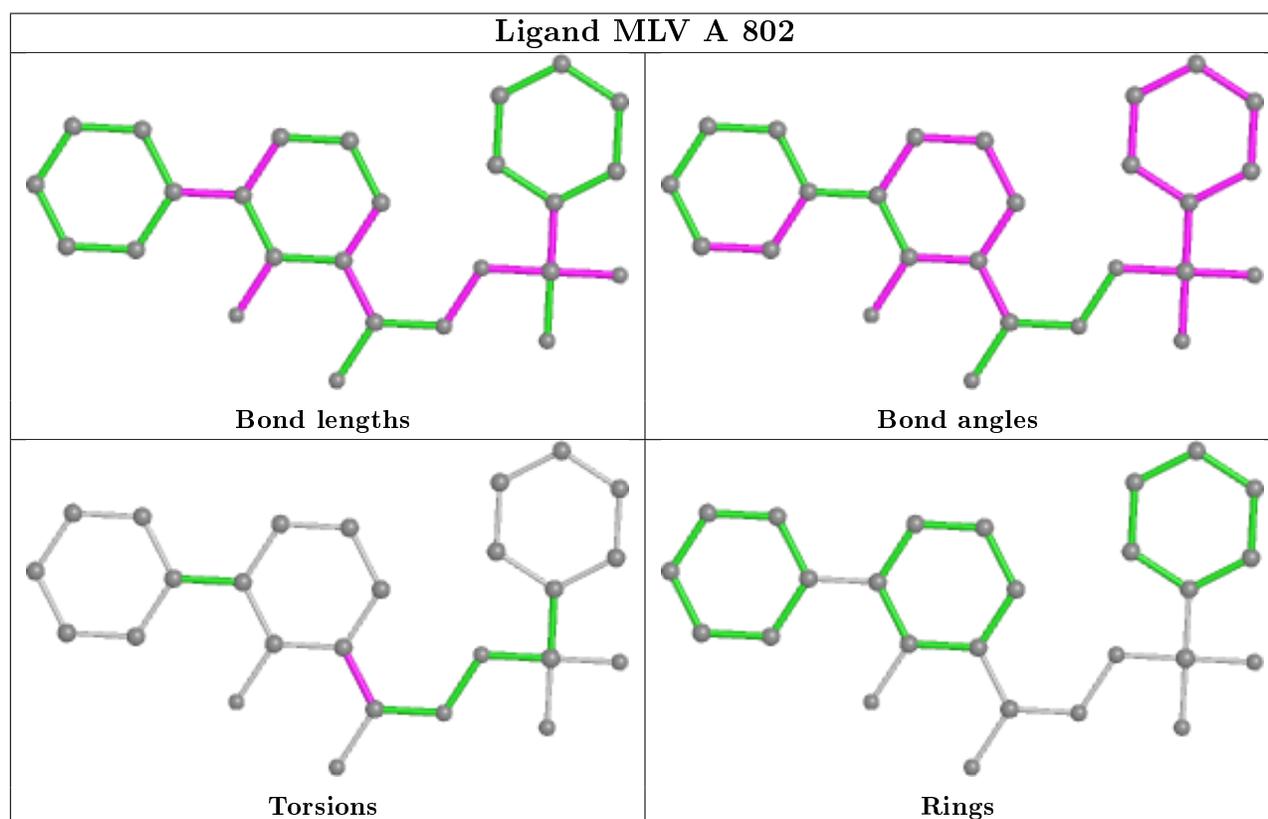
All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	803	GOL	C1-C2-C3-O3
4	A	803	GOL	O2-C2-C3-O3
3	A	802	MLV	O19-C10-C11-C13
4	A	803	GOL	O1-C1-C2-C3

There are no ring outliers.

No monomer is involved in short contacts.

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	271/273 (99%)	0.42	24 (8%) 9 12	8, 17, 42, 89	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	709	THR	7.5
1	A	707	ARG	7.0
1	A	762	LYS	5.1
1	A	759	ALA	5.0
1	A	706	PHE	4.4
1	A	760	GLN	3.8
1	A	708	GLY	3.7
1	A	697[A]	TRP	3.4
1	A	757	LYS	3.3
1	A	751	LEU	3.2
1	A	710	LEU	3.1
1	A	777	PRO	2.8
1	A	623	ARG	2.7
1	A	529	GLU	2.5
1	A	689	LEU	2.4
1	A	758[A]	SER	2.3
1	A	596	LEU	2.2
1	A	717	GLN	2.2
1	A	761	TYR	2.2
1	A	544	LEU	2.2
1	A	609	ASP	2.2
1	A	639	PRO	2.1
1	A	778	PRO	2.1
1	A	605	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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, 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(Å ²)	Q<0.9
1	ALY	A	604	12/13	0.94	0.11	12,14,22,22	0

6.3 Carbohydrates [i](#)

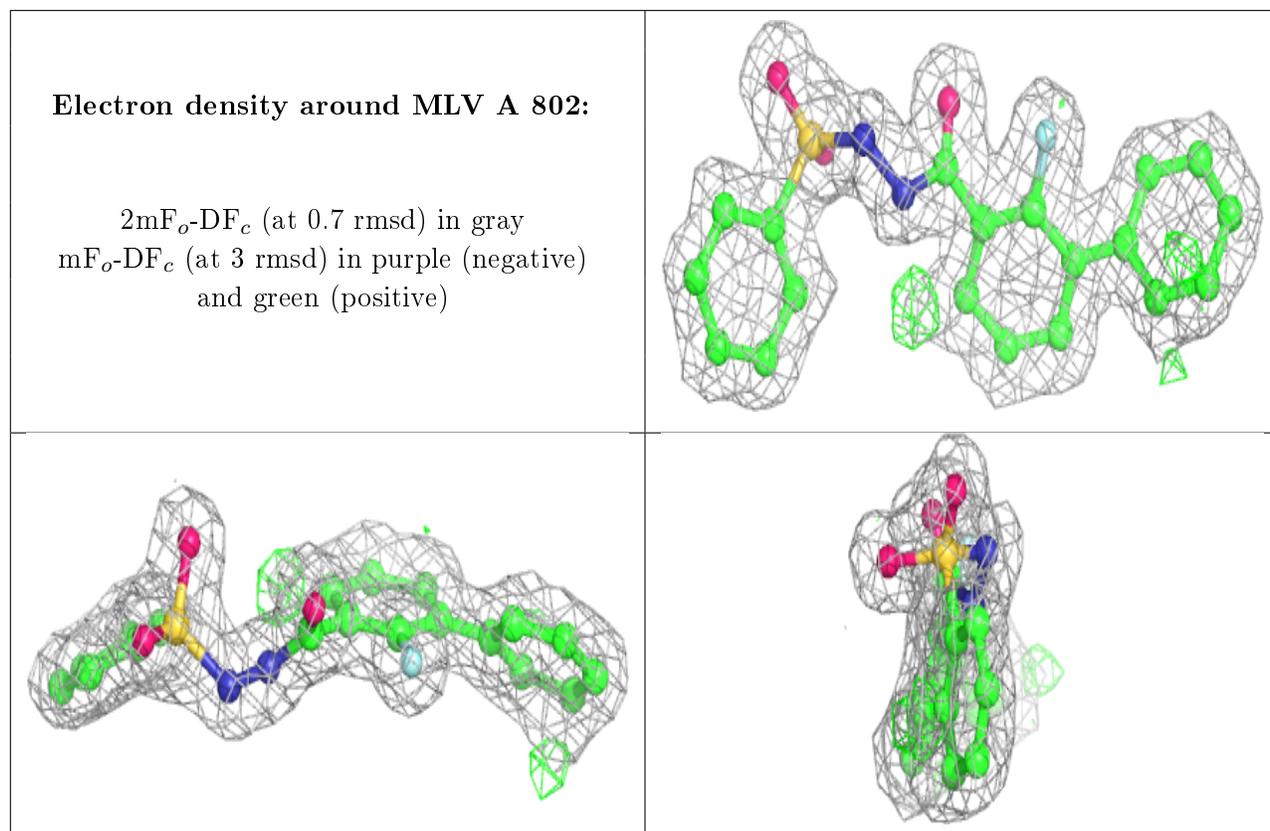
There are no carbohydrates in this entry.

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, 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(Å ²)	Q<0.9
4	GOL	A	803	6/6	0.79	0.19	32,42,45,48	0
3	MLV	A	802	26/26	0.97	0.08	13,14,16,18	0
2	ZN	A	801	1/1	1.00	0.07	10,10,10,10	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.