Full wwPDB X-ray Structure Validation Report (i)

Oct 31, 2023 – 05:10 PM JST

PDB ID	:	5GWT
Title	:	4-hydroxyisoleucine dehydrogenase mutant complexed with NADH and succinate
Authors	:	Shi, X.; Miyakawa, T.; Nakamura, A.; Tanokura, M.
Deposited on	:	2016-09-13
Resolution	:	1.90 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 (i)) were used in the production of this report:

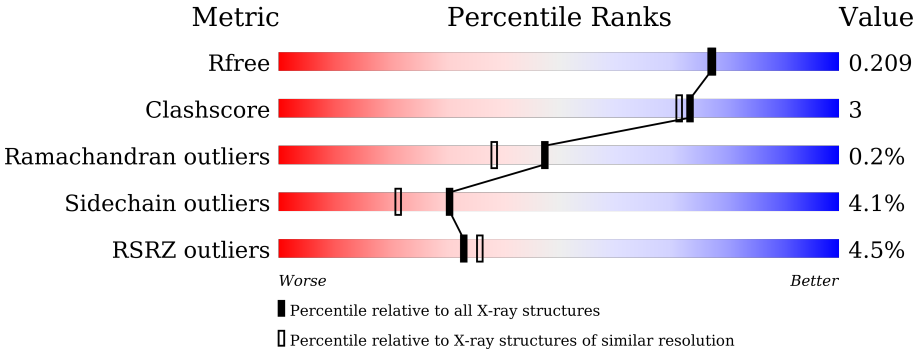
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (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-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	A	278	% 77% 9% •• 13%
1	B	278	4% 78% 8% •• 13%
1	C	278	4% 74% 11% •• 13%
1	D	278	7% 79% 6% •• 13%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SIN	C	301	-	X	-	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 8115 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 4-hydroxyisolecuine dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	243	1881	1198	318	352	13	0	0	0
1	B	243	1881	1198	318	352	13	0	0	0
1	C	243	1881	1198	318	352	13	0	0	0
1	D	243	1881	1198	318	352	13	0	0	0

There are 144 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	144	LYS	GLU	engineered mutation	UNP A0A0K0Q8K4
A	242	GLN	TRP	engineered mutation	UNP A0A0K0Q8K4
A	249	LEU	-	expression tag	UNP A0A0K0Q8K4
A	250	GLU	-	expression tag	UNP A0A0K0Q8K4
A	251	LYS	-	expression tag	UNP A0A0K0Q8K4
A	252	GLY	-	expression tag	UNP A0A0K0Q8K4
A	253	GLU	-	expression tag	UNP A0A0K0Q8K4
A	254	LEU	-	expression tag	UNP A0A0K0Q8K4
A	255	ASN	-	expression tag	UNP A0A0K0Q8K4
A	256	SER	-	expression tag	UNP A0A0K0Q8K4
A	257	LYS	-	expression tag	UNP A0A0K0Q8K4
A	258	LEU	-	expression tag	UNP A0A0K0Q8K4
A	259	GLU	-	expression tag	UNP A0A0K0Q8K4
A	260	GLY	-	expression tag	UNP A0A0K0Q8K4
A	261	LYS	-	expression tag	UNP A0A0K0Q8K4
A	262	PRO	-	expression tag	UNP A0A0K0Q8K4
A	263	ILE	-	expression tag	UNP A0A0K0Q8K4
A	264	PRO	-	expression tag	UNP A0A0K0Q8K4
A	265	ASN	-	expression tag	UNP A0A0K0Q8K4
A	266	PRO	-	expression tag	UNP A0A0K0Q8K4
A	267	LEU	-	expression tag	UNP A0A0K0Q8K4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	268	LEU	-	expression tag	UNP A0A0K0Q8K4
A	269	GLY	-	expression tag	UNP A0A0K0Q8K4
A	270	LEU	-	expression tag	UNP A0A0K0Q8K4
A	271	ASP	-	expression tag	UNP A0A0K0Q8K4
A	272	SER	-	expression tag	UNP A0A0K0Q8K4
A	273	THR	-	expression tag	UNP A0A0K0Q8K4
A	274	ARG	-	expression tag	UNP A0A0K0Q8K4
A	275	THR	-	expression tag	UNP A0A0K0Q8K4
A	276	GLY	-	expression tag	UNP A0A0K0Q8K4
A	277	HIS	-	expression tag	UNP A0A0K0Q8K4
A	278	HIS	-	expression tag	UNP A0A0K0Q8K4
A	279	HIS	-	expression tag	UNP A0A0K0Q8K4
A	280	HIS	-	expression tag	UNP A0A0K0Q8K4
A	281	HIS	-	expression tag	UNP A0A0K0Q8K4
A	282	HIS	-	expression tag	UNP A0A0K0Q8K4
B	144	LYS	GLU	engineered mutation	UNP A0A0K0Q8K4
B	242	GLN	TRP	engineered mutation	UNP A0A0K0Q8K4
B	249	LEU	-	expression tag	UNP A0A0K0Q8K4
B	250	GLU	-	expression tag	UNP A0A0K0Q8K4
B	251	LYS	-	expression tag	UNP A0A0K0Q8K4
B	252	GLY	-	expression tag	UNP A0A0K0Q8K4
B	253	GLU	-	expression tag	UNP A0A0K0Q8K4
B	254	LEU	-	expression tag	UNP A0A0K0Q8K4
B	255	ASN	-	expression tag	UNP A0A0K0Q8K4
B	256	SER	-	expression tag	UNP A0A0K0Q8K4
B	257	LYS	-	expression tag	UNP A0A0K0Q8K4
B	258	LEU	-	expression tag	UNP A0A0K0Q8K4
B	259	GLU	-	expression tag	UNP A0A0K0Q8K4
B	260	GLY	-	expression tag	UNP A0A0K0Q8K4
B	261	LYS	-	expression tag	UNP A0A0K0Q8K4
B	262	PRO	-	expression tag	UNP A0A0K0Q8K4
B	263	ILE	-	expression tag	UNP A0A0K0Q8K4
B	264	PRO	-	expression tag	UNP A0A0K0Q8K4
B	265	ASN	-	expression tag	UNP A0A0K0Q8K4
B	266	PRO	-	expression tag	UNP A0A0K0Q8K4
B	267	LEU	-	expression tag	UNP A0A0K0Q8K4
B	268	LEU	-	expression tag	UNP A0A0K0Q8K4
B	269	GLY	-	expression tag	UNP A0A0K0Q8K4
B	270	LEU	-	expression tag	UNP A0A0K0Q8K4
B	271	ASP	-	expression tag	UNP A0A0K0Q8K4
B	272	SER	-	expression tag	UNP A0A0K0Q8K4
B	273	THR	-	expression tag	UNP A0A0K0Q8K4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	274	ARG	-	expression tag	UNP A0A0K0Q8K4
B	275	THR	-	expression tag	UNP A0A0K0Q8K4
B	276	GLY	-	expression tag	UNP A0A0K0Q8K4
B	277	HIS	-	expression tag	UNP A0A0K0Q8K4
B	278	HIS	-	expression tag	UNP A0A0K0Q8K4
B	279	HIS	-	expression tag	UNP A0A0K0Q8K4
B	280	HIS	-	expression tag	UNP A0A0K0Q8K4
B	281	HIS	-	expression tag	UNP A0A0K0Q8K4
B	282	HIS	-	expression tag	UNP A0A0K0Q8K4
C	144	LYS	GLU	engineered mutation	UNP A0A0K0Q8K4
C	242	GLN	TRP	engineered mutation	UNP A0A0K0Q8K4
C	249	LEU	-	expression tag	UNP A0A0K0Q8K4
C	250	GLU	-	expression tag	UNP A0A0K0Q8K4
C	251	LYS	-	expression tag	UNP A0A0K0Q8K4
C	252	GLY	-	expression tag	UNP A0A0K0Q8K4
C	253	GLU	-	expression tag	UNP A0A0K0Q8K4
C	254	LEU	-	expression tag	UNP A0A0K0Q8K4
C	255	ASN	-	expression tag	UNP A0A0K0Q8K4
C	256	SER	-	expression tag	UNP A0A0K0Q8K4
C	257	LYS	-	expression tag	UNP A0A0K0Q8K4
C	258	LEU	-	expression tag	UNP A0A0K0Q8K4
C	259	GLU	-	expression tag	UNP A0A0K0Q8K4
C	260	GLY	-	expression tag	UNP A0A0K0Q8K4
C	261	LYS	-	expression tag	UNP A0A0K0Q8K4
C	262	PRO	-	expression tag	UNP A0A0K0Q8K4
C	263	ILE	-	expression tag	UNP A0A0K0Q8K4
C	264	PRO	-	expression tag	UNP A0A0K0Q8K4
C	265	ASN	-	expression tag	UNP A0A0K0Q8K4
C	266	PRO	-	expression tag	UNP A0A0K0Q8K4
C	267	LEU	-	expression tag	UNP A0A0K0Q8K4
C	268	LEU	-	expression tag	UNP A0A0K0Q8K4
C	269	GLY	-	expression tag	UNP A0A0K0Q8K4
C	270	LEU	-	expression tag	UNP A0A0K0Q8K4
C	271	ASP	-	expression tag	UNP A0A0K0Q8K4
C	272	SER	-	expression tag	UNP A0A0K0Q8K4
C	273	THR	-	expression tag	UNP A0A0K0Q8K4
C	274	ARG	-	expression tag	UNP A0A0K0Q8K4
C	275	THR	-	expression tag	UNP A0A0K0Q8K4
C	276	GLY	-	expression tag	UNP A0A0K0Q8K4
C	277	HIS	-	expression tag	UNP A0A0K0Q8K4
C	278	HIS	-	expression tag	UNP A0A0K0Q8K4
C	279	HIS	-	expression tag	UNP A0A0K0Q8K4

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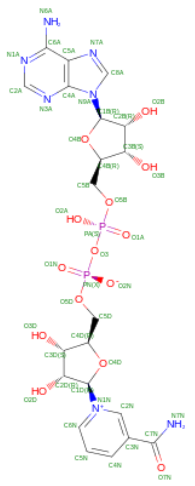
Chain	Residue	Modelled	Actual	Comment	Reference
C	280	HIS	-	expression tag	UNP A0A0K0Q8K4
C	281	HIS	-	expression tag	UNP A0A0K0Q8K4
C	282	HIS	-	expression tag	UNP A0A0K0Q8K4
D	144	LYS	GLU	engineered mutation	UNP A0A0K0Q8K4
D	242	GLN	TRP	engineered mutation	UNP A0A0K0Q8K4
D	249	LEU	-	expression tag	UNP A0A0K0Q8K4
D	250	GLU	-	expression tag	UNP A0A0K0Q8K4
D	251	LYS	-	expression tag	UNP A0A0K0Q8K4
D	252	GLY	-	expression tag	UNP A0A0K0Q8K4
D	253	GLU	-	expression tag	UNP A0A0K0Q8K4
D	254	LEU	-	expression tag	UNP A0A0K0Q8K4
D	255	ASN	-	expression tag	UNP A0A0K0Q8K4
D	256	SER	-	expression tag	UNP A0A0K0Q8K4
D	257	LYS	-	expression tag	UNP A0A0K0Q8K4
D	258	LEU	-	expression tag	UNP A0A0K0Q8K4
D	259	GLU	-	expression tag	UNP A0A0K0Q8K4
D	260	GLY	-	expression tag	UNP A0A0K0Q8K4
D	261	LYS	-	expression tag	UNP A0A0K0Q8K4
D	262	PRO	-	expression tag	UNP A0A0K0Q8K4
D	263	ILE	-	expression tag	UNP A0A0K0Q8K4
D	264	PRO	-	expression tag	UNP A0A0K0Q8K4
D	265	ASN	-	expression tag	UNP A0A0K0Q8K4
D	266	PRO	-	expression tag	UNP A0A0K0Q8K4
D	267	LEU	-	expression tag	UNP A0A0K0Q8K4
D	268	LEU	-	expression tag	UNP A0A0K0Q8K4
D	269	GLY	-	expression tag	UNP A0A0K0Q8K4
D	270	LEU	-	expression tag	UNP A0A0K0Q8K4
D	271	ASP	-	expression tag	UNP A0A0K0Q8K4
D	272	SER	-	expression tag	UNP A0A0K0Q8K4
D	273	THR	-	expression tag	UNP A0A0K0Q8K4
D	274	ARG	-	expression tag	UNP A0A0K0Q8K4
D	275	THR	-	expression tag	UNP A0A0K0Q8K4
D	276	GLY	-	expression tag	UNP A0A0K0Q8K4
D	277	HIS	-	expression tag	UNP A0A0K0Q8K4
D	278	HIS	-	expression tag	UNP A0A0K0Q8K4
D	279	HIS	-	expression tag	UNP A0A0K0Q8K4
D	280	HIS	-	expression tag	UNP A0A0K0Q8K4
D	281	HIS	-	expression tag	UNP A0A0K0Q8K4
D	282	HIS	-	expression tag	UNP A0A0K0Q8K4

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).

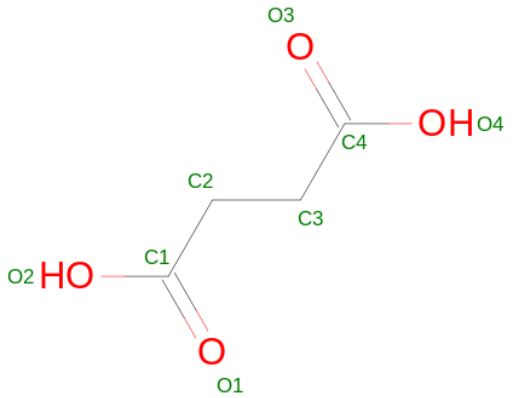
NAD



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	44	21	7	14	2	0	0
2	B	1	44	21	7	14	2	0	0
2	C	1	44	21	7	14	2	0	0
2	D	1	44	21	7	14	2	0	0

- Molecule 3 is SUCCINIC ACID (three-letter code: SIN) (formula: $C_4H_6O_4$).

SIN



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	8	4	4	0	0
3	B	1	8	4	4	0	0
3	C	1	8	4	4	0	0
3	D	1	8	4	4	0	0

- Molecule 4 is water.

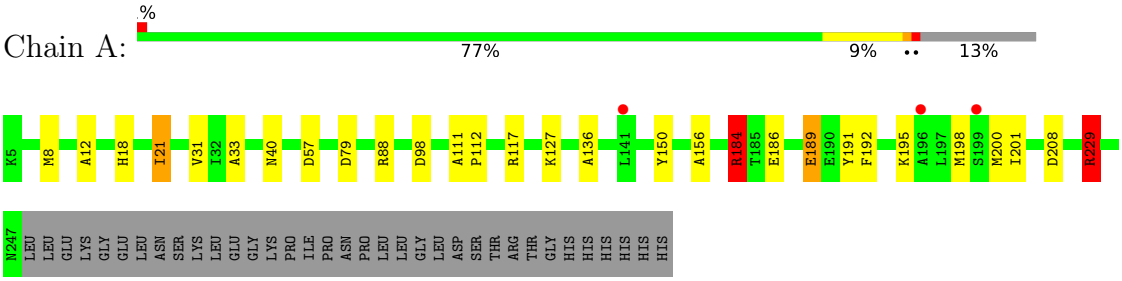
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	117	117	117	0	0
4	B	108	108	108	0	0
4	C	86	86	86	0	0
4	D	72	72	72	0	0

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.

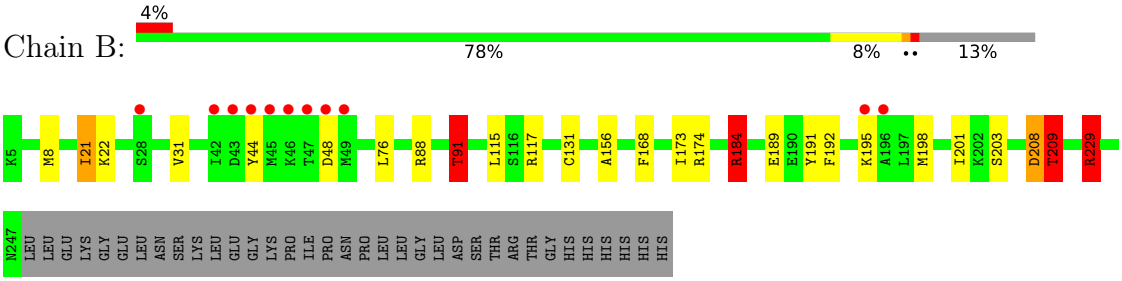
• Molecule 1: 4-hydroxyisolecuine dehydrogenase

Chain A:



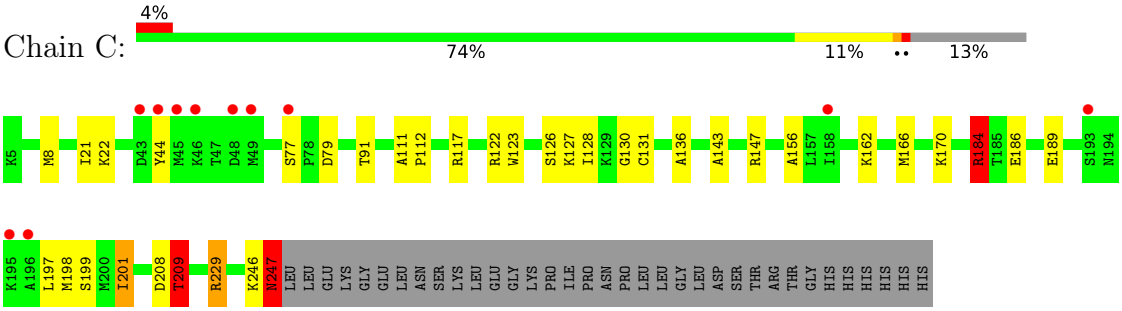
• Molecule 1: 4-hydroxyisolecuine dehydrogenase

Chain B:



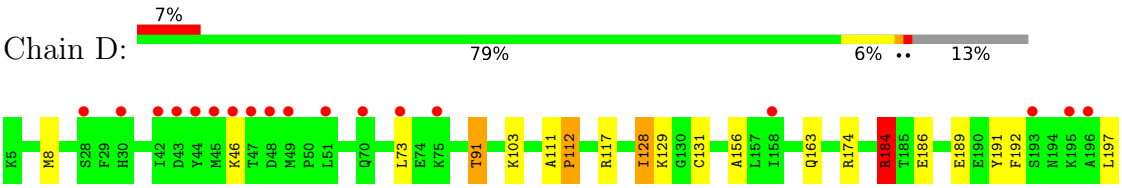
• Molecule 1: 4-hydroxyisolecuine dehydrogenase

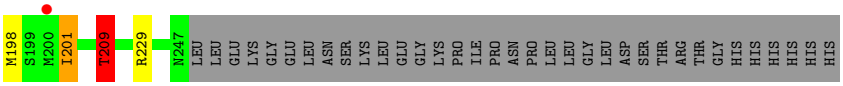
Chain C:



• Molecule 1: 4-hydroxyisolecuine dehydrogenase

Chain D:




M198
S199
M200
I201
T209
R229
N247
LEU
LEU
GLU
LYS
GLY
GLU
LEU
ASN
SER
LYS
LEU
GLU
GLY
LYS
PRO
ILE
PRO
ASN
PRO
LEU
LEU
GLY
LEU
ASP
SER
THR
ARG
THR
GLY
HIS
HIS
HIS
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HIS
HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α, β, γ	231.85Å 231.85Å 78.96Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.50 – 1.90 45.51 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.7 (45.50-1.90) 99.7 (45.51-1.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.24 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.8.0131	Depositor
R, R_{free}	0.174 , 0.201 0.184 , 0.209	Depositor DCC
R_{free} test set	2100 reflections (1.68%)	wwPDB-VP
Wilson B-factor (Å^2)	23.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 43.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.012 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	8115	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical 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.

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: SIN, NAD

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	1.20	4/1910 (0.2%)	1.28	18/2578 (0.7%)
1	B	1.09	2/1910 (0.1%)	1.15	9/2578 (0.3%)
1	C	1.10	2/1910 (0.1%)	1.16	15/2578 (0.6%)
1	D	1.08	2/1910 (0.1%)	1.17	11/2578 (0.4%)
All	All	1.12	10/7640 (0.1%)	1.19	53/10312 (0.5%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	131	CYS	CB-SG	-7.81	1.69	1.82
1	C	131	CYS	CB-SG	-6.51	1.71	1.82
1	A	150	TYR	CE2-CZ	6.39	1.46	1.38
1	A	40	ASN	CB-CG	-6.07	1.37	1.51
1	B	131	CYS	CB-SG	-6.05	1.72	1.82
1	D	186	GLU	CD-OE1	5.59	1.31	1.25
1	B	91	THR	CB-CG2	-5.24	1.35	1.52
1	A	184	ARG	CD-NE	-5.17	1.37	1.46
1	C	186	GLU	CD-OE1	5.10	1.31	1.25
1	A	186	GLU	CD-OE1	5.03	1.31	1.25

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	184	ARG	NE-CZ-NH1	19.60	130.10	120.30
1	A	229	ARG	NE-CZ-NH1	16.58	128.59	120.30
1	D	229	ARG	NE-CZ-NH1	16.25	128.42	120.30
1	B	184	ARG	NE-CZ-NH1	15.24	127.92	120.30
1	A	184	ARG	NE-CZ-NH2	-15.20	112.70	120.30
1	A	229	ARG	NE-CZ-NH2	-14.26	113.17	120.30
1	C	184	ARG	NE-CZ-NH1	13.43	127.02	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	229	ARG	NE-CZ-NH2	-13.38	113.61	120.30
1	C	229	ARG	NE-CZ-NH1	13.36	126.98	120.30
1	B	229	ARG	NE-CZ-NH1	12.97	126.78	120.30
1	D	184	ARG	NE-CZ-NH1	12.61	126.60	120.30
1	C	184	ARG	NE-CZ-NH2	-12.39	114.11	120.30
1	C	8	MET	CG-SD-CE	-12.13	80.80	100.20
1	B	184	ARG	NE-CZ-NH2	-12.10	114.25	120.30
1	D	184	ARG	NE-CZ-NH2	-12.07	114.26	120.30
1	B	229	ARG	NE-CZ-NH2	-12.06	114.27	120.30
1	C	229	ARG	NE-CZ-NH2	-8.62	115.99	120.30
1	D	8	MET	CG-SD-CE	-8.27	86.96	100.20
1	D	174	ARG	NE-CZ-NH1	7.41	124.01	120.30
1	A	184	ARG	CB-CG-CD	7.20	130.31	111.60
1	B	208	ASP	CB-CG-OD2	-7.05	111.96	118.30
1	A	57	ASP	CB-CG-OD1	6.90	124.51	118.30
1	C	162	LYS	CD-CE-NZ	6.71	127.14	111.70
1	C	79	ASP	CB-CG-OD1	-6.63	112.34	118.30
1	D	128	ILE	CB-CA-C	-6.47	98.65	111.60
1	C	117	ARG	NE-CZ-NH1	6.43	123.52	120.30
1	C	147	ARG	NE-CZ-NH1	6.38	123.49	120.30
1	D	117	ARG	NE-CZ-NH2	-6.27	117.16	120.30
1	B	209	THR	CB-CA-C	-6.16	94.97	111.60
1	B	174	ARG	NE-CZ-NH1	6.05	123.33	120.30
1	A	88	ARG	NE-CZ-NH2	-6.04	117.28	120.30
1	A	88	ARG	NE-CZ-NH1	5.93	123.26	120.30
1	C	247	ASN	CA-C-O	-5.81	107.90	120.10
1	A	57	ASP	CB-CG-OD2	-5.78	113.09	118.30
1	C	247	ASN	N-CA-CB	5.69	120.84	110.60
1	B	117	ARG	NE-CZ-NH1	5.68	123.14	120.30
1	B	88	ARG	NE-CZ-NH1	5.67	123.14	120.30
1	C	209	THR	CB-CA-C	-5.65	96.35	111.60
1	A	184	ARG	CD-NE-CZ	5.57	131.40	123.60
1	D	174	ARG	NE-CZ-NH2	-5.57	117.52	120.30
1	A	21	ILE	CA-CB-CG2	5.53	121.95	110.90
1	C	247	ASN	N-CA-C	5.53	125.92	111.00
1	D	209	THR	CB-CA-C	-5.39	97.05	111.60
1	D	117	ARG	NE-CZ-NH1	5.34	122.97	120.30
1	A	8	MET	CG-SD-CE	-5.33	91.68	100.20
1	A	98	ASP	CB-CG-OD2	-5.29	113.53	118.30
1	C	122	ARG	NE-CZ-NH1	5.28	122.94	120.30
1	A	117	ARG	NE-CZ-NH2	-5.27	117.66	120.30
1	C	117	ARG	NE-CZ-NH2	-5.26	117.67	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	79	ASP	CB-CG-OD1	-5.23	113.59	118.30
1	A	189	GLU	OE1-CD-OE2	-5.15	117.12	123.30
1	A	117	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	A	229	ARG	CD-NE-CZ	5.13	130.79	123.60

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	1881	0	1935	13	0
1	B	1881	0	1935	17	0
1	C	1881	0	1935	15	0
1	D	1881	0	1935	14	0
2	A	44	0	26	0	0
2	B	44	0	26	0	0
2	C	44	0	26	0	0
2	D	44	0	26	0	0
3	A	8	0	4	0	0
3	B	8	0	4	0	0
3	C	8	0	4	0	0
3	D	8	0	4	0	0
4	A	117	0	0	2	0
4	B	108	0	0	1	0
4	C	86	0	0	0	0
4	D	72	0	0	1	0
All	All	8115	0	7860	52	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 (52) 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:191:TYR:HB3	1:A:198:MET:HE1	1.44	1.00
1:B:8:MET:HE3	1:B:115:LEU:HD22	1.54	0.88
1:A:184:ARG:HD2	1:A:189:GLU:OE1	1.75	0.86
1:B:8:MET:CE	1:B:115:LEU:HD22	2.12	0.79
1:B:184:ARG:HD2	1:B:189:GLU:OE1	1.83	0.78
1:D:184:ARG:HD2	1:D:189:GLU:OE1	1.85	0.76
1:C:198:MET:HA	1:C:198:MET:HE2	1.67	0.76
1:B:91:THR:HG21	4:B:503:HOH:O	1.88	0.74
1:A:191:TYR:CB	1:A:198:MET:HE1	2.16	0.73
1:D:91:THR:HG21	4:D:469:HOH:O	1.88	0.73
1:D:198:MET:HA	1:D:198:MET:HE2	1.71	0.72
1:B:209:THR:HG21	1:C:229:ARG:NH2	2.09	0.66
1:B:21:ILE:HD12	1:B:31:VAL:HG11	1.77	0.65
1:B:191:TYR:HB3	1:B:198:MET:HE1	1.79	0.64
1:B:229:ARG:NH2	1:C:209:THR:HG21	2.13	0.63
1:D:128:ILE:HG22	1:D:129:LYS:O	1.99	0.63
1:C:156:ALA:HB2	1:D:156:ALA:HB2	1.84	0.59
1:A:18:HIS:HE1	4:A:468:HOH:O	1.87	0.57
1:C:184:ARG:HD2	1:C:189:GLU:OE1	2.04	0.57
1:D:184:ARG:HB2	1:D:192:PHE:CZ	2.41	0.56
1:B:198:MET:HA	1:B:198:MET:HE2	1.87	0.55
1:A:229:ARG:NH2	1:D:209:THR:HG21	2.26	0.51
1:C:198:MET:HA	1:C:198:MET:CE	2.40	0.49
1:A:156:ALA:HB2	1:B:156:ALA:HB2	1.95	0.49
1:B:209:THR:HG21	1:C:229:ARG:HH22	1.78	0.49
1:D:197:LEU:O	1:D:201:ILE:HD13	2.13	0.49
1:C:246:LYS:O	1:C:247:ASN:HB2	2.12	0.48
1:B:191:TYR:HB3	1:B:198:MET:CE	2.45	0.47
1:C:111:ALA:HB3	1:C:112:PRO:HD3	1.96	0.46
1:D:191:TYR:HB3	1:D:198:MET:CE	2.47	0.45
1:B:21:ILE:HD12	1:B:31:VAL:CG1	2.44	0.45
1:C:143:ALA:HB1	1:D:163:GLN:HA	1.99	0.44
1:C:197:LEU:O	1:C:201:ILE:HD13	2.18	0.44
1:D:111:ALA:HB3	1:D:112:PRO:HD3	2.00	0.44
1:C:166:MET:HE3	1:C:166:MET:HB3	1.91	0.43
1:A:200:MET:HG3	4:A:508:HOH:O	2.17	0.43
1:A:21:ILE:HG13	1:A:31:VAL:HG11	2.01	0.43
1:A:111:ALA:HB3	1:A:112:PRO:HD3	2.00	0.43
1:B:22:LYS:HG2	1:B:44:TYR:OH	2.19	0.43
1:A:191:TYR:HB3	1:A:198:MET:CE	2.32	0.42
1:B:184:ARG:HD3	1:B:192:PHE:CD1	2.54	0.42
1:B:198:MET:HA	1:B:198:MET:CE	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:123:TRP:CD2	1:C:130:GLY:HA3	2.55	0.41
1:D:128:ILE:HD13	1:D:128:ILE:HG21	1.85	0.41
1:A:12:ALA:CB	1:A:33:ALA:HB1	2.50	0.41
1:C:22:LYS:CD	1:C:44:TYR:OH	2.67	0.41
1:B:168:PHE:HB3	1:B:173:ILE:HB	2.02	0.41
1:C:126:SER:HB2	1:C:128:ILE:HG12	2.02	0.41
1:A:198:MET:CE	1:A:198:MET:HA	2.50	0.41
1:A:184:ARG:HD3	1:A:192:PHE:CD1	2.56	0.40
1:D:198:MET:HA	1:D:198:MET:CE	2.45	0.40
1:D:73:LEU:HD23	1:D:73:LEU:HA	1.98	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	241/278 (87%)	234 (97%)	6 (2%)	1 (0%)	34	24
1	B	241/278 (87%)	235 (98%)	6 (2%)	0	100	100
1	C	241/278 (87%)	233 (97%)	7 (3%)	1 (0%)	34	24
1	D	241/278 (87%)	232 (96%)	9 (4%)	0	100	100
All	All	964/1112 (87%)	934 (97%)	28 (3%)	2 (0%)	47	38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	136	ALA
1	C	136	ALA

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	211/242 (87%)	205 (97%)	6 (3%)	43	36
1	B	211/242 (87%)	200 (95%)	11 (5%)	23	14
1	C	211/242 (87%)	200 (95%)	11 (5%)	23	14
1	D	211/242 (87%)	204 (97%)	7 (3%)	38	29
All	All	844/968 (87%)	809 (96%)	35 (4%)	30	21

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	127	LYS
1	A	184	ARG
1	A	195	LYS
1	A	201	ILE
1	A	208	ASP
1	A	229	ARG
1	B	21	ILE
1	B	48	ASP
1	B	76	LEU
1	B	91	THR
1	B	184	ARG
1	B	195	LYS
1	B	201	ILE
1	B	203	SER
1	B	208	ASP
1	B	209	THR
1	B	229	ARG
1	C	21	ILE
1	C	77	SER
1	C	91	THR
1	C	127	LYS
1	C	170	LYS
1	C	184	ARG
1	C	199	SER

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Mol	Chain	Res	Type
1	C	201	ILE
1	C	208	ASP
1	C	209	THR
1	C	247	ASN
1	D	46	LYS
1	D	91	THR
1	D	103	LYS
1	D	112	PRO
1	D	184	ARG
1	D	201	ILE
1	D	209	THR

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

Mol	Chain	Res	Type
1	A	18	HIS
1	B	18	HIS
1	B	247	ASN
1	C	40	ASN
1	C	70	GLN
1	C	247	ASN
1	D	18	HIS
1	D	30	HIS
1	D	171	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)

8 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	$\#\|Z\| > 2$
3	SIN	C	301	-	7,7,7	1.71	1 (14%)	8,8,8	3.49	4 (50%)
2	NAD	B	301	-	42,48,48	1.24	4 (9%)	50,73,73	1.35	10 (20%)
3	SIN	D	301	-	7,7,7	1.31	0	8,8,8	2.45	3 (37%)
3	SIN	A	302	-	7,7,7	2.01	1 (14%)	8,8,8	2.77	3 (37%)
2	NAD	A	301	-	42,48,48	1.57	5 (11%)	50,73,73	1.74	10 (20%)
3	SIN	B	302	-	7,7,7	1.66	2 (28%)	8,8,8	2.37	3 (37%)
2	NAD	D	302	-	42,48,48	1.29	5 (11%)	50,73,73	1.58	11 (22%)
2	NAD	C	302	-	42,48,48	1.11	3 (7%)	50,73,73	1.32	7 (14%)

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	SIN	C	301	-	-	4/5/5/5	-
2	NAD	B	301	-	-	6/26/62/62	0/5/5/5
3	SIN	D	301	-	-	4/5/5/5	-
3	SIN	A	302	-	-	3/5/5/5	-
2	NAD	A	301	-	-	5/26/62/62	0/5/5/5
3	SIN	B	302	-	-	2/5/5/5	-
2	NAD	D	302	-	-	5/26/62/62	0/5/5/5
2	NAD	C	302	-	-	6/26/62/62	0/5/5/5

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	NAD	O4D-C1D	4.60	1.47	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	302	SIN	C3-C4	4.42	1.60	1.50
2	B	301	NAD	C4N-C3N	4.19	1.46	1.39
2	A	301	NAD	O4B-C1B	-3.68	1.35	1.41
2	C	302	NAD	C2N-N1N	3.65	1.39	1.35
2	D	302	NAD	O4D-C1D	-3.34	1.36	1.41
2	A	301	NAD	C2N-N1N	3.31	1.39	1.35
3	C	301	SIN	C3-C4	3.16	1.57	1.50
2	D	302	NAD	C4N-C3N	3.04	1.44	1.39
3	B	302	SIN	O3-C4	2.67	1.31	1.22
2	A	301	NAD	C7N-N7N	2.62	1.38	1.33
2	B	301	NAD	C5N-C4N	2.55	1.44	1.38
2	A	301	NAD	C5N-C4N	2.39	1.44	1.38
2	C	302	NAD	C5N-C4N	2.39	1.43	1.38
3	B	302	SIN	C2-C1	2.22	1.55	1.50
2	D	302	NAD	C5N-C4N	2.20	1.43	1.38
2	D	302	NAD	C2D-C3D	-2.19	1.47	1.53
2	C	302	NAD	O2D-C2D	2.17	1.48	1.43
2	B	301	NAD	C5A-C4A	2.11	1.46	1.40
2	D	302	NAD	O3B-C3B	2.11	1.47	1.43
2	B	301	NAD	C2N-N1N	2.08	1.37	1.35

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	301	SIN	O2-C1-C2	6.34	134.39	114.03
3	C	301	SIN	O1-C1-C2	-6.00	103.82	123.08
2	A	301	NAD	O4B-C1B-C2B	-5.77	98.50	106.93
3	B	302	SIN	C3-C2-C1	-5.45	101.86	113.60
3	A	302	SIN	O1-C1-C2	-4.64	108.16	123.08
3	A	302	SIN	C2-C3-C4	-4.58	103.74	113.60
3	D	301	SIN	C3-C2-C1	-4.15	104.67	113.60
3	D	301	SIN	O3-C4-C3	-4.09	109.93	123.08
2	D	302	NAD	N3A-C2A-N1A	-4.02	122.39	128.68
2	A	301	NAD	C3B-C2B-C1B	3.80	106.70	100.98
3	C	301	SIN	C2-C3-C4	-3.77	105.50	113.60
2	A	301	NAD	C3N-C7N-N7N	3.64	122.11	117.75
2	A	301	NAD	N3A-C2A-N1A	-3.50	123.20	128.68
2	C	302	NAD	C1B-N9A-C4A	-3.48	120.53	126.64
2	B	301	NAD	N3A-C2A-N1A	-3.30	123.52	128.68
2	D	302	NAD	O4B-C1B-C2B	-3.28	102.13	106.93
3	A	302	SIN	O2-C1-C2	3.24	124.43	114.03
2	B	301	NAD	O4B-C4B-C3B	3.19	111.42	105.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	302	NAD	O4B-C4B-C3B	3.13	111.31	105.11
3	D	301	SIN	O4-C4-C3	3.12	124.05	114.03
2	D	302	NAD	C6N-N1N-C2N	-3.10	119.15	121.97
2	D	302	NAD	O2N-PN-O1N	2.98	126.96	112.24
2	B	301	NAD	C4A-C5A-N7A	-2.94	106.34	109.40
2	A	301	NAD	O7N-C7N-C3N	-2.90	116.16	119.63
2	C	302	NAD	C3N-C7N-N7N	2.84	121.15	117.75
2	A	301	NAD	C2B-C3B-C4B	-2.75	97.29	102.64
2	D	302	NAD	C5N-C4N-C3N	-2.73	117.11	120.34
2	C	302	NAD	C5N-C4N-C3N	-2.63	117.23	120.34
2	B	301	NAD	C3N-C7N-N7N	2.52	120.78	117.75
2	C	302	NAD	O2A-PA-O1A	2.50	124.60	112.24
2	B	301	NAD	O2B-C2B-C1B	2.47	119.99	110.85
2	D	302	NAD	C3N-C2N-N1N	2.46	122.84	120.43
2	A	301	NAD	C2A-N1A-C6A	2.39	122.84	118.75
2	D	302	NAD	C2B-C3B-C4B	-2.37	98.04	102.64
2	C	302	NAD	C3B-C2B-C1B	2.33	104.49	100.98
2	B	301	NAD	C1B-N9A-C4A	-2.31	122.58	126.64
3	B	302	SIN	O1-C1-C2	-2.31	115.67	123.08
2	D	302	NAD	C3B-C2B-C1B	2.29	104.43	100.98
2	A	301	NAD	O4D-C4D-C3D	2.28	109.63	105.11
2	D	302	NAD	C4A-C5A-N7A	-2.28	107.03	109.40
2	B	301	NAD	O2A-PA-O1A	2.26	123.44	112.24
2	C	302	NAD	O4B-C1B-C2B	-2.17	103.75	106.93
2	A	301	NAD	O2B-C2B-C1B	2.16	118.83	110.85
2	C	302	NAD	N6A-C6A-N1A	2.15	123.05	118.57
2	A	301	NAD	O4B-C4B-C3B	2.11	109.30	105.11
2	B	301	NAD	O7N-C7N-N7N	-2.09	119.61	122.58
2	D	302	NAD	C1B-N9A-C4A	-2.08	122.99	126.64
2	B	301	NAD	O4B-C4B-C5B	-2.08	102.54	109.37
2	B	301	NAD	C3B-C2B-C1B	2.06	104.08	100.98
3	C	301	SIN	O3-C4-C3	-2.02	116.58	123.08
3	B	302	SIN	O2-C1-C2	2.02	120.52	114.03

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	NAD	C5D-O5D-PN-O1N
2	A	301	NAD	C5D-O5D-PN-O2N
2	B	301	NAD	C5D-O5D-PN-O2N
2	B	301	NAD	O4D-C1D-N1N-C2N

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Mol	Chain	Res	Type	Atoms
2	C	302	NAD	C5D-O5D-PN-O2N
2	C	302	NAD	O4D-C1D-N1N-C2N
2	D	302	NAD	C5D-O5D-PN-O1N
2	D	302	NAD	C5D-O5D-PN-O2N
2	A	301	NAD	C5D-O5D-PN-O3
2	B	301	NAD	C5D-O5D-PN-O1N
2	C	302	NAD	C5D-O5D-PN-O1N
3	A	302	SIN	C2-C3-C4-O4
2	A	301	NAD	O4B-C4B-C5B-O5B
3	D	301	SIN	C2-C3-C4-O4
3	D	301	SIN	O2-C1-C2-C3
3	A	302	SIN	C2-C3-C4-O3
3	C	301	SIN	C2-C3-C4-O4
3	D	301	SIN	C2-C3-C4-O3
3	B	302	SIN	O2-C1-C2-C3
3	D	301	SIN	O1-C1-C2-C3
3	C	301	SIN	C2-C3-C4-O3
3	C	301	SIN	O2-C1-C2-C3
3	B	302	SIN	O1-C1-C2-C3
3	A	302	SIN	O2-C1-C2-C3
2	B	301	NAD	C5D-O5D-PN-O3
2	C	302	NAD	C5D-O5D-PN-O3
2	D	302	NAD	C5D-O5D-PN-O3
2	B	301	NAD	O4B-C4B-C5B-O5B
2	C	302	NAD	O4B-C4B-C5B-O5B
2	A	301	NAD	PA-O3-PN-O1N
2	B	301	NAD	PA-O3-PN-O2N
2	C	302	NAD	PA-O3-PN-O1N
2	D	302	NAD	PA-O3-PN-O1N
2	D	302	NAD	O4B-C4B-C5B-O5B
3	C	301	SIN	O1-C1-C2-C3

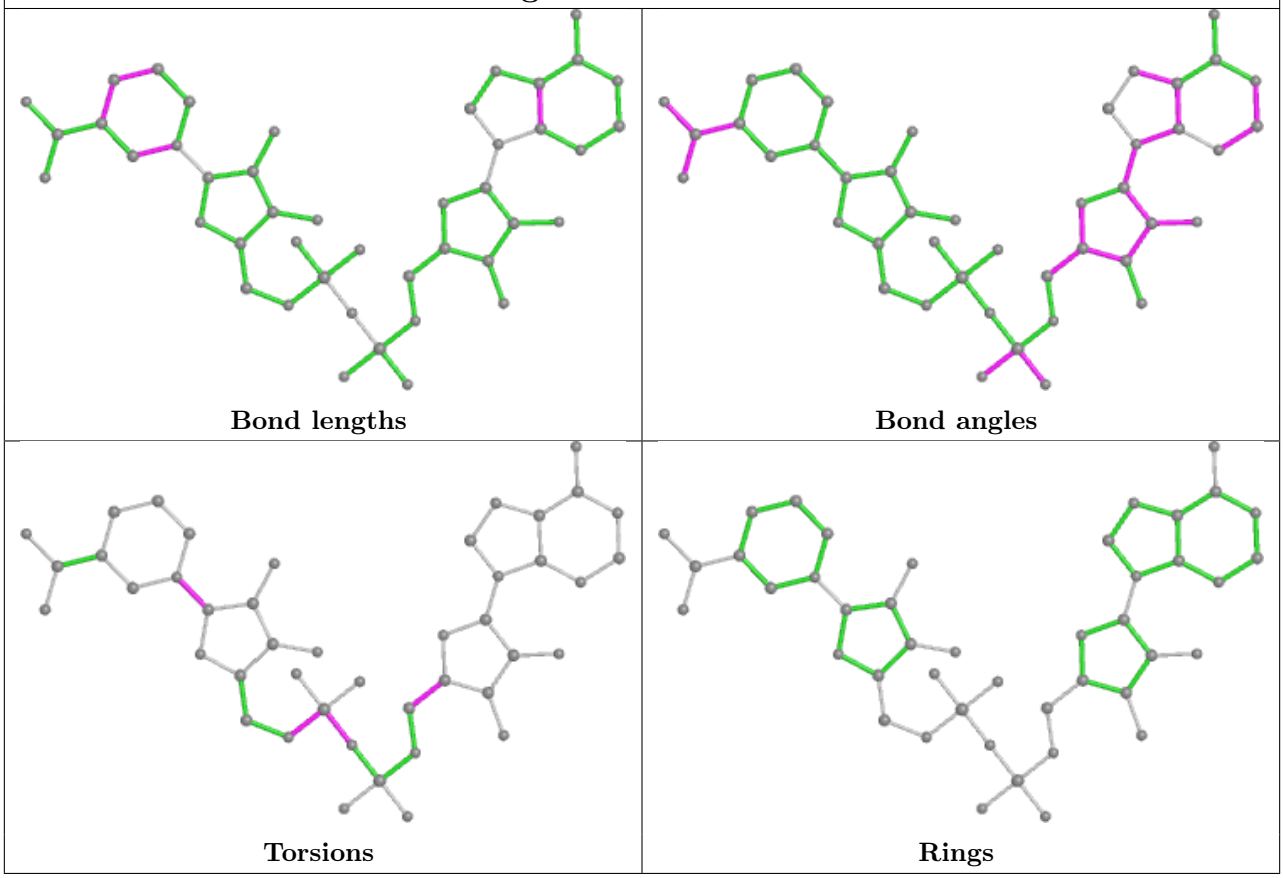
There are no ring outliers.

No monomer is involved in short contacts.

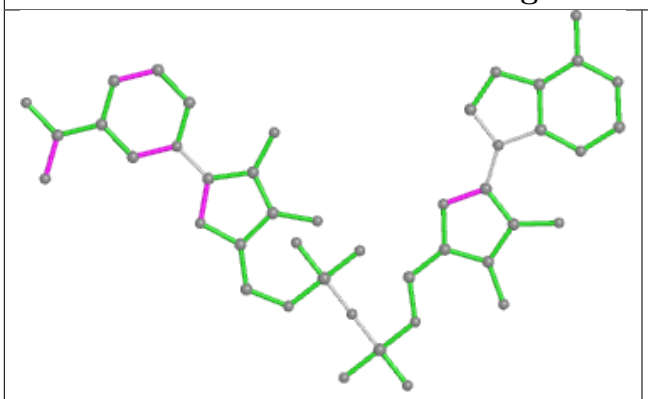
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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.

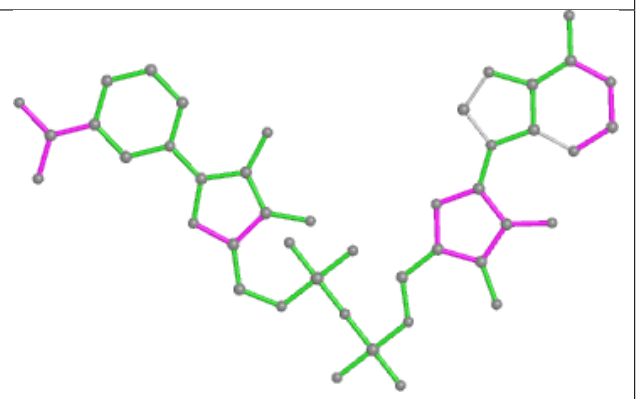
Ligand NAD B 301



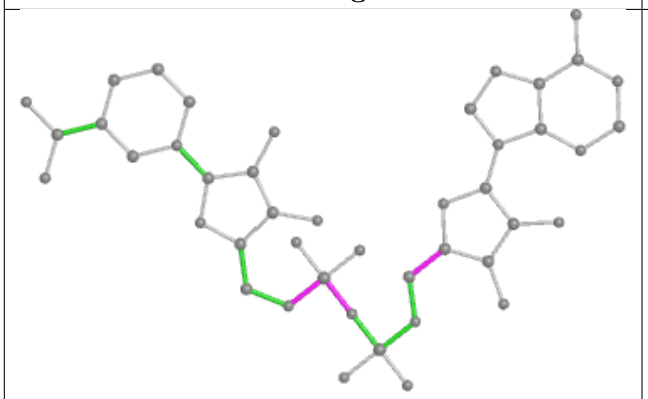
Ligand NAD A 301



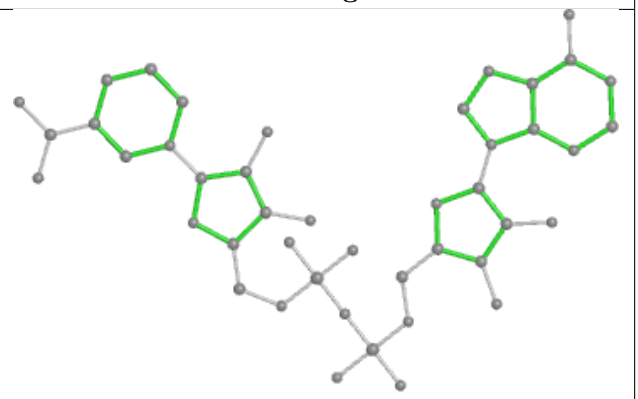
Bond lengths



Bond angles

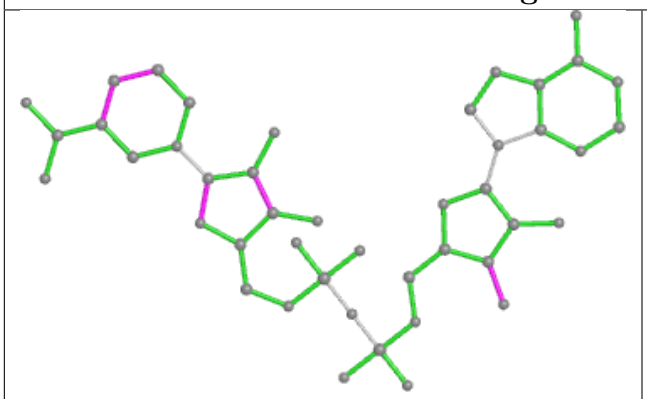


Torsions

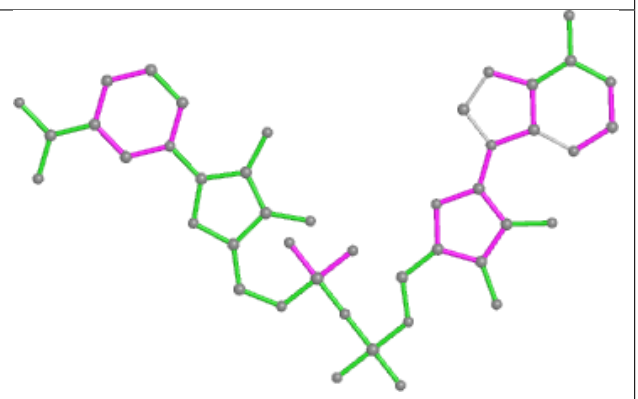


Rings

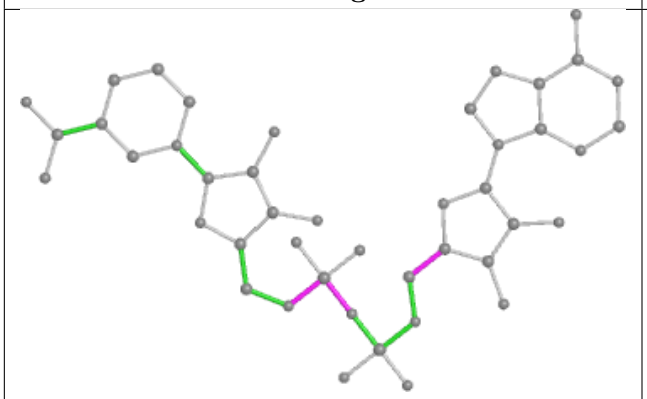
Ligand NAD D 302



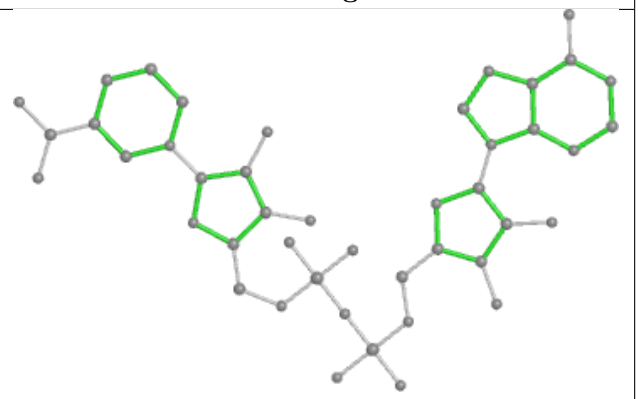
Bond lengths



Bond angles

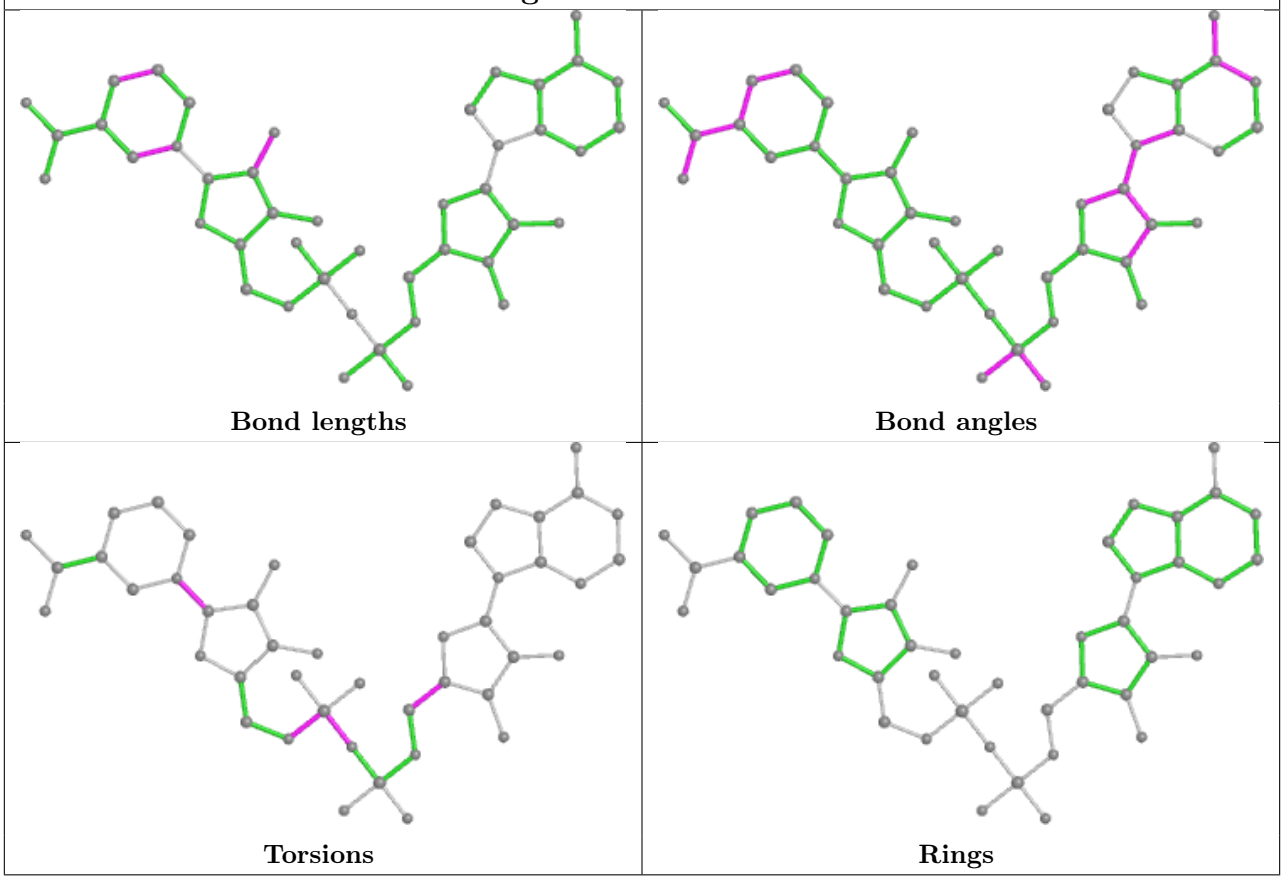


Torsions



Rings

Ligand NAD C 302



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, 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	243/278 (87%)	-0.27	3 (1%)	79	81	10, 17, 36, 47	0
1	B	243/278 (87%)	-0.04	11 (4%)	33	36	11, 20, 47, 78	0
1	C	243/278 (87%)	-0.02	11 (4%)	33	36	13, 22, 49, 81	0
1	D	243/278 (87%)	0.26	19 (7%)	13	14	14, 25, 52, 91	0
All	All	972/1112 (87%)	-0.02	44 (4%)	33	36	10, 21, 48, 91	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	46	LYS	6.2
1	D	42	ILE	5.6
1	D	45	MET	4.9
1	B	46	LYS	4.7
1	C	46	LYS	4.2
1	D	44	TYR	4.1
1	C	48	ASP	4.0
1	D	48	ASP	4.0
1	B	48	ASP	3.8
1	D	73	LEU	3.8
1	D	49	MET	3.7
1	B	196	ALA	3.6
1	D	43	ASP	3.5
1	C	195	LYS	3.4
1	B	45	MET	3.4
1	D	196	ALA	3.2
1	D	193	SER	3.2
1	D	70	GLN	3.1
1	C	196	ALA	3.0
1	D	195	LYS	2.9
1	B	49	MET	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	45	MET	2.9
1	D	47	THR	2.8
1	A	196	ALA	2.6
1	B	43	ASP	2.5
1	D	200	MET	2.5
1	D	75	LYS	2.5
1	B	195	LYS	2.4
1	A	199	SER	2.4
1	C	43	ASP	2.4
1	D	30	HIS	2.4
1	C	49	MET	2.4
1	D	28	SER	2.3
1	B	44	TYR	2.3
1	C	44	TYR	2.3
1	D	51	LEU	2.3
1	D	158	ILE	2.3
1	B	42	ILE	2.2
1	B	47	THR	2.2
1	C	77	SER	2.2
1	B	28	SER	2.1
1	C	158	ILE	2.1
1	C	193	SER	2.0
1	A	141	LEU	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, 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.

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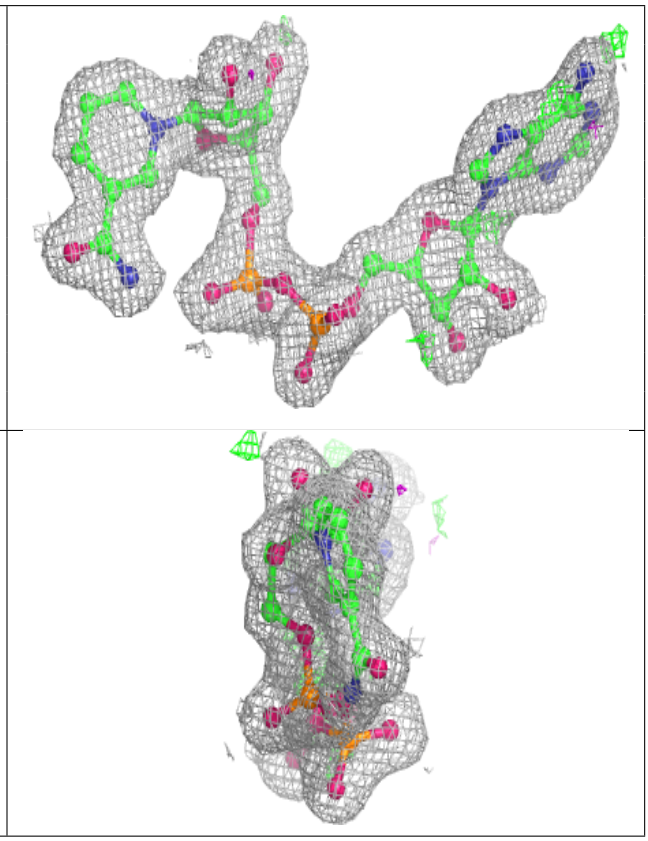
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	SIN	C	301	8/8	0.93	0.11	19,24,28,33	0
3	SIN	A	302	8/8	0.96	0.10	15,17,22,24	0
3	SIN	D	301	8/8	0.96	0.07	20,25,29,30	0
2	NAD	D	302	44/44	0.97	0.07	16,21,26,27	0
3	SIN	B	302	8/8	0.97	0.07	17,22,26,27	0
2	NAD	C	302	44/44	0.98	0.07	15,18,20,22	0
2	NAD	A	301	44/44	0.98	0.07	10,12,14,16	0
2	NAD	B	301	44/44	0.98	0.06	14,17,20,21	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.

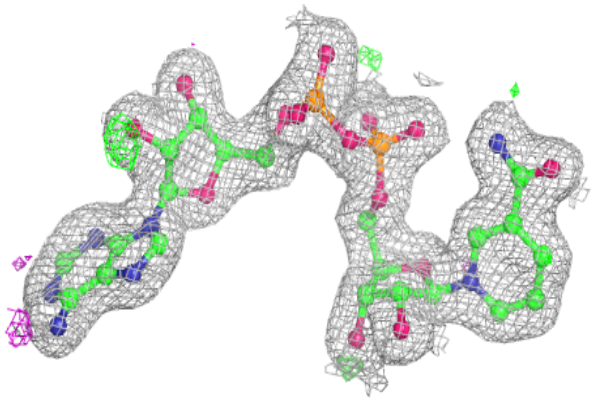
Electron density around NAD D 302:

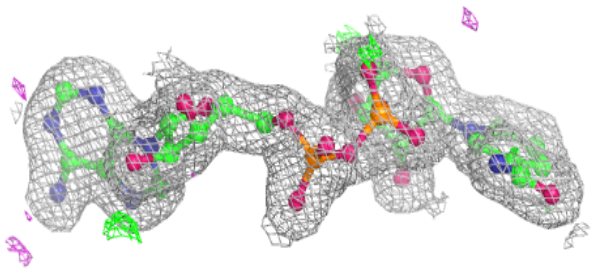
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)

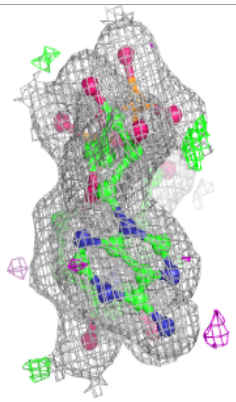


Electron density around NAD C 302:

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

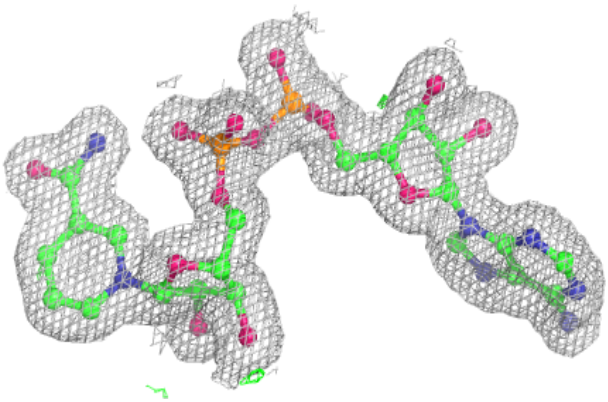


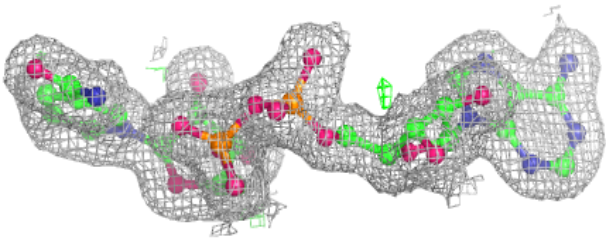


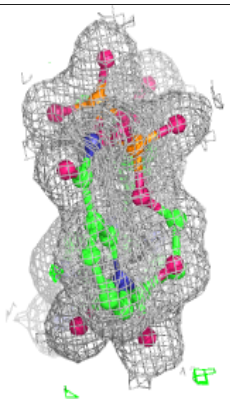


Electron density around NAD A 301:

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

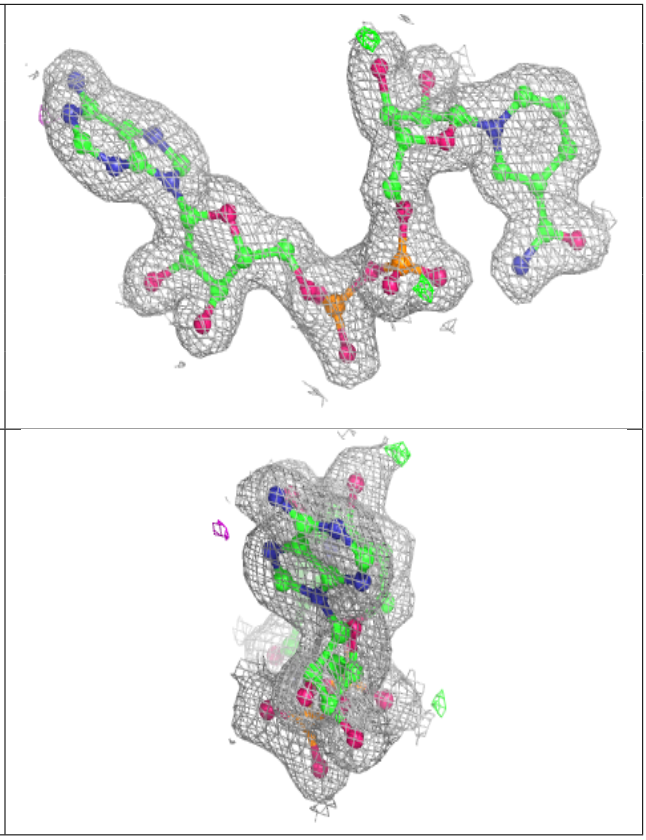


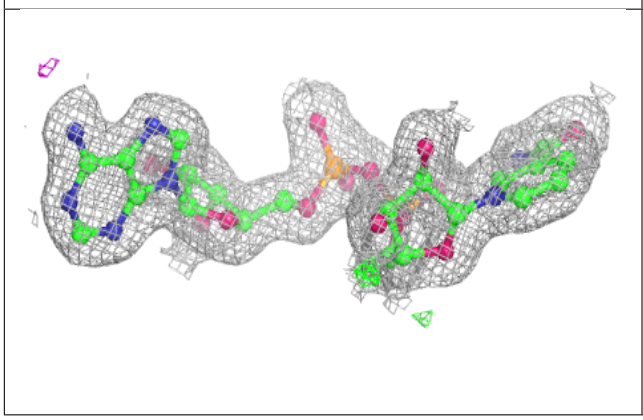




Electron density around NAD B 301:

$2mF_o$-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.