



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 02:56 pm GMT

PDB ID : 3FNF  
Title : Crystal structure of InhA bound to triclosan derivative  
Authors : Wang, F.  
Deposited on : 2008-12-24  
Resolution : 2.30 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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

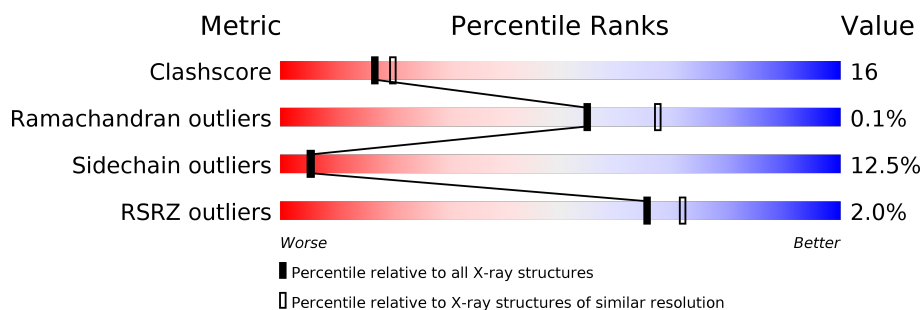
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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))
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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	269	<div> <div>0%</div> <div> <div></div> <div>71%</div> <div>24%</div> <div>5%</div> </div> </div>
1	B	269	<div> <div>0%</div> <div> <div></div> <div>67%</div> <div>19%</div> <div>7%</div> <div>7%</div> </div> </div>
1	C	269	<div> <div>2%</div> <div> <div></div> <div>71%</div> <div>21%</div> <div>• • •</div> </div> </div>
1	D	269	<div> <div>4%</div> <div> <div></div> <div>67%</div> <div>25%</div> <div>• •</div> </div> </div>

## 2 Entry composition [i](#)

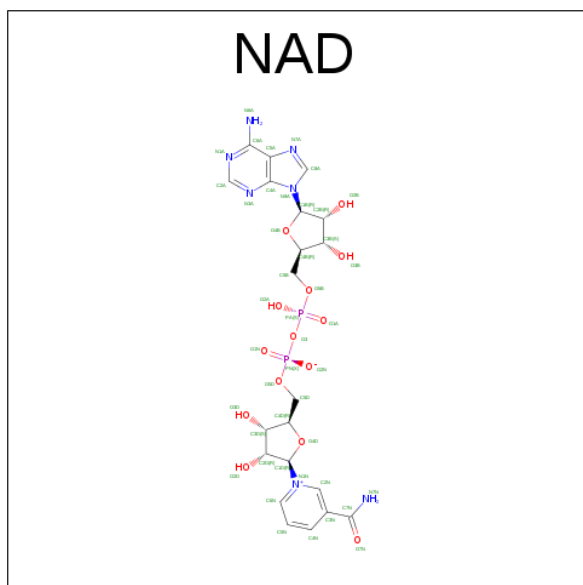
There are 4 unique types of molecules in this entry. The entry contains 8260 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Enoyl-[acyl-carrier-protein] reductase [NADH].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	0	0	0
			1994	1263	348	373	10			
1	B	249	Total	C	N	O	S	0	0	0
			1866	1183	327	346	10			
1	C	261	Total	C	N	O	S	0	0	0
			1953	1238	341	364	10			
1	D	257	Total	C	N	O	S	0	0	0
			1926	1221	337	358	10			

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



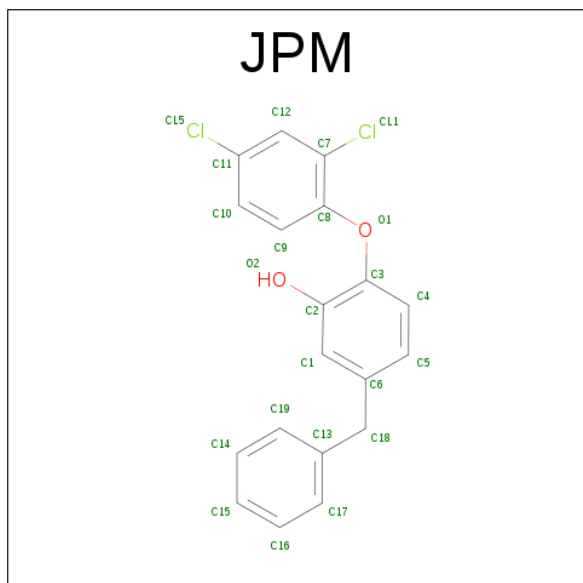
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
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 5-BENZYL-2-(2,4-DICHLOROPHENOXY)PHENOL (three-letter code: JPM) (formula: C<sub>19</sub>H<sub>14</sub>Cl<sub>2</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Cl	O	0	0
			23	19	2	2		
3	B	1	Total	C	Cl	O	0	0
			23	19	2	2		
3	C	1	Total	C	Cl	O	0	0
			23	19	2	2		
3	D	1	Total	C	Cl	O	0	0
			23	19	2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	68	Total	O	0	0
			68	68		
4	B	64	Total	O	0	0
			64	64		
4	C	63	Total	O	0	0
			63	63		

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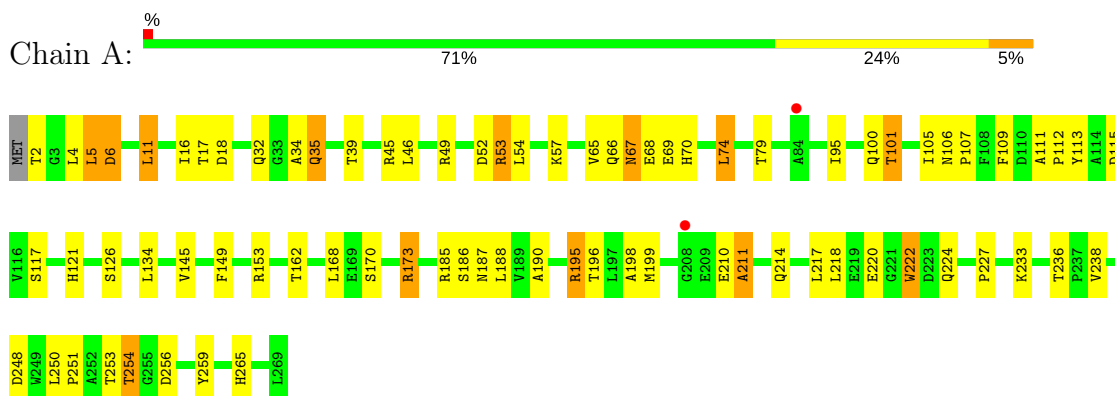
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	58	Total	O	0	0
			58	58		

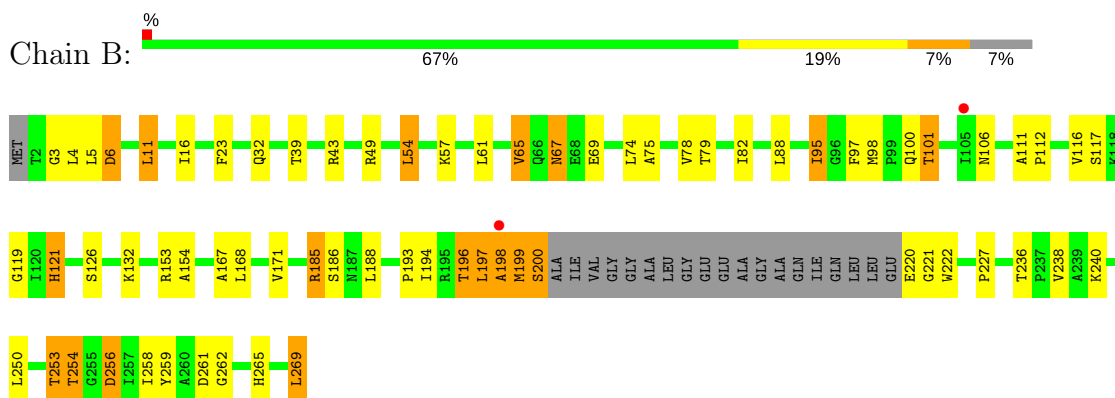
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

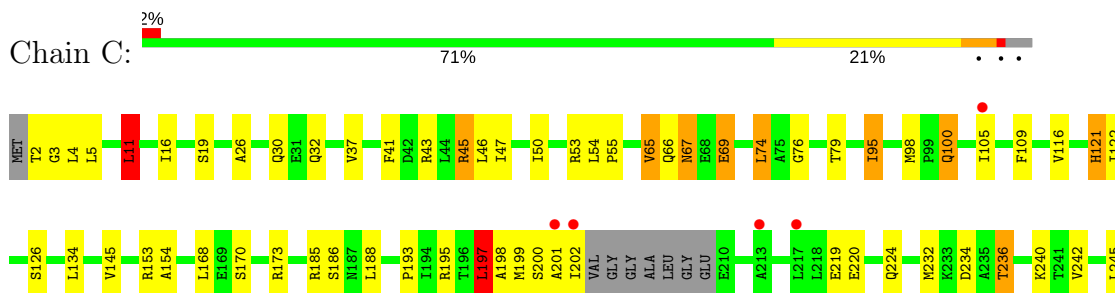
- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]

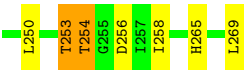


- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]

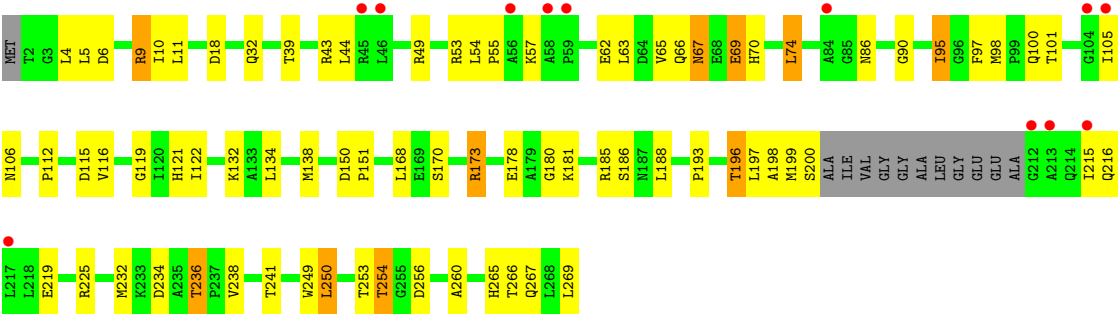


- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]





● Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	125.65Å 92.38Å 102.41Å 90.00° 106.51° 90.00°	Depositor
Resolution (Å)	19.85 – 2.30 40.16 – 1.21	Depositor EDS
% Data completeness (in resolution range)	98.8 (19.85-2.30) 2.1 (40.16-1.21)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.82 (at 1.21Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.200 , 0.260 (Not available) , (Not available)	Depositor DCC
$R_{free}$ test set	NotAvailable	DCC
Wilson B-factor (Å <sup>2</sup> )	5.0	Xtriage
Anisotropy	5.794	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 5.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.41	EDS
Total number of atoms	8260	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.35% 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, 0.333 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: JPM, 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	0.69	0/2032	0.83	3/2758 (0.1%)
1	B	0.67	0/1903	0.80	4/2582 (0.2%)
1	C	0.65	0/1990	0.82	4/2700 (0.1%)
1	D	0.61	0/1963	0.78	2/2663 (0.1%)
All	All	0.65	0/7888	0.81	13/10703 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	173	ARG	NE-CZ-NH2	-9.48	115.56	120.30
1	C	11	LEU	CA-CB-CG	7.45	132.43	115.30
1	C	269	LEU	CA-CB-CG	7.24	131.96	115.30
1	C	173	ARG	NE-CZ-NH2	-7.02	116.79	120.30
1	B	11	LEU	CA-CB-CG	6.84	131.04	115.30
1	A	173	ARG	NE-CZ-NH1	5.45	123.03	120.30
1	A	11	LEU	CA-CB-CG	5.42	127.77	115.30
1	B	185	ARG	NE-CZ-NH2	-5.36	117.62	120.30
1	B	269	LEU	CA-CB-CG	5.32	127.54	115.30
1	B	198	ALA	N-CA-C	5.18	124.97	111.00
1	D	173	ARG	NE-CZ-NH2	-5.11	117.75	120.30
1	C	197	LEU	CA-CB-CG	5.05	126.92	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	18	ASP	CB-CG-OD1	5.00	122.81	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	211	ALA	Peptide

## 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	1994	0	2008	74	0
1	B	1866	0	1877	63	0
1	C	1953	0	1967	65	0
1	D	1926	0	1940	68	0
2	A	44	0	26	2	0
2	B	44	0	26	2	0
2	C	44	0	26	1	0
2	D	44	0	26	3	0
3	A	23	0	13	0	0
3	B	23	0	13	2	0
3	C	23	0	13	1	0
3	D	23	0	13	3	0
4	A	68	0	0	6	0
4	B	64	0	0	5	0
4	C	63	0	0	9	0
4	D	58	0	0	3	0
All	All	8260	0	7948	250	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 16.

All (250) 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:16:ILE:HG23	1:B:197:LEU:CD2	1.73	1.18
1:B:197:LEU:HD12	1:B:197:LEU:H	1.15	1.09
1:B:106:ASN:HB2	4:B:304:HOH:O	1.50	1.08
1:A:211:ALA:HA	1:A:214:GLN:HB2	1.36	1.05
1:C:186:SER:H	1:C:254:THR:HG22	1.25	0.97
1:B:197:LEU:H	1:B:197:LEU:CD1	1.75	0.96
1:B:16:ILE:CG2	1:B:197:LEU:CD2	2.47	0.93
1:D:186:SER:H	1:D:254:THR:HG22	1.32	0.92
1:C:100:GLN:HG3	4:C:327:HOH:O	1.69	0.91
1:D:106:ASN:HB2	4:D:279:HOH:O	1.69	0.91
1:C:4:LEU:H	1:C:32:GLN:HE21	1.17	0.90
1:A:254:THR:HG21	4:A:273:HOH:O	1.72	0.88
1:B:16:ILE:HG23	1:B:197:LEU:HD21	1.57	0.86
1:B:256:ASP:OD2	1:C:265:HIS:HE1	1.58	0.86
1:B:196:THR:HG23	1:B:198:ALA:H	1.44	0.83
1:B:221:GLY:HA2	4:B:284:HOH:O	1.77	0.83
1:B:101:THR:HG21	1:B:112:PRO:HD2	1.60	0.82
1:D:95:ILE:HD11	1:D:122:ILE:HG23	1.60	0.82
1:C:220:GLU:O	1:C:224:GLN:HG3	1.80	0.81
1:A:195:ARG:HH11	1:A:195:ARG:HG3	1.47	0.80
1:D:95:ILE:CD1	1:D:122:ILE:HG23	2.11	0.79
1:B:197:LEU:HD12	1:B:197:LEU:N	1.96	0.79
1:C:65:VAL:HG11	1:C:126:SER:HB3	1.65	0.79
1:A:186:SER:H	1:A:254:THR:CG2	1.96	0.79
1:C:65:VAL:CG1	1:C:126:SER:HB3	2.13	0.79
1:B:16:ILE:HG23	1:B:197:LEU:HD23	1.64	0.77
1:D:74:LEU:HD13	1:D:134:LEU:HD21	1.67	0.77
1:D:196:THR:HG22	1:D:199:MET:H	1.49	0.77
1:C:245:LEU:HD11	1:C:258:ILE:HD13	1.68	0.76
1:A:153:ARG:HD2	4:A:331:HOH:O	1.84	0.76
1:C:4:LEU:H	1:C:32:GLN:NE2	1.84	0.75
1:A:4:LEU:H	1:A:32:GLN:HE21	1.33	0.75
1:B:4:LEU:H	1:B:32:GLN:HE21	1.34	0.75
1:D:74:LEU:CD1	1:D:134:LEU:HD21	2.18	0.74
1:A:105:ILE:HD11	1:A:210:GLU:O	1.87	0.74
1:A:153:ARG:NH2	1:C:153:ARG:NH2	2.37	0.72
1:B:254:THR:HG21	4:B:272:HOH:O	1.88	0.72
1:A:186:SER:H	1:A:254:THR:HG22	1.55	0.72
1:B:253:THR:HG23	4:C:272:HOH:O	1.88	0.72
1:A:256:ASP:OD2	1:D:265:HIS:HE1	1.73	0.72
1:A:196:THR:HG23	1:A:198:ALA:H	1.55	0.71
1:C:45:ARG:HE	1:C:45:ARG:H	1.36	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:ALA:O	1:B:79:THR:HG23	1.90	0.71
1:B:16:ILE:CG2	1:B:197:LEU:HD22	2.20	0.71
1:A:68:GLU:HB2	4:A:314:HOH:O	1.92	0.70
1:C:65:VAL:HG11	1:C:126:SER:CB	2.21	0.70
1:D:44:LEU:HD21	1:D:62:GLU:HB2	1.73	0.70
1:A:259:TYR:O	1:D:253:THR:HG22	1.92	0.70
1:B:256:ASP:OD2	1:C:265:HIS:CE1	2.45	0.70
1:D:185:ARG:HA	1:D:254:THR:CG2	2.22	0.69
1:D:196:THR:HG23	1:D:198:ALA:H	1.56	0.69
1:B:101:THR:HG23	1:B:111:ALA:HA	1.72	0.69
1:C:195:ARG:HE	1:C:232:MET:HE2	1.57	0.68
1:A:101:THR:HG23	1:A:111:ALA:HA	1.76	0.68
1:A:2:THR:O	1:A:6:ASP:OD2	2.11	0.68
1:B:196:THR:HG21	2:B:310:NAD:O1N	1.93	0.67
4:A:271:HOH:O	1:D:253:THR:HG23	1.94	0.67
1:B:153:ARG:HD2	1:D:265:HIS:O	1.93	0.67
1:B:253:THR:CG2	4:C:272:HOH:O	2.44	0.66
1:C:234:ASP:OD1	1:C:236:THR:HG23	1.96	0.65
1:A:35:GLN:HB3	4:A:294:HOH:O	1.96	0.65
1:A:196:THR:HG21	2:A:300:NAD:O1N	1.97	0.65
1:C:186:SER:N	1:C:254:THR:HG22	2.08	0.64
1:C:185:ARG:HD3	4:C:289:HOH:O	1.98	0.64
1:C:254:THR:HG21	4:C:299:HOH:O	1.98	0.64
1:A:153:ARG:CD	4:A:331:HOH:O	2.46	0.63
1:D:215:ILE:HG22	1:D:219:GLU:OE1	1.99	0.63
1:A:117:SER:HG	1:B:117:SER:HG	1.45	0.63
1:C:199:MET:O	1:C:201:ALA:HA	1.98	0.63
1:B:186:SER:H	1:B:254:THR:HG23	1.62	0.63
1:A:67:ASN:HD22	1:A:67:ASN:C	2.03	0.62
1:D:241:THR:HG21	1:D:260:ALA:HB2	1.82	0.62
1:B:153:ARG:HD2	1:D:266:THR:HA	1.80	0.62
1:A:227:PRO:O	1:D:180:GLY:HA3	2.00	0.62
1:B:265:HIS:HE1	1:C:256:ASP:OD2	1.83	0.61
1:D:196:THR:HG21	2:D:330:NAD:O1N	2.01	0.61
1:D:67:ASN:HD21	1:D:69:GLU:HB3	1.65	0.61
1:A:101:THR:HG21	1:A:115:ASP:OD2	2.01	0.60
1:A:185:ARG:HA	1:A:254:THR:HG23	1.84	0.60
1:B:200:SER:O	1:B:200:SER:OG	2.16	0.60
1:C:185:ARG:HA	1:C:254:THR:CG2	2.31	0.60
1:C:79:THR:HG23	4:C:296:HOH:O	2.02	0.59
1:D:4:LEU:H	1:D:32:GLN:HE21	1.49	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:43:ARG:O	1:C:47:ILE:HG13	2.02	0.59
1:C:202:ILE:HG22	1:C:202:ILE:O	2.03	0.58
1:B:220:GLU:HA	4:B:298:HOH:O	2.03	0.58
1:B:67:ASN:HD21	1:B:69:GLU:HG2	1.68	0.58
1:C:65:VAL:CG1	1:C:126:SER:CB	2.81	0.57
1:B:185:ARG:NH2	4:B:334:HOH:O	2.37	0.57
1:D:4:LEU:HB3	1:D:32:GLN:HG3	1.86	0.57
1:A:49:ARG:O	1:A:53:ARG:NE	2.38	0.57
1:A:173:ARG:HB3	1:B:154:ALA:HB2	1.85	0.57
1:B:222:TRP:HE1	1:B:261:ASP:HB2	1.69	0.57
1:C:16:ILE:HD11	1:C:43:ARG:HD2	1.86	0.56
1:C:195:ARG:HE	1:C:232:MET:CE	2.19	0.56
1:A:100:GLN:H	1:A:100:GLN:CD	2.09	0.56
1:D:4:LEU:H	1:D:32:GLN:HG3	1.71	0.56
1:D:70:HIS:O	1:D:74:LEU:HB2	2.05	0.56
1:B:258:ILE:HD12	1:B:258:ILE:N	2.22	0.55
1:C:95:ILE:CD1	1:C:122:ILE:HG23	2.36	0.55
1:A:186:SER:H	1:A:254:THR:HG23	1.71	0.55
1:D:98:MET:HE3	1:D:115:ASP:O	2.06	0.55
1:D:101:THR:HG21	1:D:115:ASP:OD2	2.07	0.55
1:B:153:ARG:CD	1:D:266:THR:HA	2.37	0.55
1:D:134:LEU:O	1:D:138:MET:HG3	2.06	0.55
1:A:265:HIS:HE1	1:D:256:ASP:OD2	1.90	0.54
1:D:234:ASP:OD1	1:D:236:THR:HG23	2.07	0.54
1:A:16:ILE:HG23	1:A:17:THR:HG23	1.89	0.54
1:C:65:VAL:HG13	1:C:126:SER:HB3	1.87	0.54
1:A:256:ASP:OD2	1:D:265:HIS:CE1	2.58	0.54
1:B:186:SER:H	1:B:254:THR:CG2	2.21	0.54
1:D:215:ILE:O	1:D:219:GLU:HG2	2.08	0.53
1:B:65:VAL:HG11	1:B:95:ILE:HD11	1.89	0.53
1:D:198:ALA:HB1	3:D:430:JPM:C7	2.39	0.53
1:C:76:GLY:HA3	4:C:326:HOH:O	2.09	0.53
1:A:46:LEU:HD12	1:A:49:ARG:NH2	2.23	0.53
1:A:45:ARG:HE	1:A:49:ARG:HH21	1.57	0.53
1:A:67:ASN:HD21	1:A:69:GLU:HB2	1.73	0.53
1:C:67:ASN:C	1:C:67:ASN:HD22	2.12	0.53
1:C:95:ILE:HD11	1:C:122:ILE:HG23	1.91	0.52
1:D:9:ARG:HH21	1:D:86:ASN:CG	2.13	0.52
1:B:194:ILE:O	1:B:200:SER:HB3	2.09	0.52
1:D:199:MET:O	1:D:200:SER:CB	2.57	0.52
1:A:101:THR:HG21	1:A:112:PRO:HD2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:GLN:H	1:A:100:GLN:NE2	2.08	0.51
1:D:4:LEU:CB	1:D:32:GLN:HG3	2.40	0.51
1:B:101:THR:CG2	1:B:111:ALA:HA	2.39	0.51
1:C:30:GLN:OE1	1:C:55:PRO:HD2	2.09	0.51
1:D:254:THR:HG21	4:D:276:HOH:O	2.09	0.51
1:C:11:LEU:HA	1:C:37:VAL:O	2.11	0.51
1:B:196:THR:HG22	1:B:199:MET:O	2.11	0.51
1:B:4:LEU:H	1:B:32:GLN:NE2	2.05	0.51
1:A:45:ARG:NH2	1:A:49:ARG:HH22	2.09	0.51
1:B:23:PHE:CE2	1:B:54:LEU:HD13	2.46	0.51
1:B:227:PRO:HD2	1:B:262:GLY:O	2.10	0.51
1:A:195:ARG:CG	1:A:195:ARG:HH11	2.21	0.50
1:B:78:VAL:O	1:B:82:ILE:HG12	2.11	0.50
1:D:178:GLU:HA	1:D:181:LYS:HE2	1.93	0.50
1:D:198:ALA:HB1	3:D:430:JPM:C8	2.41	0.50
1:D:95:ILE:CD1	1:D:122:ILE:CG2	2.87	0.50
1:A:45:ARG:NE	1:A:49:ARG:HH21	2.10	0.49
1:D:67:ASN:HD22	1:D:70:HIS:H	1.60	0.49
1:A:211:ALA:CA	1:A:214:GLN:HB2	2.25	0.49
1:C:197:LEU:O	1:C:200:SER:O	2.29	0.49
1:A:65:VAL:HG13	1:A:126:SER:CB	2.42	0.49
1:D:101:THR:HG21	1:D:112:PRO:HD2	1.95	0.49
1:A:46:LEU:HD12	1:A:49:ARG:HH22	1.77	0.49
1:A:66:GLN:HE21	1:A:121:HIS:CD2	2.31	0.49
1:B:236:THR:HG22	1:B:240:LYS:HE3	1.94	0.48
1:D:10:ILE:HG12	1:D:90:GLY:HA3	1.93	0.48
1:D:97:PHE:O	1:D:119:GLY:HA2	2.13	0.48
1:A:65:VAL:HG13	1:A:126:SER:HB3	1.95	0.48
1:B:194:ILE:O	1:B:200:SER:CB	2.61	0.48
1:C:43:ARG:NE	4:C:319:HOH:O	2.47	0.48
1:A:113:TYR:CE2	1:B:121:HIS:HB2	2.48	0.48
1:C:154:ALA:HB2	1:D:173:ARG:HB3	1.94	0.48
1:D:225:ARG:HD2	1:D:267:GLN:O	2.14	0.48
1:C:26:ALA:O	1:C:30:GLN:HG3	2.14	0.48
1:B:98:MET:CE	1:B:116:VAL:HA	2.44	0.47
1:A:153:ARG:HH21	1:C:153:ARG:NH2	2.12	0.47
1:D:66:GLN:HE21	1:D:121:HIS:CD2	2.32	0.47
1:D:4:LEU:N	1:D:32:GLN:HG3	2.29	0.47
1:A:65:VAL:CG1	1:A:126:SER:CB	2.93	0.47
1:A:185:ARG:HA	1:A:254:THR:CG2	2.44	0.47
1:C:193:PRO:HA	2:C:320:NAD:O7N	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:186:SER:N	1:D:254:THR:HG22	2.15	0.47
1:D:249:TRP:O	1:D:250:LEU:HD13	2.15	0.47
1:A:185:ARG:CA	1:A:254:THR:HG23	2.44	0.46
1:D:49:ARG:NE	4:D:293:HOH:O	2.40	0.46
1:D:185:ARG:HA	1:D:254:THR:HG23	1.96	0.46
1:C:67:ASN:HD22	1:C:69:GLU:H	1.63	0.46
1:D:196:THR:HG23	1:D:198:ALA:N	2.29	0.46
1:B:98:MET:HE3	1:B:116:VAL:HA	1.96	0.46
1:C:185:ARG:CG	1:C:254:THR:HG23	2.46	0.46
1:A:52:ASP:HB2	1:A:53:ARG:HH21	1.80	0.45
1:A:11:LEU:HD21	1:A:39:THR:CG2	2.46	0.45
1:D:4:LEU:HB3	1:D:32:GLN:CG	2.45	0.45
1:A:220:GLU:O	1:A:224:GLN:HG3	2.17	0.45
1:A:65:VAL:HB	2:A:300:NAD:N1A	2.31	0.45
1:C:236:THR:O	1:C:240:LYS:HG3	2.16	0.45
1:A:65:VAL:CG1	1:A:126:SER:HB3	2.47	0.45
1:A:211:ALA:HA	1:A:214:GLN:CB	2.26	0.45
1:C:46:LEU:O	1:C:50:ILE:HG12	2.16	0.45
1:C:98:MET:CE	1:C:116:VAL:HA	2.46	0.45
1:A:105:ILE:HD12	1:A:105:ILE:HA	1.76	0.45
1:C:234:ASP:OD1	1:C:236:THR:CG2	2.63	0.45
1:A:52:ASP:HB2	1:A:53:ARG:HE	1.82	0.45
1:C:198:ALA:HB1	3:C:420:JPM:C7	2.47	0.45
1:B:259:TYR:O	1:C:253:THR:HB	2.17	0.44
1:A:187:ASN:ND2	1:A:256:ASP:H	2.15	0.44
1:A:49:ARG:O	1:A:53:ARG:CZ	2.65	0.44
1:B:65:VAL:CG2	1:B:126:SER:HB2	2.48	0.44
1:A:210:GLU:O	1:A:211:ALA:HB3	2.17	0.44
1:A:74:LEU:HD13	1:A:134:LEU:HD21	2.00	0.44
1:B:23:PHE:CZ	1:B:54:LEU:HD13	2.52	0.44
1:A:190:ALA:HB3	1:A:259:TYR:CD2	2.53	0.44
1:A:5:LEU:HB3	1:A:34:ALA:HB2	1.98	0.44
1:C:67:ASN:ND2	1:C:69:GLU:H	2.16	0.44
1:D:185:ARG:HA	1:D:254:THR:HG21	1.99	0.44
1:B:198:ALA:HB1	3:B:410:JPM:C7	2.48	0.44
1:C:43:ARG:NH1	4:C:309:HOH:O	2.51	0.44
1:D:199:MET:HB2	3:D:430:JPM:H5	1.98	0.44
1:A:106:ASN:HA	1:A:107:PRO:HD3	1.89	0.43
1:B:3:GLY:O	1:B:6:ASP:HB2	2.17	0.43
1:D:150:ASP:HA	1:D:151:PRO:HD3	1.90	0.43
1:C:74:LEU:HD13	1:C:134:LEU:HD21	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:67:ASN:HD22	1:C:69:GLU:N	2.16	0.43
1:B:97:PHE:O	1:B:119:GLY:HA2	2.19	0.43
1:C:109:PHE:HB3	1:D:132:LYS:HD2	2.00	0.43
1:A:196:THR:HG23	1:A:198:ALA:N	2.29	0.42
1:C:232:MET:HE3	1:C:232:MET:HB3	1.80	0.42
1:D:63:LEU:HD13	1:D:74:LEU:HG	2.01	0.42
1:A:195:ARG:NH1	1:A:195:ARG:HG3	2.25	0.42
1:A:248:ASP:O	1:A:251:PRO:HG3	2.20	0.42
1:C:219:GLU:HG2	1:C:232:MET:SD	2.60	0.42
1:C:66:GLN:HE21	1:C:121:HIS:CD2	2.37	0.42
1:B:167:ALA:O	1:B:171:VAL:HG23	2.19	0.42
1:D:11:LEU:HD13	1:D:11:LEU:C	2.40	0.42
1:B:153:ARG:HA	1:B:153:ARG:HD3	1.94	0.42
1:A:145:VAL:HA	1:A:187:ASN:O	2.20	0.41
1:D:186:SER:H	1:D:254:THR:CG2	2.17	0.41
1:A:45:ARG:NH2	1:A:49:ARG:NH2	2.68	0.41
1:C:105:ILE:HG13	1:C:105:ILE:O	2.20	0.41
1:D:98:MET:CE	1:D:116:VAL:HA	2.50	0.41
1:A:67:ASN:ND2	1:A:70:HIS:H	2.19	0.41
1:C:145:VAL:HG11	1:C:242:VAL:HG13	2.02	0.41
1:C:41:PHE:CD1	1:C:41:PHE:C	2.93	0.41
1:C:201:ALA:HB1	1:C:202:ILE:HD13	2.03	0.41
1:B:101:THR:HG21	1:B:112:PRO:CD	2.41	0.41
1:B:78:VAL:HG11	1:B:88:LEU:HD11	2.02	0.41
1:C:46:LEU:C	1:C:46:LEU:HD12	2.40	0.41
1:D:199:MET:O	1:D:200:SER:OG	2.37	0.41
1:D:193:PRO:HA	2:D:330:NAD:O7N	2.21	0.41
1:D:11:LEU:HD21	1:D:39:THR:HG23	2.03	0.41
1:C:95:ILE:CD1	1:C:122:ILE:CG2	2.99	0.41
1:B:185:ARG:HA	1:B:254:THR:HG23	2.02	0.40
1:B:193:PRO:HA	2:B:310:NAD:O7N	2.21	0.40
1:B:39:THR:HA	1:B:61:LEU:O	2.22	0.40
1:C:195:ARG:HH21	1:C:219:GLU:HG3	1.85	0.40
1:D:219:GLU:HB2	1:D:232:MET:SD	2.61	0.40
1:A:109:PHE:HB3	1:B:132:LYS:HG3	2.03	0.40
1:C:4:LEU:N	1:C:32:GLN:HE21	1.98	0.40
1:A:149:PHE:CZ	1:A:222:TRP:CH2	3.09	0.40
1:A:195:ARG:NH1	1:A:195:ARG:CG	2.83	0.40
1:A:11:LEU:HD21	1:A:39:THR:HG23	2.03	0.40
1:A:67:ASN:ND2	1:A:67:ASN:C	2.71	0.40
1:D:185:ARG:CA	1:D:254:THR:CG2	2.98	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:95:ILE:HD13	2:D:330:NAD:C4A	2.50	0.40
1:B:198:ALA:HB1	3:B:410:JPM:C8	2.52	0.40
1:C:2:THR:HA	1:C:3:GLY:HA3	1.90	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	266/269 (99%)	251 (94%)	14 (5%)	1 (0%)	38	47
1	B	245/269 (91%)	229 (94%)	16 (6%)	0	100	100
1	C	257/269 (96%)	242 (94%)	15 (6%)	0	100	100
1	D	253/269 (94%)	235 (93%)	18 (7%)	0	100	100
All	All	1021/1076 (95%)	957 (94%)	63 (6%)	1 (0%)	55	67

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	195	ARG

### 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	203/205 (99%)	177 (87%)	26 (13%)	5	5
1	B	192/205 (94%)	166 (86%)	26 (14%)	4	4
1	C	200/205 (98%)	179 (90%)	21 (10%)	8	9
1	D	198/205 (97%)	172 (87%)	26 (13%)	5	5
All	All	793/820 (97%)	694 (88%)	99 (12%)	5	5

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LEU
1	A	6	ASP
1	A	18	ASP
1	A	35	GLN
1	A	53	ARG
1	A	54	LEU
1	A	57	LYS
1	A	67	ASN
1	A	74	LEU
1	A	79	THR
1	A	95	ILE
1	A	101	THR
1	A	162	THR
1	A	168	LEU
1	A	170	SER
1	A	188	LEU
1	A	199	MET
1	A	217	LEU
1	A	218	LEU
1	A	222	TRP
1	A	233	LYS
1	A	236	THR
1	A	238	VAL
1	A	250	LEU
1	A	253	THR
1	A	254	THR
1	B	5	LEU
1	B	6	ASP
1	B	11	LEU
1	B	43	ARG
1	B	49	ARG
1	B	54	LEU
1	B	57	LYS

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Mol	Chain	Res	Type
1	B	65	VAL
1	B	67	ASN
1	B	74	LEU
1	B	95	ILE
1	B	100	GLN
1	B	101	THR
1	B	121	HIS
1	B	168	LEU
1	B	188	LEU
1	B	196	THR
1	B	197	LEU
1	B	199	MET
1	B	200	SER
1	B	238	VAL
1	B	250	LEU
1	B	253	THR
1	B	254	THR
1	B	256	ASP
1	B	269	LEU
1	C	5	LEU
1	C	11	LEU
1	C	19	SER
1	C	45	ARG
1	C	53	ARG
1	C	54	LEU
1	C	65	VAL
1	C	67	ASN
1	C	69	GLU
1	C	74	LEU
1	C	95	ILE
1	C	100	GLN
1	C	121	HIS
1	C	168	LEU
1	C	170	SER
1	C	188	LEU
1	C	197	LEU
1	C	236	THR
1	C	250	LEU
1	C	253	THR
1	C	254	THR
1	D	5	LEU
1	D	6	ASP

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Mol	Chain	Res	Type
1	D	9	ARG
1	D	43	ARG
1	D	53	ARG
1	D	54	LEU
1	D	55	PRO
1	D	57	LYS
1	D	65	VAL
1	D	67	ASN
1	D	69	GLU
1	D	74	LEU
1	D	95	ILE
1	D	100	GLN
1	D	105	ILE
1	D	168	LEU
1	D	170	SER
1	D	188	LEU
1	D	196	THR
1	D	197	LEU
1	D	216	GLN
1	D	236	THR
1	D	238	VAL
1	D	250	LEU
1	D	254	THR
1	D	269	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	67	ASN
1	A	86	ASN
1	A	121	HIS
1	A	187	ASN
1	A	265	HIS
1	B	32	GLN
1	B	66	GLN
1	B	67	ASN
1	B	86	ASN
1	B	187	ASN
1	B	265	HIS
1	C	32	GLN
1	C	67	ASN

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Mol	Chain	Res	Type
1	C	86	ASN
1	C	100	GLN
1	C	121	HIS
1	C	187	ASN
1	C	265	HIS
1	D	32	GLN
1	D	67	ASN
1	D	86	ASN
1	D	121	HIS
1	D	187	ASN
1	D	265	HIS

### 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 carbohydrates 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
2	NAD	A	300	-	41,48,48	1.70	3 (7%)	43,73,73	2.10	5 (11%)
3	JPM	A	400	-	25,25,25	2.90	4 (16%)	34,34,34	1.33	5 (14%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAD	B	310	-	41,48,48	1.60	3 (7%)	43,73,73	2.13	4 (9%)
3	JPM	B	410	-	25,25,25	2.93	2 (8%)	34,34,34	1.27	3 (8%)
2	NAD	C	320	-	41,48,48	1.62	3 (7%)	43,73,73	2.33	6 (13%)
3	JPM	C	420	-	25,25,25	2.80	3 (12%)	34,34,34	1.21	2 (5%)
2	NAD	D	330	-	41,48,48	1.62	3 (7%)	43,73,73	2.10	5 (11%)
3	JPM	D	430	-	25,25,25	3.00	2 (8%)	34,34,34	1.09	4 (11%)

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	NAD	A	300	-	-	0/22/62/62	0/5/5/5
3	JPM	A	400	-	-	0/8/8/8	0/3/3/3
2	NAD	B	310	-	-	0/22/62/62	0/5/5/5
3	JPM	B	410	-	-	0/8/8/8	0/3/3/3
2	NAD	C	320	-	-	0/22/62/62	0/5/5/5
3	JPM	C	420	-	-	0/8/8/8	0/3/3/3
2	NAD	D	330	-	-	0/22/62/62	0/5/5/5
3	JPM	D	430	-	-	0/8/8/8	0/3/3/3

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	410	JPM	C11-CL5	-10.72	1.51	1.74
3	D	430	JPM	C11-CL5	-10.43	1.51	1.74
3	A	400	JPM	C11-CL5	-10.41	1.51	1.74
3	C	420	JPM	C11-CL5	-10.35	1.52	1.74
3	D	430	JPM	C7-CL1	-9.84	1.50	1.73
3	B	410	JPM	C7-CL1	-9.19	1.51	1.73
3	A	400	JPM	C7-CL1	-9.13	1.51	1.73
3	C	420	JPM	C7-CL1	-8.30	1.53	1.73
2	B	310	NAD	C2A-N1A	2.01	1.37	1.33
3	A	400	JPM	C12-C7	2.04	1.42	1.38
3	A	400	JPM	C12-C11	2.09	1.41	1.38
3	C	420	JPM	C10-C11	2.15	1.42	1.38
2	C	320	NAD	C2A-N1A	2.60	1.38	1.33
2	D	330	NAD	C2A-N1A	2.69	1.39	1.33
2	A	300	NAD	C2A-N1A	2.88	1.39	1.33
2	B	310	NAD	C2A-N3A	3.69	1.38	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	330	NAD	C2A-N3A	3.94	1.38	1.32
2	C	320	NAD	C2A-N3A	3.97	1.38	1.32
2	A	300	NAD	C2A-N3A	4.39	1.39	1.32
2	D	330	NAD	O7N-C7N	7.76	1.40	1.24
2	C	320	NAD	O7N-C7N	7.86	1.40	1.24
2	A	300	NAD	O7N-C7N	7.99	1.40	1.24
2	B	310	NAD	O7N-C7N	8.14	1.41	1.24

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	310	NAD	N3A-C2A-N1A	-11.78	118.60	128.86
2	A	300	NAD	N3A-C2A-N1A	-11.33	118.99	128.86
2	C	320	NAD	N3A-C2A-N1A	-11.33	118.99	128.86
2	D	330	NAD	N3A-C2A-N1A	-10.36	119.83	128.86
2	C	320	NAD	C4B-O4B-C1B	-6.48	102.87	109.77
2	D	330	NAD	O7N-C7N-C3N	-3.74	115.25	119.62
2	B	310	NAD	C4B-O4B-C1B	-3.45	106.10	109.77
2	A	300	NAD	C4B-O4B-C1B	-3.00	106.57	109.77
2	C	320	NAD	C4A-C5A-N7A	-2.99	106.52	109.41
3	A	400	JPM	C10-C11-C12	-2.83	117.69	121.51
3	B	410	JPM	C10-C11-C12	-2.80	117.72	121.51
3	C	420	JPM	C10-C11-C12	-2.60	118.00	121.51
2	C	320	NAD	C1B-N9A-C4A	-2.56	122.22	126.64
2	D	330	NAD	C4B-O4B-C1B	-2.53	107.08	109.77
3	D	430	JPM	C10-C11-C12	-2.50	118.13	121.51
2	C	320	NAD	O7N-C7N-N7N	-2.30	119.30	122.58
2	A	300	NAD	C5B-C4B-C3B	-2.15	107.08	115.29
2	D	330	NAD	O5B-C5B-C4B	-2.10	101.57	109.00
2	A	300	NAD	O5B-C5B-C4B	-2.07	101.67	109.00
2	B	310	NAD	O2N-PN-O1N	2.04	122.83	112.28
3	D	430	JPM	C9-C10-C11	2.23	121.61	119.24
3	A	400	JPM	C12-C7-CL1	2.29	122.11	118.50
3	D	430	JPM	C7-C12-C11	2.44	121.39	118.72
3	D	430	JPM	C8-O1-C3	2.76	124.66	117.98
3	A	400	JPM	C9-C10-C11	2.86	122.28	119.24
2	B	310	NAD	C3N-C7N-N7N	2.97	121.17	117.77
3	A	400	JPM	C12-C11-CL5	3.16	123.09	119.14
3	C	420	JPM	C7-C12-C11	3.16	122.17	118.72
2	A	300	NAD	C3N-C7N-N7N	3.32	121.57	117.77
3	B	410	JPM	C8-O1-C3	3.50	126.46	117.98
2	C	320	NAD	C3N-C7N-N7N	3.60	121.88	117.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	400	JPM	C8-O1-C3	3.72	127.00	117.98
3	B	410	JPM	C7-C12-C11	3.93	123.01	118.72
2	D	330	NAD	C3N-C7N-N7N	4.96	123.44	117.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	300	NAD	2	0
2	B	310	NAD	2	0
3	B	410	JPM	2	0
2	C	320	NAD	1	0
3	C	420	JPM	1	0
2	D	330	NAD	3	0
3	D	430	JPM	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	268/269 (99%)	-0.43	2 (0%) 87 90	27, 37, 57, 71	0
1	B	249/269 (92%)	-0.47	2 (0%) 86 89	27, 36, 53, 66	0
1	C	261/269 (97%)	-0.34	5 (1%) 67 73	29, 38, 55, 74	0
1	D	257/269 (95%)	-0.17	12 (4%) 32 39	30, 40, 59, 66	0
All	All	1035/1076 (96%)	-0.35	21 (2%) 65 72	27, 38, 57, 74	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	202	ILE	5.8
1	D	105	ILE	3.7
1	D	59	PRO	3.2
1	C	213	ALA	3.1
1	D	46	LEU	3.0
1	D	212	GLY	3.0
1	D	58	ALA	2.7
1	D	215	ILE	2.6
1	C	217	LEU	2.6
1	D	56	ALA	2.5
1	D	213	ALA	2.5
1	D	45	ARG	2.4
1	C	201	ALA	2.3
1	D	84	ALA	2.3
1	D	217	LEU	2.3
1	B	105	ILE	2.3
1	B	198	ALA	2.2
1	C	105	ILE	2.2
1	D	104	GLY	2.2
1	A	208	GLY	2.1
1	A	84	ALA	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
2	NAD	C	320	44/44	0.97	0.09	-0.25	28,34,37,38	0
3	JPM	A	400	23/23	0.99	0.10	-0.30	21,32,43,44	0
3	JPM	C	420	23/23	0.99	0.11	-0.30	21,34,39,41	0
3	JPM	D	430	23/23	0.96	0.10	-0.48	32,41,43,43	0
3	JPM	B	410	23/23	0.97	0.09	-0.60	36,42,48,49	0
2	NAD	A	300	44/44	0.97	0.08	-0.72	28,33,37,38	0
2	NAD	B	310	44/44	0.96	0.07	-0.82	36,41,44,44	0
2	NAD	D	330	44/44	0.94	0.08	-0.97	39,42,45,47	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.