



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 05:11 AM UTC

PDB ID : 8J2B / pdb_00008j2b
Title : Structure of the C-terminal subenzyme of the malonyl-CoA reductase from *Chloroflexus aurantiacus*, in complex with NADP⁺ and malonate
Authors : Ma, Q.; Liu, C.
Deposited on : 2023-04-14
Resolution : 2.13 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

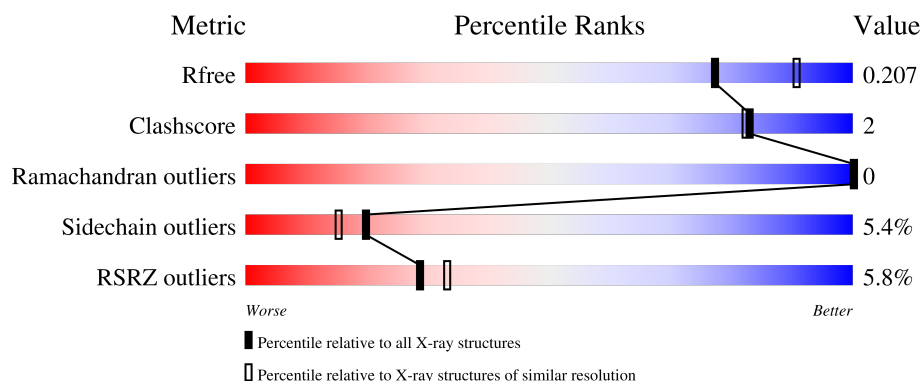
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.13 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	3689 (2.16-2.12)
Clashscore	190562	3812 (2.16-2.12)
Ramachandran outliers	187476	3773 (2.16-2.12)
Sidechain outliers	187428	3772 (2.16-2.12)
RSRZ outliers	180081	3691 (2.16-2.12)

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	674	6% 86% 11% ..

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	MLA	A	2002	-	X	-	-

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.

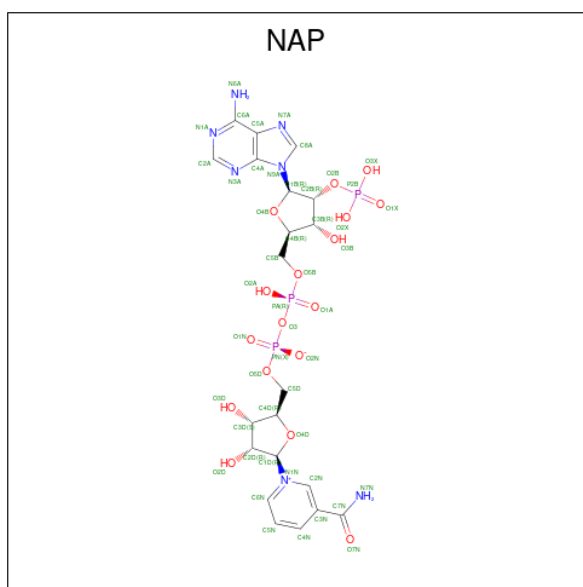
- Molecule 1 is a protein called Short-chain dehydrogenase/reductase SDR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	653	Total	C	N	O	S	0	0	0
			5040	3165	918	938	19			

There are 4 discrepancies between the modelled and reference sequences:

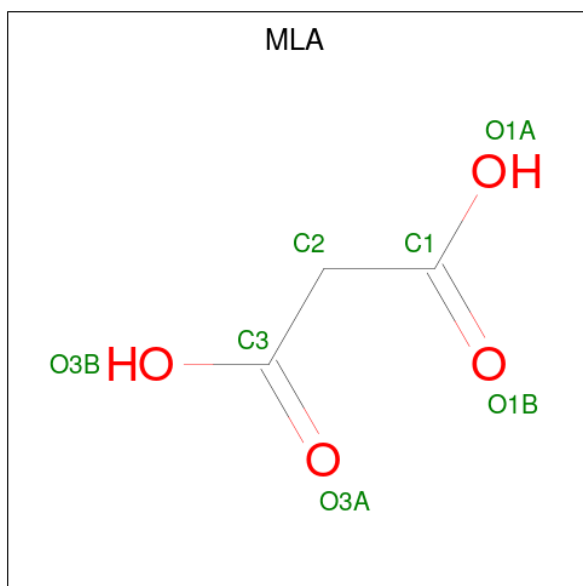
Chain	Residue	Modelled	Actual	Comment	Reference
A	546	GLY	-	expression tag	UNP A9WIU3
A	547	ALA	-	expression tag	UNP A9WIU3
A	548	MET	-	expression tag	UNP A9WIU3
A	549	ALA	-	expression tag	UNP A9WIU3

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is MALONIC ACID (CCD ID: MLA) (formula: $C_3H_4O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	3	4		

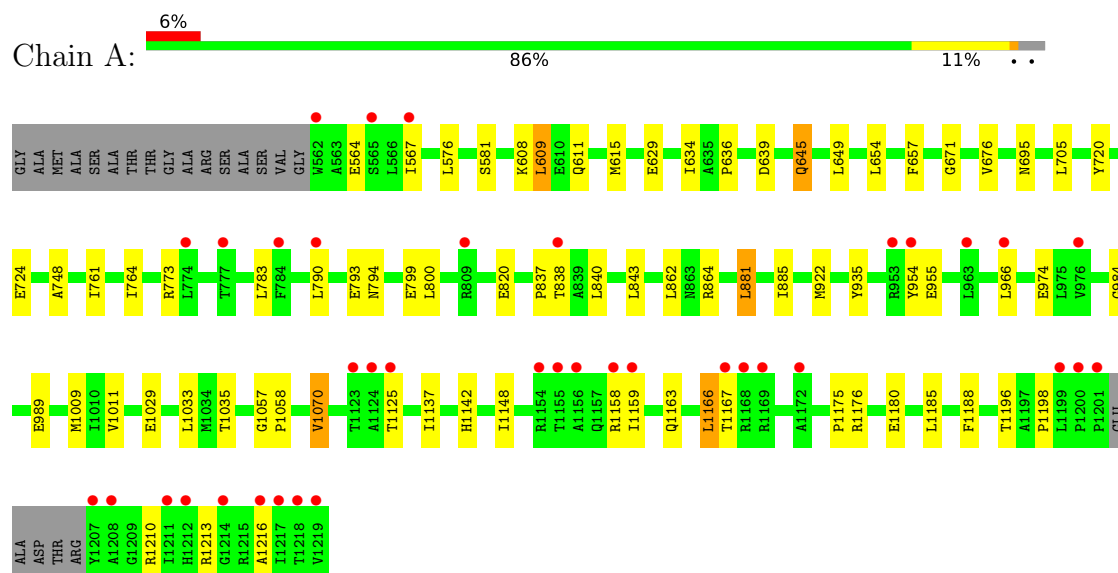
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	117	Total	O	0	0
			117	117		

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: Short-chain dehydrogenase/reductase SDR



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	100.96Å 125.22Å 73.81Å 90.00° 105.59° 90.00°	Depositor
Resolution (Å)	62.61 – 2.13 62.61 – 2.13	Depositor EDS
% Data completeness (in resolution range)	96.9 (62.61-2.13) 96.9 (62.61-2.13)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 2.12Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.185 , 0.203 0.188 , 0.207	Depositor DCC
R_{free} test set	2392 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	49.0	Xtriage
Anisotropy	0.217	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 37.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5212	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MLA, NAP

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	0.89	2/5134 (0.0%)	1.33	13/6963 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	922	MET	SD-CE	-7.31	1.61	1.79
1	A	1070	VAL	CA-C	5.93	1.58	1.52

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	793	GLU	CA-C-N	5.80	128.05	120.28
1	A	793	GLU	C-N-CA	5.80	128.05	120.28
1	A	1137	ILE	CA-C-N	5.62	128.13	120.54
1	A	1137	ILE	C-N-CA	5.62	128.13	120.54
1	A	1210	ARG	CA-C-N	5.33	127.38	120.56
1	A	1210	ARG	C-N-CA	5.33	127.38	120.56
1	A	1070	VAL	CB-CA-C	5.31	114.53	109.33
1	A	724	GLU	CB-CG-CD	5.26	121.54	112.60
1	A	657	PHE	CA-CB-CG	5.14	118.94	113.80
1	A	1057	GLY	N-CA-C	-5.10	101.94	112.34
1	A	764	ILE	N-CA-CB	5.02	120.50	111.21
1	A	671	GLY	CA-C-N	5.00	127.31	120.46
1	A	671	GLY	C-N-CA	5.00	127.31	120.46

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	5040	0	5070	24	0
2	A	48	0	25	1	0
3	A	7	0	2	0	0
4	A	117	0	0	1	0
All	All	5212	0	5097	24	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 2.

All (24) 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:581:SER:HB3	1:A:609:LEU:HD13	1.60	0.82
1:A:881:LEU:HD23	1:A:885:ILE:HD13	1.77	0.67
1:A:974:GLU:HB3	1:A:1198:PRO:HG2	1.82	0.61
1:A:790:LEU:O	1:A:794:ASN:HB2	2.06	0.56
1:A:935:TYR:HB2	4:A:2207:HOH:O	2.06	0.55
1:A:799:GLU:HB3	1:A:843:LEU:HD11	1.91	0.52
1:A:564:GLU:HA	1:A:567:ILE:HD12	1.92	0.51
1:A:1058:PRO:HB3	1:A:1196:THR:HB	1.93	0.51
1:A:1163:GLN:O	1:A:1216:ALA:HA	2.12	0.49
1:A:639:ASP:H	1:A:645:GLN:HE21	1.61	0.48
1:A:720:TYR:OH	1:A:954:TYR:HB3	2.13	0.48
1:A:837:PRO:HD2	1:A:840:LEU:HD12	1.95	0.47
1:A:1142:HIS:CD2	1:A:1216:ALA:CB	2.98	0.47
1:A:1009:MET:HB3	1:A:1035:THR:HG22	1.98	0.45
1:A:609:LEU:HG	1:A:634:ILE:HB	2.00	0.44
1:A:576:LEU:HD11	1:A:649:LEU:HD11	2.00	0.44
1:A:748:ALA:HA	1:A:761:ILE:HG12	2.00	0.44
1:A:1166:LEU:HD11	1:A:1188:PHE:CD1	2.53	0.43
1:A:1175:PRO:HB3	1:A:1180:GLU:HB3	2.01	0.42
1:A:676:VAL:HG11	1:A:1148:ILE:HD12	2.02	0.42
1:A:773:ARG:HD2	2:A:2001:NAP:O1N	2.20	0.42
1:A:609:LEU:HD23	1:A:636:PRO:HA	2.02	0.41
1:A:611:GLN:O	1:A:615:MET:HG3	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:984:GLY:O	1:A:1011:VAL:HA	2.22	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	649/674 (96%)	636 (98%)	13 (2%)	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	522/534 (98%)	494 (95%)	28 (5%)	20	15

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	608	LYS
1	A	609	LEU
1	A	629	GLU
1	A	645	GLN
1	A	654	LEU

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Mol	Chain	Res	Type
1	A	695	ASN
1	A	705	LEU
1	A	783	LEU
1	A	800	LEU
1	A	820	GLU
1	A	838	THR
1	A	862	LEU
1	A	864	ARG
1	A	881	LEU
1	A	955	GLU
1	A	966	LEU
1	A	989	GLU
1	A	1029	GLU
1	A	1033	LEU
1	A	1070	VAL
1	A	1125	THR
1	A	1158	ARG
1	A	1159	ILE
1	A	1166	LEU
1	A	1167	THR
1	A	1176	ARG
1	A	1185	LEU
1	A	1213	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	607	HIS
1	A	611	GLN
1	A	645	GLN
1	A	733	ASN
1	A	816	HIS
1	A	986	HIS
1	A	1162	ASN

5.3.3 RNA ⓘ

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](#)

2 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).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAP	A	2001	-	50,52,52	0.39	0	71,80,80	0.72	2 (2%)
3	MLA	A	2002	-	6,6,6	2.37	4 (66%)	7,7,7	1.91	4 (57%)

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	NAP	A	2001	-	-	6/35/67/67	0/5/5/5
3	MLA	A	2002	-	-	1/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	2002	MLA	O1B-C1	3.04	1.32	1.22
3	A	2002	MLA	O1A-C1	-3.02	1.20	1.30
3	A	2002	MLA	O3B-C3	-2.97	1.21	1.30
3	A	2002	MLA	O3A-C3	2.57	1.30	1.22

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2001	NAP	O2B-C2B-C1B	2.73	119.65	110.05
3	A	2002	MLA	O1A-C1-C2	2.53	122.36	114.51
3	A	2002	MLA	O3B-C3-C2	2.50	122.26	114.51
3	A	2002	MLA	O1B-C1-C2	-2.50	115.00	122.11
3	A	2002	MLA	O3A-C3-C2	-2.48	115.05	122.11
2	A	2001	NAP	O2A-PA-O1A	2.08	122.10	112.44

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	2001	NAP	C5D-O5D-PN-O3
2	A	2001	NAP	C5D-O5D-PN-O2N
2	A	2001	NAP	O4D-C1D-N1N-C2N
2	A	2001	NAP	C5D-O5D-PN-O1N
2	A	2001	NAP	C1B-C2B-O2B-P2B
2	A	2001	NAP	O4D-C1D-N1N-C6N
3	A	2002	MLA	C1-C2-C3-O3A

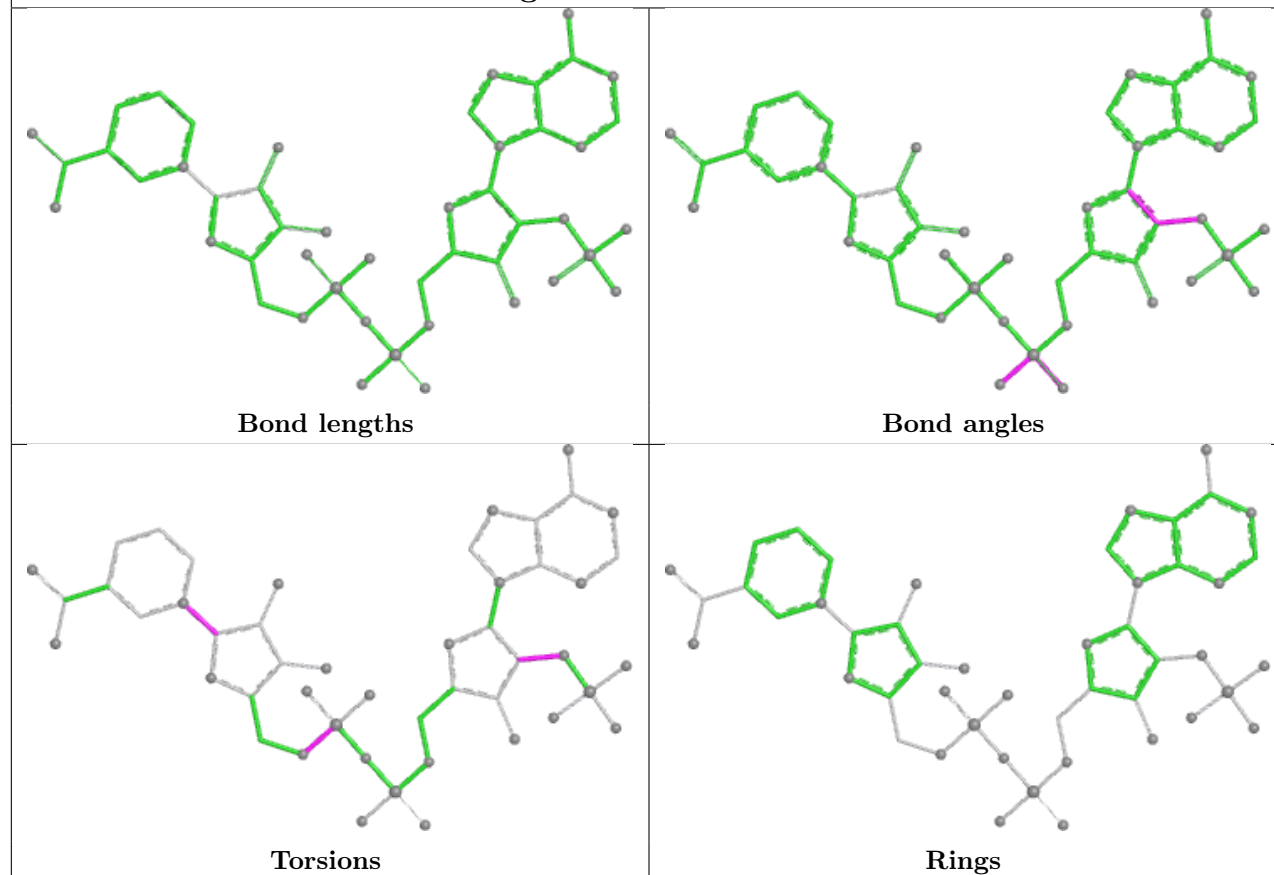
There are no ring outliers.

1 monomer is involved in 1 short contact:

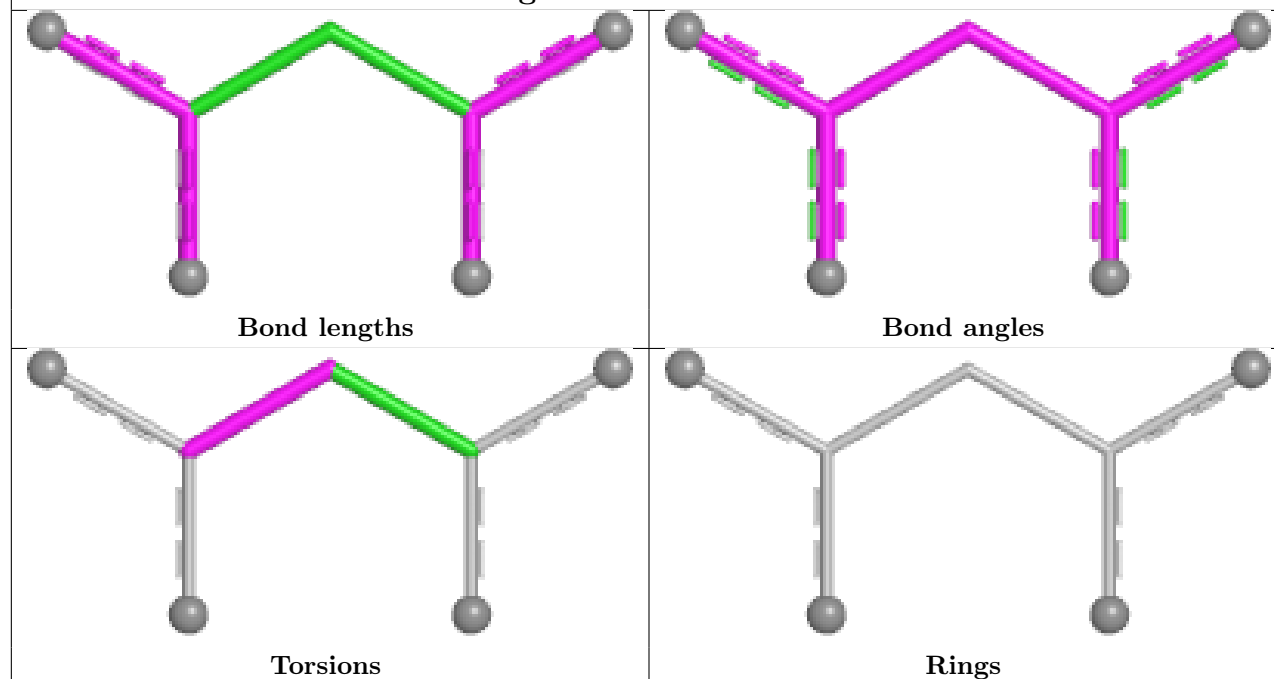
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2001	NAP	1	0

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 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.

Ligand NAP A 2001



Ligand MLA A 2002



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	653/674 (96%)	0.37	38 (5%) 29 33	36, 57, 91, 117	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1217	ILE	5.2
1	A	1172	ALA	4.7
1	A	966	LEU	4.2
1	A	1124	ALA	4.0
1	A	1201	PRO	4.0
1	A	1211	ILE	3.5
1	A	562	TRP	3.5
1	A	1199	LEU	3.4
1	A	1214	GLY	3.3
1	A	1218	THR	3.2
1	A	963	LEU	3.2
1	A	1208	ALA	3.1
1	A	1156	ALA	3.0
1	A	954	TYR	2.9
1	A	1216	ALA	2.8
1	A	1212	HIS	2.8
1	A	1168	ARG	2.7
1	A	1167	THR	2.7
1	A	1169	ARG	2.7
1	A	777	THR	2.6
1	A	1155	THR	2.6
1	A	1159	ILE	2.6
1	A	1207	TYR	2.5
1	A	774	LEU	2.5
1	A	1219	VAL	2.5
1	A	1200	PRO	2.5
1	A	953	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	567	ILE	2.4
1	A	565	SER	2.4
1	A	976	VAL	2.3
1	A	838	THR	2.2
1	A	1125	THR	2.2
1	A	790	LEU	2.1
1	A	1123	THR	2.1
1	A	784	PHE	2.1
1	A	1158	ARG	2.1
1	A	809	ARG	2.1
1	A	1154	ARG	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](#)

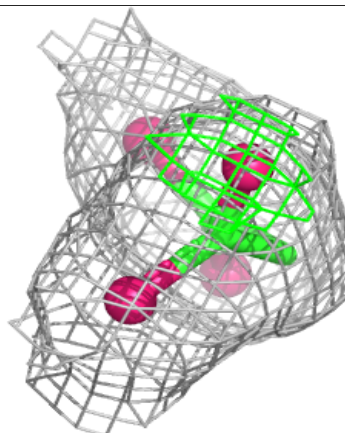
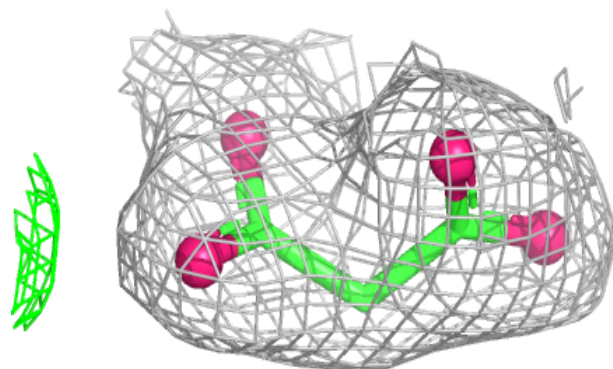
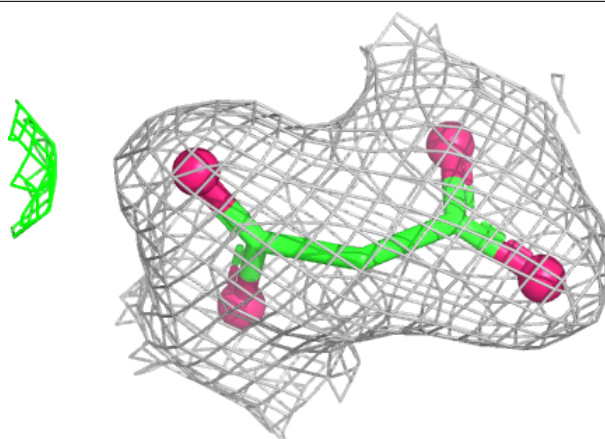
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	MLA	A	2002	7/7	0.96	0.06	43,44,45,46	0
2	NAP	A	2001	48/48	0.97	0.06	36,41,51,56	0

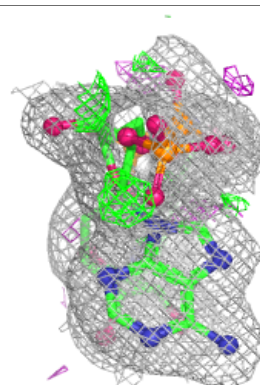
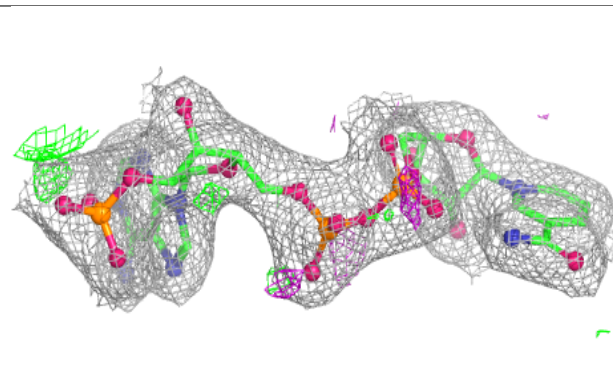
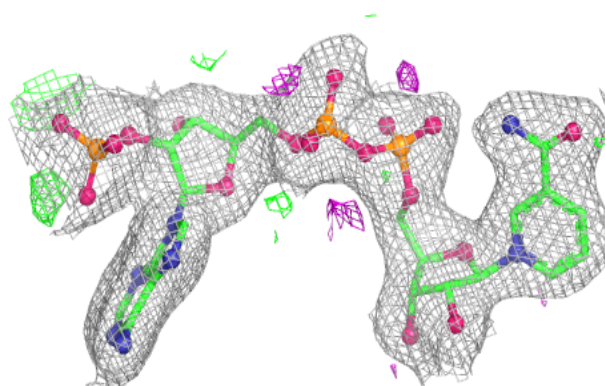
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around MLA A 2002:

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

**Electron density around NAP A 2001:**

$2mF_o-DF_c$ (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.