



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 07:58 PM GMT

PDB ID : 4WLO  
Title : Crystal structure of oxaloacetate and NADH bound MDH2  
Authors : Eo, Y.M.; Han, B.G.; Ahn, H.C.  
Deposited on : 2014-10-07  
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026982  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20026982

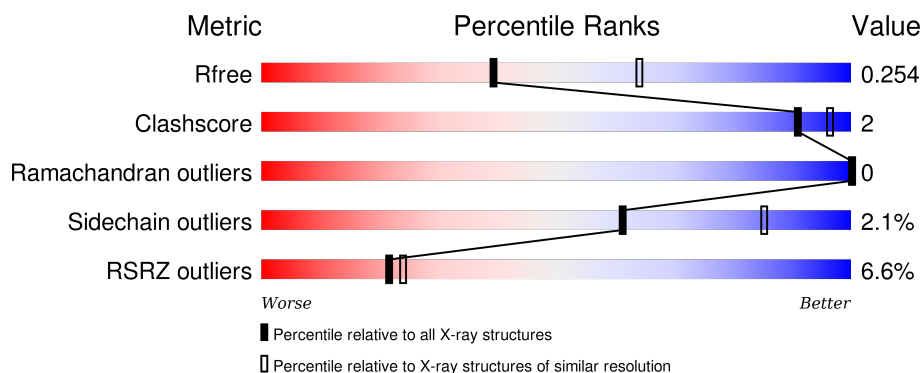
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	340	<div> <div>4%</div> <div>86% 6% • 8%</div> </div>
1	B	340	<div> <div>3%</div> <div>88% • 8%</div> </div>
1	C	340	<div> <div>12%</div> <div>87% • • 8%</div> </div>
1	D	340	<div> <div>5%</div> <div>88% • 8%</div> </div>

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	OAA	A	402	-	-	X	-

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9488 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 Malate dehydrogenase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	0	0	0
			2312	1472	390	437	13			
1	B	314	Total	C	N	O	S	0	0	0
			2312	1472	390	437	13			
1	C	314	Total	C	N	O	S	0	0	0
			2312	1472	390	437	13			
1	D	314	Total	C	N	O	S	0	1	0
			2315	1474	390	438	13			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	expression tag	UNP P40926
A	0	GLY	-	expression tag	UNP P40926
A	1	SER	-	expression tag	UNP P40926
A	2	SER	-	expression tag	UNP P40926
A	3	HIS	-	expression tag	UNP P40926
A	4	HIS	-	expression tag	UNP P40926
A	5	HIS	-	expression tag	UNP P40926
A	6	HIS	-	expression tag	UNP P40926
A	7	HIS	-	expression tag	UNP P40926
A	8	HIS	-	expression tag	UNP P40926
A	9	SER	-	expression tag	UNP P40926
A	10	SER	-	expression tag	UNP P40926
A	11	GLY	-	expression tag	UNP P40926
A	12	LEU	-	expression tag	UNP P40926
A	13	VAL	-	expression tag	UNP P40926
A	14	PRO	-	expression tag	UNP P40926
A	15	ARG	-	expression tag	UNP P40926
A	16	GLY	-	expression tag	UNP P40926
A	17	SER	-	expression tag	UNP P40926
A	18	HIS	-	expression tag	UNP P40926
A	19	MET	-	expression tag	UNP P40926

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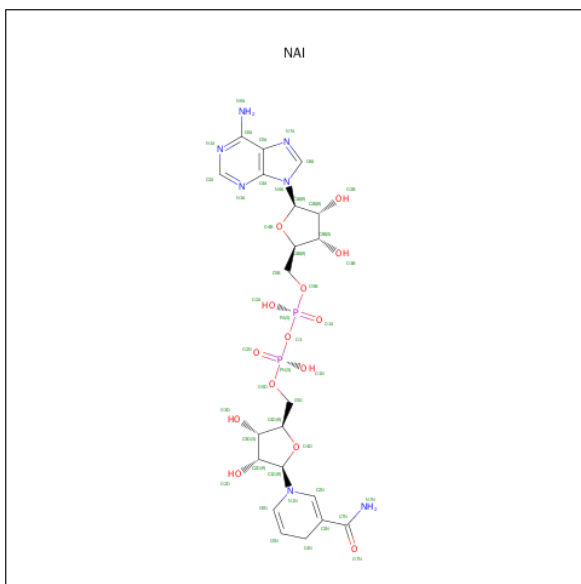
Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	MET	-	expression tag	UNP P40926
B	0	GLY	-	expression tag	UNP P40926
B	1	SER	-	expression tag	UNP P40926
B	2	SER	-	expression tag	UNP P40926
B	3	HIS	-	expression tag	UNP P40926
B	4	HIS	-	expression tag	UNP P40926
B	5	HIS	-	expression tag	UNP P40926
B	6	HIS	-	expression tag	UNP P40926
B	7	HIS	-	expression tag	UNP P40926
B	8	HIS	-	expression tag	UNP P40926
B	9	SER	-	expression tag	UNP P40926
B	10	SER	-	expression tag	UNP P40926
B	11	GLY	-	expression tag	UNP P40926
B	12	LEU	-	expression tag	UNP P40926
B	13	VAL	-	expression tag	UNP P40926
B	14	PRO	-	expression tag	UNP P40926
B	15	ARG	-	expression tag	UNP P40926
B	16	GLY	-	expression tag	UNP P40926
B	17	SER	-	expression tag	UNP P40926
B	18	HIS	-	expression tag	UNP P40926
B	19	MET	-	expression tag	UNP P40926
C	-1	MET	-	expression tag	UNP P40926
C	0	GLY	-	expression tag	UNP P40926
C	1	SER	-	expression tag	UNP P40926
C	2	SER	-	expression tag	UNP P40926
C	3	HIS	-	expression tag	UNP P40926
C	4	HIS	-	expression tag	UNP P40926
C	5	HIS	-	expression tag	UNP P40926
C	6	HIS	-	expression tag	UNP P40926
C	7	HIS	-	expression tag	UNP P40926
C	8	HIS	-	expression tag	UNP P40926
C	9	SER	-	expression tag	UNP P40926
C	10	SER	-	expression tag	UNP P40926
C	11	GLY	-	expression tag	UNP P40926
C	12	LEU	-	expression tag	UNP P40926
C	13	VAL	-	expression tag	UNP P40926
C	14	PRO	-	expression tag	UNP P40926
C	15	ARG	-	expression tag	UNP P40926
C	16	GLY	-	expression tag	UNP P40926
C	17	SER	-	expression tag	UNP P40926
C	18	HIS	-	expression tag	UNP P40926
C	19	MET	-	expression tag	UNP P40926

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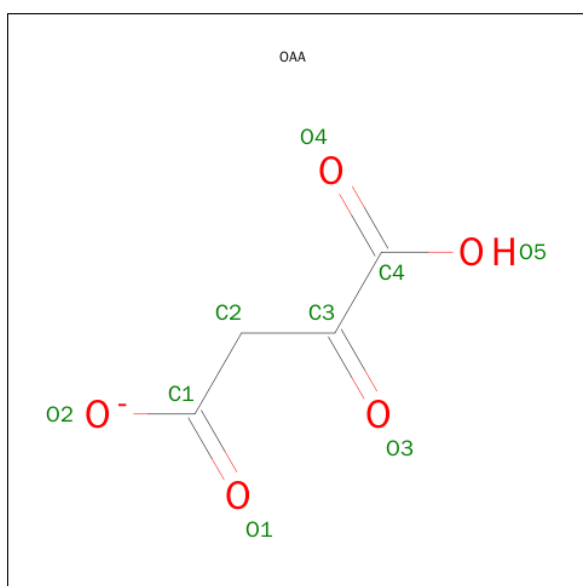
Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	MET	-	expression tag	UNP P40926
D	0	GLY	-	expression tag	UNP P40926
D	1	SER	-	expression tag	UNP P40926
D	2	SER	-	expression tag	UNP P40926
D	3	HIS	-	expression tag	UNP P40926
D	4	HIS	-	expression tag	UNP P40926
D	5	HIS	-	expression tag	UNP P40926
D	6	HIS	-	expression tag	UNP P40926
D	7	HIS	-	expression tag	UNP P40926
D	8	HIS	-	expression tag	UNP P40926
D	9	SER	-	expression tag	UNP P40926
D	10	SER	-	expression tag	UNP P40926
D	11	GLY	-	expression tag	UNP P40926
D	12	LEU	-	expression tag	UNP P40926
D	13	VAL	-	expression tag	UNP P40926
D	14	PRO	-	expression tag	UNP P40926
D	15	ARG	-	expression tag	UNP P40926
D	16	GLY	-	expression tag	UNP P40926
D	17	SER	-	expression tag	UNP P40926
D	18	HIS	-	expression tag	UNP P40926
D	19	MET	-	expression tag	UNP P40926

- Molecule 2 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (three-letter code: NAI) (formula:  $C_{21}H_{29}N_7O_{14}P_2$ ).



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

- Molecule 3 is OXALOACETATE ION (three-letter code: OAA) (formula:  $C_4H_3O_5$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			9	4	5		
3	B	1	Total	C	O	0	0
			9	4	5		
3	C	1	Total	C	O	0	0
			9	4	5		
3	D	1	Total	C	O	0	0
			9	4	5		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	O	0	0
			3	3		
4	B	16	Total	O	0	0
			16	16		

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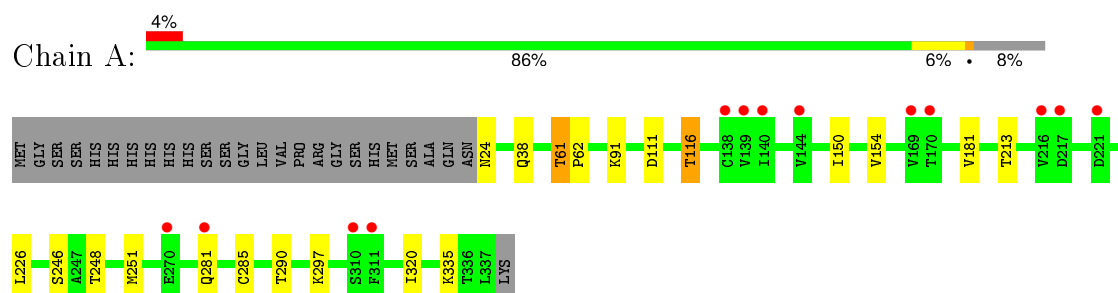
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	2	Total	O	0	0
			2	2		
4	D	4	Total	O	0	0
			4	4		

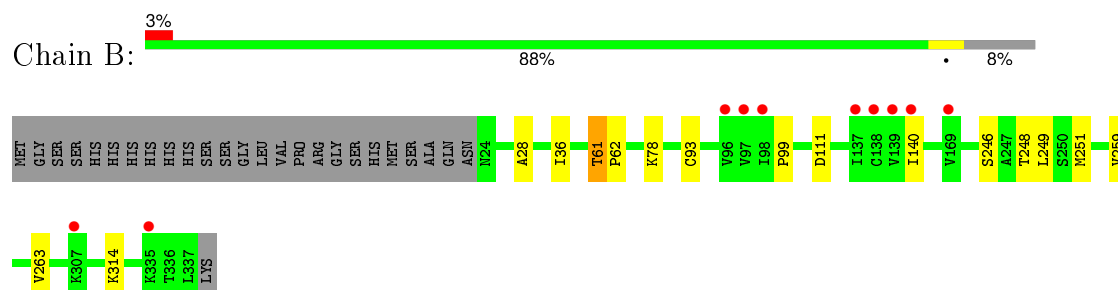
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

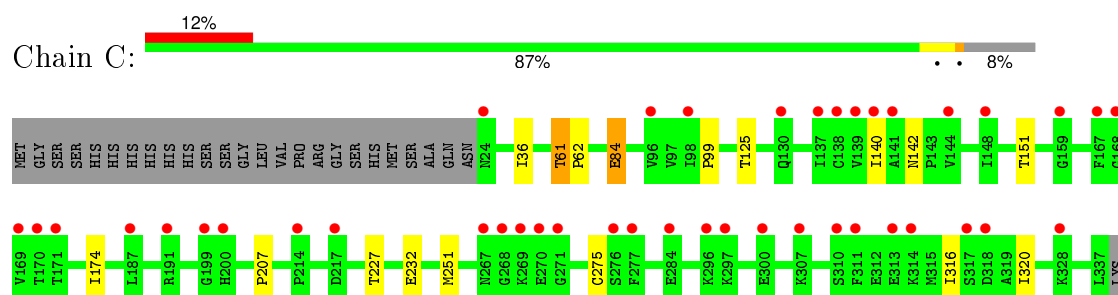
- Molecule 1: Malate dehydrogenase, mitochondrial



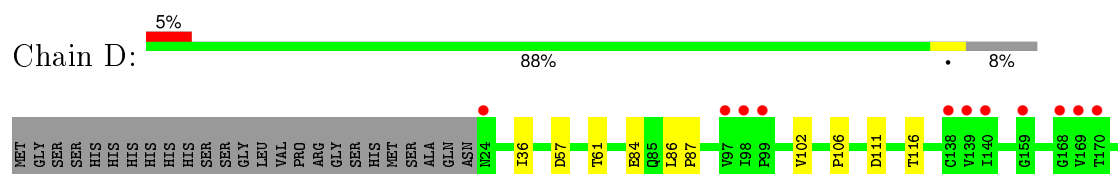
- Molecule 1: Malate dehydrogenase, mitochondrial

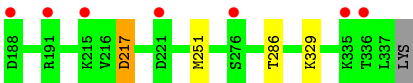


- Molecule 1: Malate dehydrogenase, mitochondrial



- Molecule 1: Malate dehydrogenase, mitochondrial





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	59.99Å 152.19Å 155.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.35 – 2.50 28.35 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.9 (28.35-2.50) 99.0 (28.35-2.50)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	8.30 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, $R_{free}$	0.212 , 0.255 0.216 , 0.254	Depositor DCC
$R_{free}$ test set	2531 reflections (5.36%)	DCC
Wilson B-factor (Å <sup>2</sup> )	43.5	Xtriage
Anisotropy	0.070	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 36.9	EDS
Estimated twinning fraction	0.009 for -h,l,k	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 49982 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9488	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.55% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: NAI, OAA

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	0.33	0/2351	0.51	0/3192
1	B	0.33	0/2351	0.51	0/3192
1	C	0.30	0/2351	0.47	0/3192
1	D	0.31	0/2358	0.48	0/3202
All	All	0.32	0/9411	0.50	0/12778

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2312	0	2400	8	0
1	B	2312	0	2400	9	0
1	C	2312	0	2400	11	0
1	D	2315	0	2403	8	0
2	A	44	0	27	5	0
2	B	44	0	27	5	0
2	C	44	0	27	4	0
2	D	44	0	27	6	0
3	A	9	0	2	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	9	0	2	1	0
3	C	9	0	2	2	0
3	D	9	0	2	3	0
4	A	3	0	0	0	0
4	B	16	0	0	0	0
4	C	2	0	0	0	0
4	D	4	0	0	0	0
All	All	9488	0	9719	43	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 (43) 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:251:MET:SD	2:B:401:NAI:H4N	2.23	0.79
1:A:116:THR:HG23	1:D:106:PRO:O	1.97	0.64
1:A:251:MET:SD	2:A:401:NAI:H4N	2.41	0.60
1:C:316:ILE:O	1:C:320:ILE:HD12	2.05	0.57
2:B:401:NAI:H42N	3:B:402:OAA:C3	2.38	0.54
1:C:84:GLU:H	1:C:84:GLU:CD	2.13	0.52
2:A:401:NAI:H42N	3:A:402:OAA:C3	2.39	0.52
1:D:251:MET:SD	2:D:401:NAI:H4N	2.51	0.51
1:B:99:PRO:O	2:B:401:NAI:H51N	2.11	0.50
1:C:125:THR:HG21	1:C:151:THR:HG23	1.94	0.50
1:D:111:ASP:OD2	1:D:329:LYS:NZ	2.45	0.50
2:A:401:NAI:H42N	3:A:402:OAA:C4	2.42	0.49
1:C:99:PRO:O	2:C:401:NAI:H51N	2.15	0.47
1:C:174:ILE:HD11	1:C:275:CYS:HB2	1.98	0.46
1:D:102:VAL:HG21	1:D:116:THR:HG21	1.97	0.46
1:A:150:ILE:O	1:A:154:VAL:HG23	2.16	0.46
1:B:36:ILE:HD11	2:B:401:NAI:C6N	2.47	0.45
2:A:401:NAI:C5N	3:A:402:OAA:C2	2.94	0.45
2:C:401:NAI:H42N	3:C:402:OAA:C4	2.46	0.45
2:D:401:NAI:H42N	3:D:402:OAA:C3	2.46	0.45
1:B:140:ILE:O	2:B:401:NAI:H2N	2.17	0.45
1:A:285:CYS:SG	1:A:320:ILE:HG23	2.57	0.45
1:D:86:LEU:HB3	1:D:87:PRO:HD3	1.99	0.44
1:A:181:VAL:HG22	1:A:226:LEU:HD21	1.99	0.44
2:D:401:NAI:H42N	3:D:402:OAA:C4	2.47	0.44
1:C:36:ILE:HD11	2:C:401:NAI:C6N	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:57:ASP:OD2	2:D:401:NAI:O3B	2.35	0.44
1:C:142:ASN:ND2	3:C:402:OAA:O3	2.51	0.43
1:B:28:ALA:HB2	1:B:93:CYS:SG	2.59	0.42
1:D:217:ASP:OD1	1:D:217:ASP:N	2.52	0.42
1:A:246:SER:O	1:A:248:THR:HG23	2.20	0.42
2:A:401:NAI:C5N	3:A:402:OAA:H21	2.49	0.42
1:C:99:PRO:HA	1:C:140:ILE:HB	2.02	0.42
1:D:36:ILE:HD11	2:D:401:NAI:C6N	2.50	0.41
1:B:246:SER:O	1:B:248:THR:HG23	2.19	0.41
1:C:251:MET:SD	2:C:401:NAI:H4N	2.60	0.41
1:B:259:VAL:O	1:B:263:VAL:HG23	2.19	0.41
1:B:61:THR:N	1:B:62:PRO:CD	2.84	0.41
1:C:207:PRO:CG	1:C:227:THR:HG23	2.51	0.41
2:D:401:NAI:C5N	3:D:402:OAA:C2	2.98	0.41
1:A:38:GLN:HB3	1:B:249:LEU:HD21	2.04	0.40
1:A:61:THR:N	1:A:62:PRO:CD	2.84	0.40
1:C:61:THR:N	1:C:62:PRO:CD	2.84	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	312/340 (92%)	305 (98%)	7 (2%)	0	100	100
1	B	312/340 (92%)	303 (97%)	9 (3%)	0	100	100
1	C	312/340 (92%)	304 (97%)	8 (3%)	0	100	100
1	D	313/340 (92%)	307 (98%)	6 (2%)	0	100	100
All	All	1249/1360 (92%)	1219 (98%)	30 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains ⓘ

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	253/275 (92%)	243 (96%)	10 (4%)	38	64
1	B	253/275 (92%)	249 (98%)	4 (2%)	70	90
1	C	253/275 (92%)	250 (99%)	3 (1%)	78	93
1	D	254/275 (92%)	249 (98%)	5 (2%)	63	86
All	All	1013/1100 (92%)	991 (98%)	22 (2%)	61	84

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	61	THR
1	A	91	LYS
1	A	111	ASP
1	A	116	THR
1	A	213	THR
1	A	281	GLN
1	A	290	THR
1	A	297	LYS
1	A	335	LYS
1	B	61	THR
1	B	78	LYS
1	B	111	ASP
1	B	314	LYS
1	C	61	THR
1	C	84	GLU
1	C	232	GLU
1	D	61	THR
1	D	84	GLU
1	D	217	ASP
1	D	286[A]	THR
1	D	286[B]	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	164	ASN
1	A	267	ASN
1	B	164	ASN
1	B	193	ASN
1	B	267	ASN
1	C	24	ASN
1	C	267	ASN
1	D	24	ASN
1	D	164	ASN
1	D	267	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

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	$\# Z  > 2$
2	NAI	A	401	-	41,48,48	1.67	7 (17%)	46,73,73	2.32	4 (8%)
3	OAA	A	402	-	2,8,8	1.18	0	2,10,10	2.68	1 (50%)
2	NAI	B	401	-	41,48,48	1.64	7 (17%)	46,73,73	2.29	5 (10%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	OAA	B	402	-	2,8,8	1.24	0	2,10,10	2.60	1 (50%)
2	NAI	C	401	-	41,48,48	1.63	7 (17%)	46,73,73	2.28	4 (8%)
3	OAA	C	402	-	2,8,8	1.66	1 (50%)	2,10,10	1.26	0
2	NAI	D	401	-	41,48,48	1.54	7 (17%)	46,73,73	2.40	5 (10%)
3	OAA	D	402	-	2,8,8	0.75	0	2,10,10	2.75	1 (50%)

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	NAI	A	401	-	-	0/25/72/72	0/5/5/5
3	OAA	A	402	-	-	0/2/8/8	0/0/0/0
2	NAI	B	401	-	-	0/25/72/72	0/5/5/5
3	OAA	B	402	-	-	0/2/8/8	0/0/0/0
2	NAI	C	401	-	-	0/25/72/72	0/5/5/5
3	OAA	C	402	-	-	0/2/8/8	0/0/0/0
2	NAI	D	401	-	-	0/25/72/72	0/5/5/5
3	OAA	D	402	-	-	0/2/8/8	0/0/0/0

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	NAI	C4N-C5N	-4.16	1.40	1.49
2	D	401	NAI	C4N-C5N	-3.93	1.40	1.49
2	C	401	NAI	C4N-C5N	-3.82	1.40	1.49
2	B	401	NAI	C4N-C5N	-3.64	1.41	1.49
2	A	401	NAI	C5A-C4A	-3.06	1.33	1.40
2	D	401	NAI	C5A-C4A	-2.95	1.33	1.40
2	B	401	NAI	C5A-C4A	-2.88	1.34	1.40
2	C	401	NAI	C5A-C4A	-2.75	1.34	1.40
2	D	401	NAI	C2N-C3N	2.09	1.40	1.34
2	C	401	NAI	C2N-C3N	2.12	1.40	1.34
2	D	401	NAI	O4B-C1B	2.19	1.44	1.41
2	A	401	NAI	C2N-C3N	2.21	1.40	1.34
2	B	401	NAI	C2N-C3N	2.21	1.40	1.34
3	C	402	OAA	C2-C3	2.30	1.53	1.51
2	B	401	NAI	O4B-C1B	2.66	1.45	1.41
2	A	401	NAI	C2A-N1A	3.10	1.39	1.33
2	C	401	NAI	O4B-C1B	3.20	1.45	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	NAI	C6N-C5N	3.24	1.39	1.33
2	D	401	NAI	C2A-N1A	3.32	1.40	1.33
2	D	401	NAI	C6N-C5N	3.39	1.39	1.33
2	C	401	NAI	C2A-N1A	3.45	1.40	1.33
2	C	401	NAI	C6N-C5N	3.49	1.39	1.33
2	B	401	NAI	C2A-N1A	3.56	1.40	1.33
2	B	401	NAI	C6N-C5N	3.63	1.39	1.33
2	A	401	NAI	O4B-C1B	3.90	1.46	1.41
2	D	401	NAI	C2A-N3A	3.97	1.39	1.32
2	C	401	NAI	C2A-N3A	4.41	1.40	1.32
2	A	401	NAI	C2A-N3A	4.51	1.40	1.32
2	B	401	NAI	C2A-N3A	4.69	1.40	1.32

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	NAI	N3A-C2A-N1A	-13.47	118.29	128.87
2	D	401	NAI	N3A-C2A-N1A	-13.27	118.45	128.87
2	B	401	NAI	N3A-C2A-N1A	-13.18	118.51	128.87
2	C	401	NAI	N3A-C2A-N1A	-12.93	118.71	128.87
2	D	401	NAI	C1B-N9A-C4A	-5.51	120.66	126.81
2	D	401	NAI	C4B-O4B-C1B	-4.24	105.14	109.64
2	C	401	NAI	C1B-N9A-C4A	-4.08	122.26	126.81
2	A	401	NAI	C1B-N9A-C4A	-4.06	122.27	126.81
3	D	402	OAA	C1-C2-C3	-3.87	108.47	115.52
3	A	402	OAA	C1-C2-C3	-3.73	108.73	115.52
2	B	401	NAI	C1B-N9A-C4A	-3.71	122.66	126.81
3	B	402	OAA	C1-C2-C3	-3.63	108.91	115.52
2	A	401	NAI	C4B-O4B-C1B	-3.43	106.01	109.64
2	B	401	NAI	C4B-O4B-C1B	-3.08	106.38	109.64
2	C	401	NAI	C4B-O4B-C1B	-2.79	106.69	109.64
2	D	401	NAI	C5B-C4B-C3B	-2.16	106.83	115.20
2	B	401	NAI	O4D-C1D-N1N	2.06	112.01	108.09
2	D	401	NAI	O4B-C1B-N9A	2.64	113.09	108.11
2	A	401	NAI	O4B-C1B-N9A	3.58	114.88	108.11
2	C	401	NAI	O4B-C1B-N9A	3.84	115.37	108.11
2	B	401	NAI	O4B-C1B-N9A	4.03	115.72	108.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	NAI	5	0
3	A	402	OAA	4	0
2	B	401	NAI	5	0
3	B	402	OAA	1	0
2	C	401	NAI	4	0
3	C	402	OAA	2	0
2	D	401	NAI	6	0
3	D	402	OAA	3	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	314/340 (92%)	0.31	13 (4%) 41 46	25, 43, 62, 74	0
1	B	314/340 (92%)	0.13	10 (3%) 51 56	21, 33, 48, 69	0
1	C	314/340 (92%)	0.74	42 (13%) 4 4	36, 56, 76, 81	0
1	D	314/340 (92%)	0.38	18 (5%) 27 31	33, 45, 64, 75	0
All	All	1256/1360 (92%)	0.39	83 (6%) 22 24	21, 44, 69, 81	0

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	139	VAL	5.9
1	C	169	VAL	5.5
1	C	310	SER	4.3
1	C	168	GLY	4.3
1	D	139	VAL	4.1
1	B	169	VAL	4.0
1	B	139	VAL	3.7
1	D	98	ILE	3.7
1	C	270	GLU	3.7
1	C	276	SER	3.6
1	C	24	ASN	3.5
1	C	271	GLY	3.5
1	A	270	GLU	3.5
1	D	169	VAL	3.5
1	D	24	ASN	3.4
1	C	140	ILE	3.4
1	C	170	THR	3.4
1	B	335	LYS	3.3
1	C	191	ARG	3.2
1	C	144	VAL	3.2
1	C	98	ILE	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	191	ARG	3.1
1	D	140	ILE	3.1
1	B	140	ILE	3.1
1	B	138	CYS	3.0
1	C	214	PRO	3.0
1	C	171	THR	2.9
1	C	141	ALA	2.9
1	C	284	GLU	2.9
1	A	169	VAL	2.9
1	D	97	VAL	2.8
1	C	317	SER	2.8
1	D	221	ASP	2.8
1	B	98	ILE	2.8
1	C	307	LYS	2.8
1	C	159	GLY	2.8
1	D	335	LYS	2.8
1	C	138	CYS	2.7
1	A	139	VAL	2.7
1	D	215	LYS	2.7
1	D	336	THR	2.7
1	D	138	CYS	2.7
1	A	138	CYS	2.6
1	D	170	THR	2.6
1	C	314	LYS	2.6
1	A	216	VAL	2.5
1	A	221	ASP	2.5
1	C	130	GLN	2.5
1	C	217	ASP	2.5
1	C	268	GLY	2.5
1	C	328	LYS	2.5
1	B	96	VAL	2.4
1	A	140	ILE	2.4
1	C	199	GLY	2.4
1	D	168	GLY	2.4
1	C	200	HIS	2.4
1	C	296	LYS	2.4
1	A	144	VAL	2.4
1	C	148	ILE	2.3
1	C	277	PHE	2.3
1	A	310	SER	2.3
1	B	97	VAL	2.3
1	C	313	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	300	GLU	2.2
1	C	167	PHE	2.2
1	D	99	PRO	2.2
1	A	170	THR	2.2
1	C	311	PHE	2.2
1	C	269	LYS	2.2
1	C	318	ASP	2.2
1	A	217	ASP	2.2
1	B	137	ILE	2.2
1	C	96	VAL	2.1
1	A	311	PHE	2.1
1	C	137	ILE	2.1
1	C	297	LYS	2.1
1	A	281	GLN	2.1
1	D	188	ASP	2.1
1	D	276	SER	2.1
1	B	307	LYS	2.1
1	D	159	GLY	2.0
1	C	267	ASN	2.0
1	C	187	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 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	OAA	C	402	9/9	0.76	0.29	1.79	60,63,64,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	OAA	B	402	9/9	0.90	0.20	1.47	36,38,41,41	0
3	OAA	D	402	9/9	0.90	0.20	1.13	49,53,55,58	0
3	OAA	A	402	9/9	0.91	0.20	0.55	39,46,49,51	0
2	NAI	B	401	44/44	0.89	0.20	0.34	52,56,65,68	0
2	NAI	C	401	44/44	0.89	0.22	0.26	51,65,93,94	0
2	NAI	D	401	44/44	0.93	0.16	-0.43	38,44,58,59	0
2	NAI	A	401	44/44	0.93	0.15	-0.43	36,42,48,49	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.