



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:06 PM UTC

PDB ID : 4UCK / pdb_00004uck
Title : X-ray structure and activities of an essential Mononegavirales L- protein domain
Authors : Paesen, G.C.; Collet, A.; Sallamand, C.; Debart, F.; Vasseur, J.J.; Canard, B.; Decroly, E.; Grimes, J.M.
Deposited on : 2014-12-03
Resolution : 2.66 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 types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

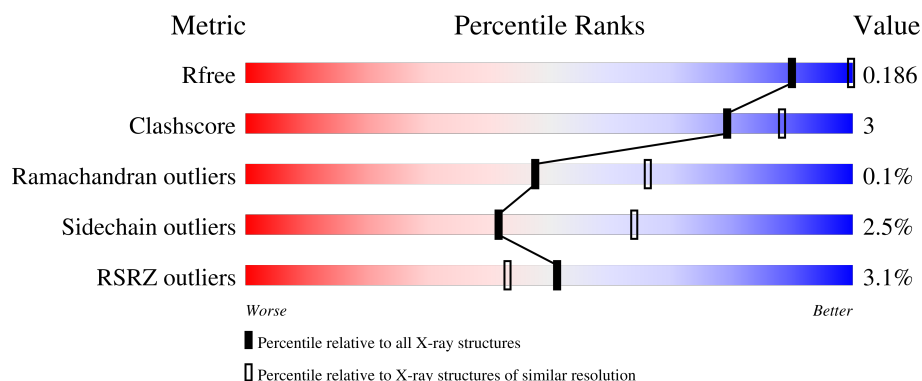
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.66 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}	180053	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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 $\leq 5\%$. 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	415	3% 79% 10% 11%
1	B	415	3% 82% 8% 9%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6276 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 RNA-DIRECTED RNA POLYMERASE L.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	370	Total	C	N	O	S	0	0	0
			3011	1937	513	541	20			
1	B	376	Total	C	N	O	S	0	0	0
			3054	1961	520	553	20			

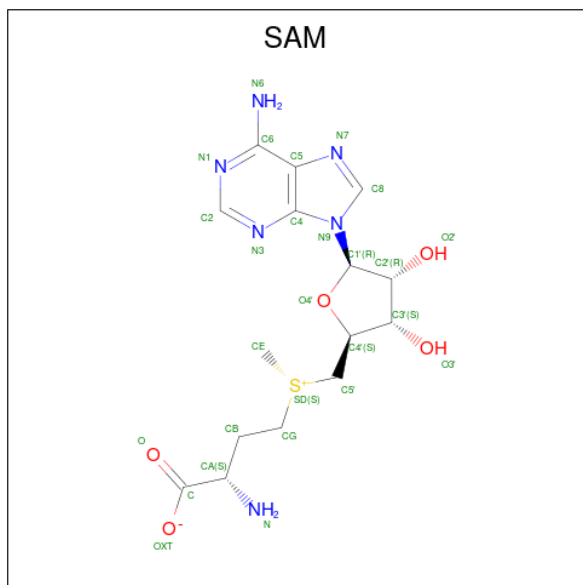
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1599	MET	-	expression tag	UNP Q91L20
A	2006	SER	-	expression tag	UNP Q91L20
A	2007	GLY	-	expression tag	UNP Q91L20
A	2008	HIS	-	expression tag	UNP Q91L20
A	2009	HIS	-	expression tag	UNP Q91L20
A	2010	HIS	-	expression tag	UNP Q91L20
A	2011	HIS	-	expression tag	UNP Q91L20
A	2012	HIS	-	expression tag	UNP Q91L20
A	2013	HIS	-	expression tag	UNP Q91L20
A	1995	GLU	LYS	engineered mutation	UNP Q91L20
B	1599	MET	-	expression tag	UNP Q91L20
B	2006	SER	-	expression tag	UNP Q91L20
B	2007	GLY	-	expression tag	UNP Q91L20
B	2008	HIS	-	expression tag	UNP Q91L20
B	2009	HIS	-	expression tag	UNP Q91L20
B	2010	HIS	-	expression tag	UNP Q91L20
B	2011	HIS	-	expression tag	UNP Q91L20
B	2012	HIS	-	expression tag	UNP Q91L20
B	2013	HIS	-	expression tag	UNP Q91L20
B	1995	GLU	LYS	engineered mutation	UNP Q91L20

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 3 is S-ADENOSYLMETHIONINE (CCD ID: SAM) (formula: $C_{15}H_{22}N_6O_5S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
3	B	1	Total	C	N	O	S	0	0
			27	15	6	5	1		

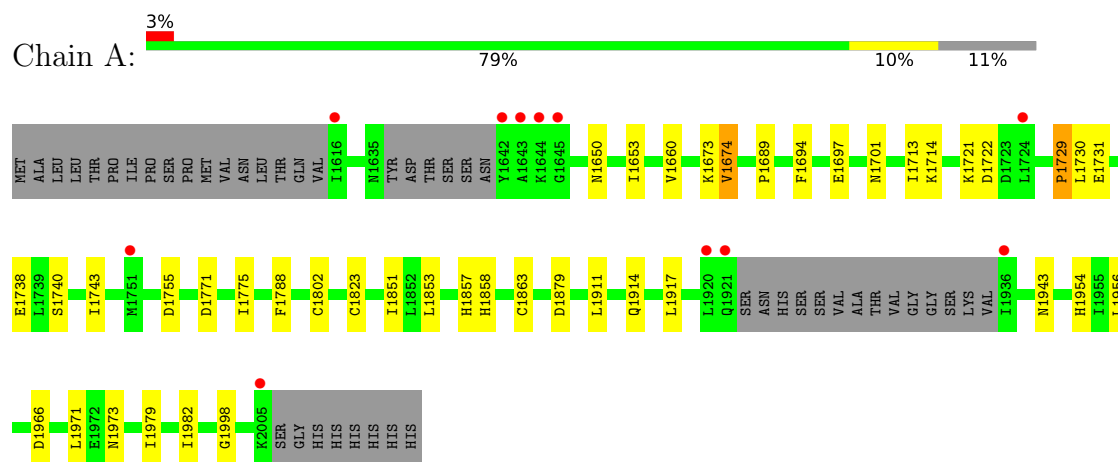
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	74	Total	O	0	0
			74	74		
4	B	81	Total	O	0	0
			81	81		

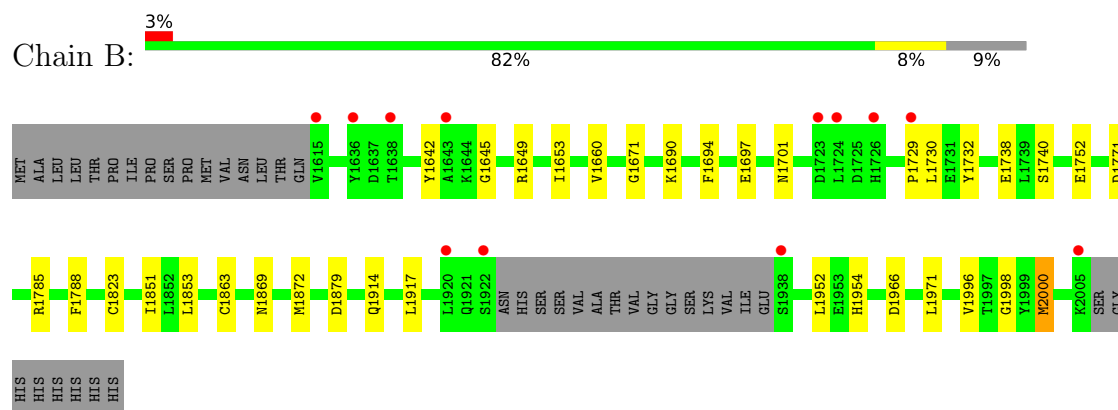
3 Residue-property plots [i](#)

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: RNA-DIRECTED RNA POLYMERASE L



• Molecule 1: RNA-DIRECTED RNA POLYMERASE L



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.34Å 84.35Å 181.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	76.51 – 2.66 76.51 – 2.66	Depositor EDS
% Data completeness (in resolution range)	98.5 (76.51-2.66) 98.7 (76.51-2.66)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.62 (at 2.65Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.186 , 0.204 0.194 , 0.186	Depositor DCC
R_{free} test set	1809 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	74.7	Xtriage
Anisotropy	0.623	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 66.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.026 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6276	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% 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: ZN, SAM

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	0.70	0/3075	1.31	8/4145 (0.2%)
1	B	0.71	1/3120 (0.0%)	1.29	1/4209 (0.0%)
All	All	0.71	1/6195 (0.0%)	1.30	9/8354 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	2000	MET	SD-CE	-6.23	1.64	1.79

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1802	CYS	CA-C-N	6.10	128.74	120.38
1	A	1802	CYS	C-N-CA	6.10	128.74	120.38
1	A	1729	PRO	CA-C-N	5.82	128.35	120.38
1	A	1729	PRO	C-N-CA	5.82	128.35	120.38
1	A	1755	ASP	CA-CB-CG	5.53	118.13	112.60
1	A	1943	ASN	CA-C-N	5.23	127.29	120.28
1	A	1943	ASN	C-N-CA	5.23	127.29	120.28
1	B	1879	ASP	CA-CB-CG	5.07	117.67	112.60
1	A	1879	ASP	CA-CB-CG	5.02	117.62	112.60

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	3011	0	3053	17	0
1	B	3054	0	3087	17	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	27	0	22	0	0
3	B	27	0	22	0	0
4	A	74	0	0	0	0
4	B	81	0	0	0	0
All	All	6276	0	6184	32	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 3.

All (32) 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:B:1671:GLY:H	1:B:1701:ASN:HD22	1.23	0.87
1:B:1952:LEU:HD22	1:B:2000:MET:HE1	1.73	0.71
1:A:1701:ASN:HD22	1:A:1731:GLU:HG2	1.71	0.55
1:B:1671:GLY:N	1:B:1701:ASN:HD22	2.01	0.53
1:A:1788:PHE:HZ	1:A:1851:ILE:HD11	1.74	0.52
1:B:1851:ILE:HD12	1:B:1853:LEU:HD21	1.93	0.51
1:B:1788:PHE:HZ	1:B:1851:ILE:HD11	1.75	0.51
1:B:1823:CYS:HB3	1:B:1863:CYS:CB	2.41	0.51
1:A:1823:CYS:HB3	1:A:1863:CYS:CB	2.40	0.50
1:B:1914:GLN:HA	1:B:1917:LEU:HD12	1.93	0.50
1:A:1851:ILE:HD12	1:A:1853:LEU:HD21	1.93	0.50
1:A:1973:ASN:HB3	1:B:1642:TYR:HB2	1.93	0.50
1:B:1671:GLY:H	1:B:1701:ASN:ND2	2.00	0.50
1:A:1914:GLN:HA	1:A:1917:LEU:HD12	1.93	0.49
1:B:1694:PHE:HB3	1:B:1697:GLU:HG2	1.96	0.47
1:A:1979:ILE:HD12	1:A:1982:ILE:HD11	1.96	0.47
1:A:1660:VAL:O	1:A:1998:GLY:HA3	2.15	0.46
1:A:1694:PHE:HB3	1:A:1697:GLU:HG2	1.97	0.46
1:A:1689:PRO:HG3	1:A:1775:ILE:HD12	1.98	0.46
1:A:1650:ASN:O	1:B:1645:GLY:HA3	2.16	0.46
1:B:1729:PRO:HB2	1:B:1732:TYR:HB3	1.99	0.45
1:B:1649:ARG:HB3	1:B:1872:MET:HE2	2.00	0.44
1:A:1857:HIS:H	1:A:1857:HIS:CD2	2.36	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1714:LYS:HE2	1:A:1743:ILE:HG21	2.00	0.43
1:B:1996:VAL:HG12	1:B:2000:MET:HE2	2.00	0.43
1:B:1660:VAL:O	1:B:1998:GLY:HA3	2.19	0.43
1:B:1738:GLU:HG2	1:B:1740:SER:H	1.83	0.42
1:A:1673:LYS:HG2	1:A:1674:VAL:HG22	2.02	0.42
1:A:1738:GLU:HG2	1:A:1740:SER:H	1.84	0.41
1:B:1954:HIS:HE1	1:B:1971:LEU:O	2.04	0.41
1:A:1954:HIS:HE1	1:A:1971:LEU:O	2.04	0.41
1:A:1911:LEU:HD12	1:A:1956:LEU:HD22	2.02	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	364/415 (88%)	350 (96%)	13 (4%)	1 (0%)	36	52
1	B	372/415 (90%)	360 (97%)	12 (3%)	0	100	100
All	All	736/830 (89%)	710 (96%)	25 (3%)	1 (0%)	48	66

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1729	PRO

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	334/374 (89%)	325 (97%)	9 (3%)	39	61
1	B	340/374 (91%)	332 (98%)	8 (2%)	43	65
All	All	674/748 (90%)	657 (98%)	17 (2%)	42	64

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1653	ILE
1	A	1674	VAL
1	A	1713	ILE
1	A	1721	LYS
1	A	1722	ASP
1	A	1730	LEU
1	A	1771	ASP
1	A	1858	HIS
1	A	1966	ASP
1	B	1653	ILE
1	B	1690	LYS
1	B	1730	LEU
1	B	1752	GLU
1	B	1771	ASP
1	B	1785	ARG
1	B	1869	ASN
1	B	1966	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1635	ASN
1	A	1659	HIS
1	A	1824	ASN
1	A	1857	HIS
1	A	1868	GLN
1	A	1869	ASN
1	A	1894	ASN
1	A	1914	GLN
1	A	1954	HIS
1	A	1973	ASN
1	B	1621	GLN

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Mol	Chain	Res	Type
1	B	1635	ASN
1	B	1641	ASN
1	B	1659	HIS
1	B	1701	ASN
1	B	1824	ASN
1	B	1858	HIS
1	B	1912	ASN
1	B	1954	HIS
1	B	1973	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are 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	$\# Z > 2$
3	SAM	A	2409	-	27,29,29	0.47	0	34,42,42	0.46	0
3	SAM	B	2409	-	27,29,29	0.47	0	34,42,42	0.46	0

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
3	SAM	A	2409	-	-	2/17/33/33	0/3/3/3
3	SAM	B	2409	-	-	3/17/33/33	0/3/3/3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

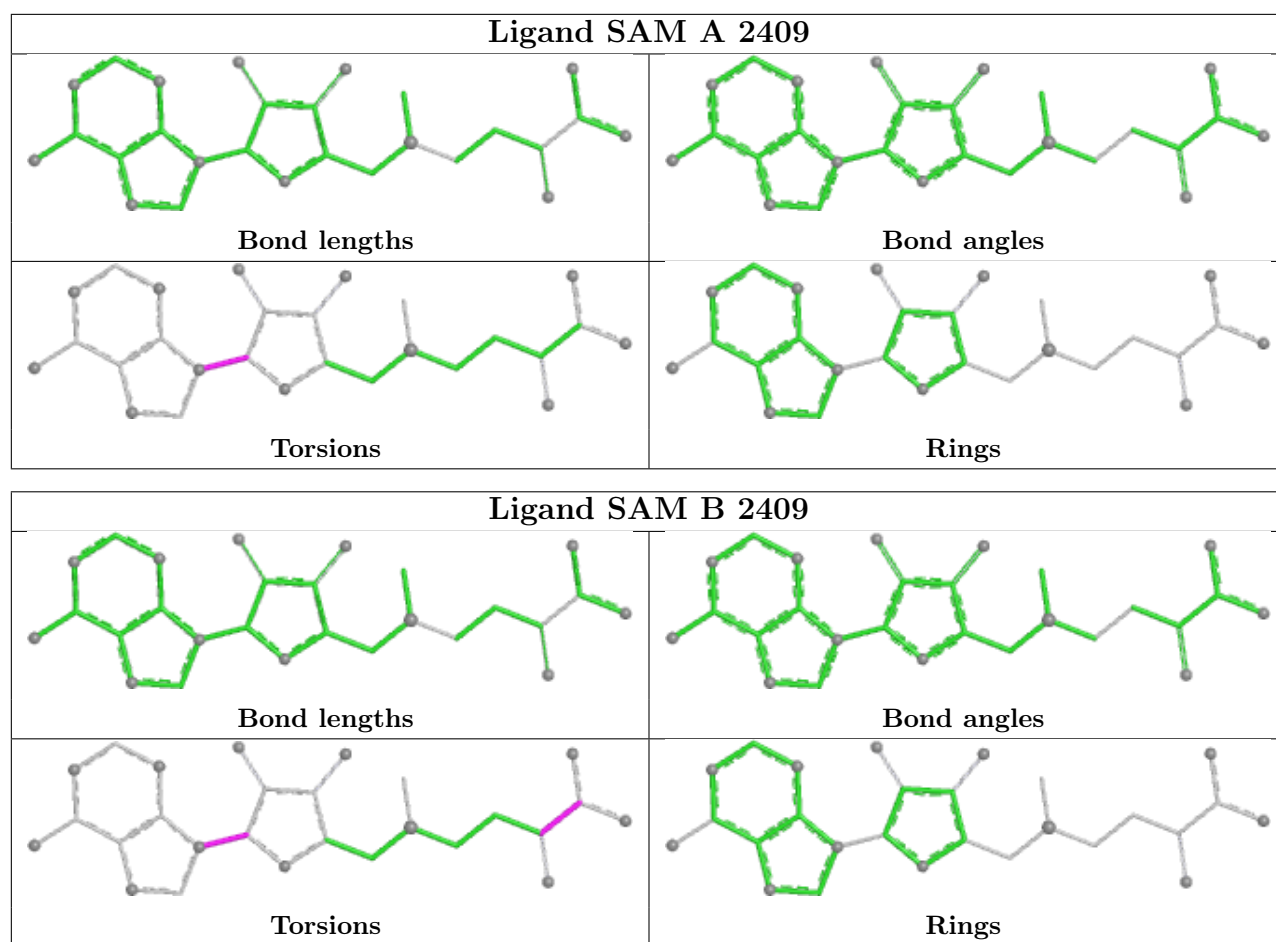
All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	2409	SAM	C2'-C1'-N9-C8
3	B	2409	SAM	OXT-C-CA-CB
3	B	2409	SAM	C2'-C1'-N9-C8
3	A	2409	SAM	O4'-C1'-N9-C8
3	B	2409	SAM	O-C-CA-CB

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	370/415 (89%)	0.15	11 (2%) 52 45	65, 85, 132, 173	0
1	B	376/415 (90%)	0.14	12 (3%) 50 42	63, 83, 124, 185	0
All	All	746/830 (89%)	0.14	23 (3%) 51 43	63, 84, 130, 185	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1922	SER	5.5
1	A	1936	ILE	4.8
1	A	2005	LYS	3.9
1	B	2005	LYS	3.8
1	B	1615	VAL	3.6
1	B	1920	LEU	3.3
1	B	1724	LEU	3.3
1	A	1642	TYR	3.2
1	A	1644	LYS	3.2
1	A	1724	LEU	3.2
1	A	1616	ILE	3.0
1	A	1920	LEU	2.6
1	B	1726	HIS	2.6
1	A	1643	ALA	2.6
1	B	1638	THR	2.3
1	B	1636	TYR	2.3
1	B	1643	ALA	2.2
1	A	1645	GLY	2.2
1	A	1921	GLN	2.2
1	A	1751	MET	2.1
1	B	1938	SER	2.1
1	B	1723	ASP	2.0
1	B	1729	PRO	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

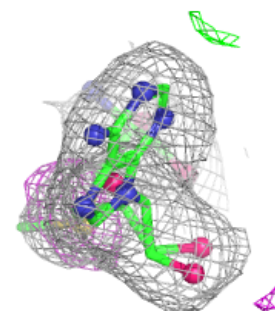
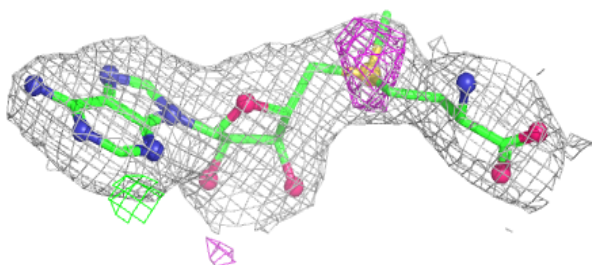
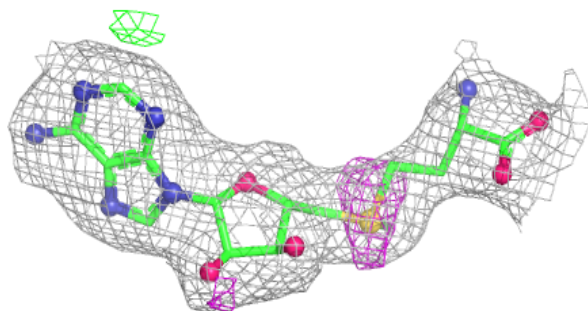
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SAM	A	2409	27/27	0.94	0.09	68,75,84,86	0
3	SAM	B	2409	27/27	0.94	0.09	69,80,92,94	0
2	ZN	A	2408	1/1	1.00	0.02	78,78,78,78	0
2	ZN	B	2408	1/1	1.00	0.02	77,77,77,77	0

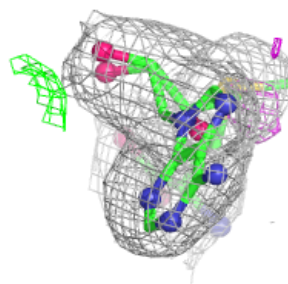
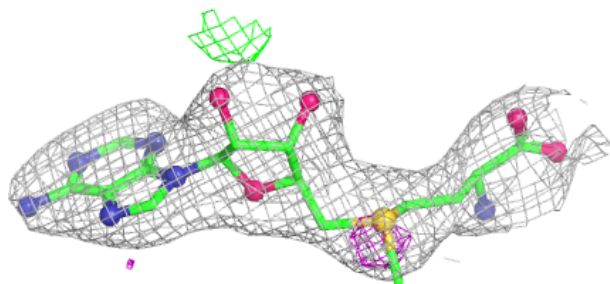
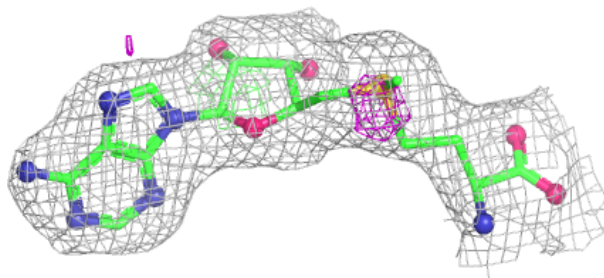
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around SAM A 2409:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around SAM B 2409:**

$2mF_o-DF_c$ (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.