



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 03:20 pm GMT

PDB ID : 4R9S  
Title : Mycobacterium tuberculosis InhA bound to NITD-916  
Authors : Noble, C.G.  
Deposited on : 2014-09-07  
Resolution : 3.20 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

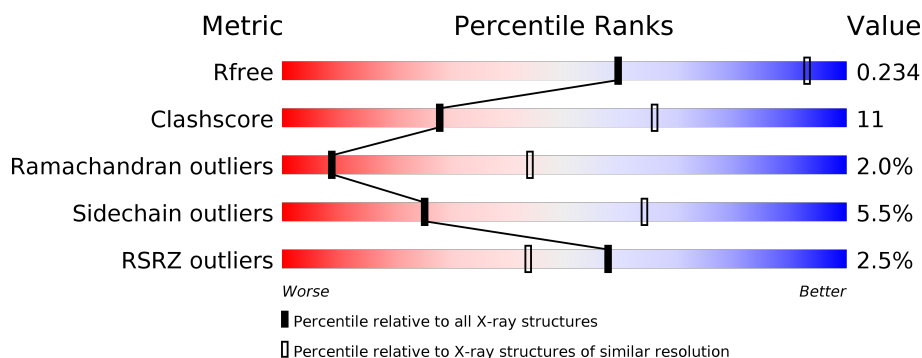
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.20 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))
$R_{free}$	100719	1015 (3.22-3.18)
Clashscore	112137	1009 (3.20-3.20)
Ramachandran outliers	110173	1118 (3.22-3.18)
Sidechain outliers	110143	1117 (3.22-3.18)
RSRZ outliers	101464	1020 (3.22-3.18)

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	272	<div> <div>71%</div> <div>26%</div> <div>..</div> </div>
1	C	272	<div> <div>74%</div> <div>22%</div> <div>..</div> </div>
1	E	272	<div> <div>4%</div> <div>76%</div> <div>21%</div> <div>..</div> </div>
1	G	272	<div> <div>5%</div> <div>63%</div> <div>34%</div> <div>..</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	3KY	A	302	-	-	-	X
3	3KY	E	302	-	-	X	-
3	3KY	G	302	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8271 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.

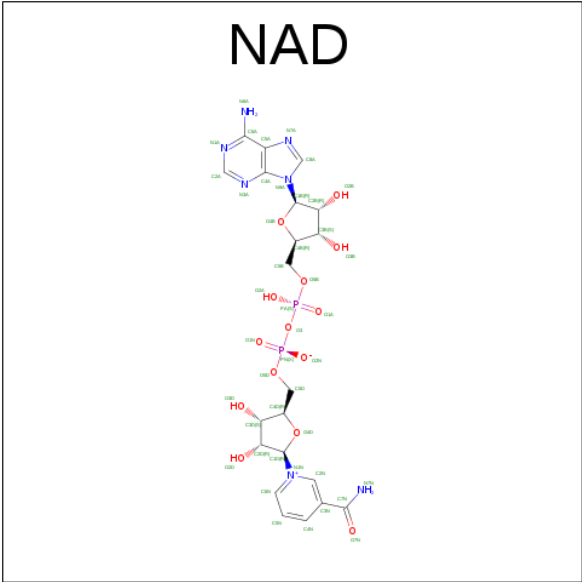
- Molecule 1 is a protein called Enoyl-[acyl-carrier-protein] reductase [NADH].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	0	0	0
			1996	1264	348	374	10			
1	C	268	Total	C	N	O	S	0	0	0
			1996	1264	348	374	10			
1	E	268	Total	C	N	O	S	0	0	0
			1996	1264	348	374	10			
1	G	268	Total	C	N	O	S	0	0	0
			1996	1264	348	374	10			

There are 12 discrepancies between the modelled and reference sequences:

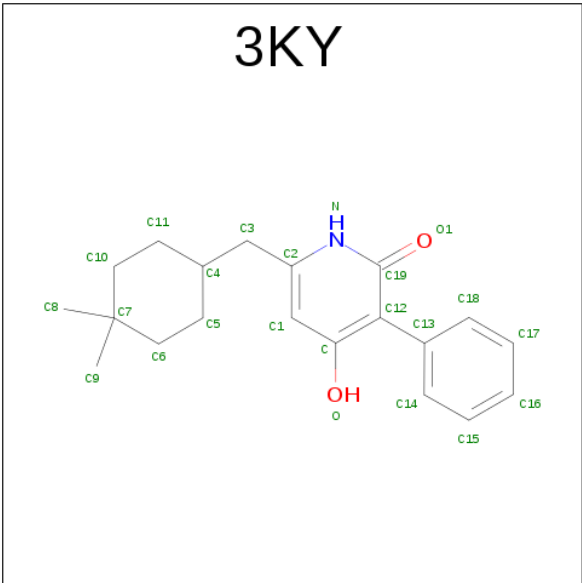
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP I6Y6N7
A	-1	SER	-	EXPRESSION TAG	UNP I6Y6N7
A	0	HIS	-	EXPRESSION TAG	UNP I6Y6N7
C	-2	GLY	-	EXPRESSION TAG	UNP I6Y6N7
C	-1	SER	-	EXPRESSION TAG	UNP I6Y6N7
C	0	HIS	-	EXPRESSION TAG	UNP I6Y6N7
E	-2	GLY	-	EXPRESSION TAG	UNP I6Y6N7
E	-1	SER	-	EXPRESSION TAG	UNP I6Y6N7
E	0	HIS	-	EXPRESSION TAG	UNP I6Y6N7
G	-2	GLY	-	EXPRESSION TAG	UNP I6Y6N7
G	-1	SER	-	EXPRESSION TAG	UNP I6Y6N7
G	0	HIS	-	EXPRESSION TAG	UNP I6Y6N7

- 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	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	E	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	G	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is 6-[(4,4-DIMETHYLCYCLOHEXYL)METHYL]-4-HYDROXY-3-PHENYLPYRIDIN-2(1H)-ONE (three-letter code: 3KY) (formula: C<sub>20</sub>H<sub>25</sub>NO<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			23	20	1	2		
3	C	1	Total	C	N	O	0	0
			23	20	1	2		
3	E	1	Total	C	N	O	0	0
			23	20	1	2		
3	G	1	Total	C	N	O	0	0
			23	20	1	2		

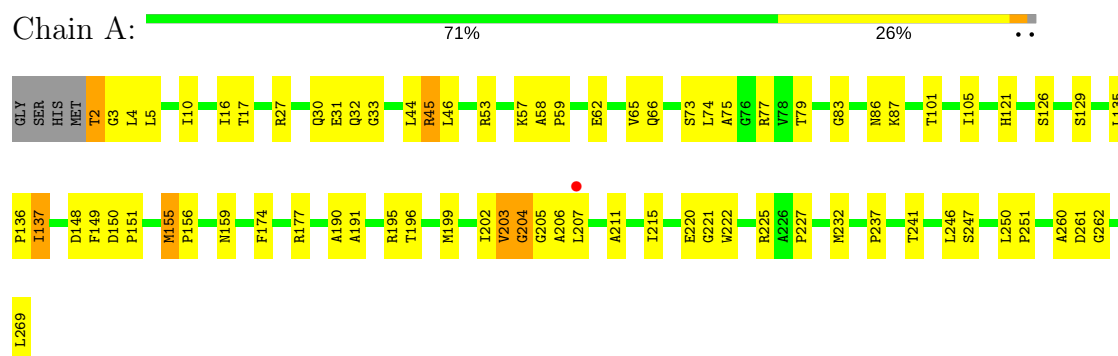
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	O	0	0
			3	3		
4	C	3	Total	O	0	0
			3	3		
4	E	8	Total	O	0	0
			8	8		
4	G	5	Total	O	0	0
			5	5		

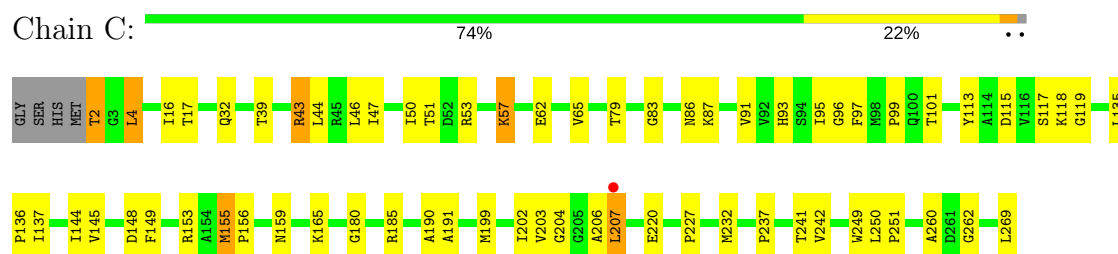
### 3 Residue-property plots [i](#)

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

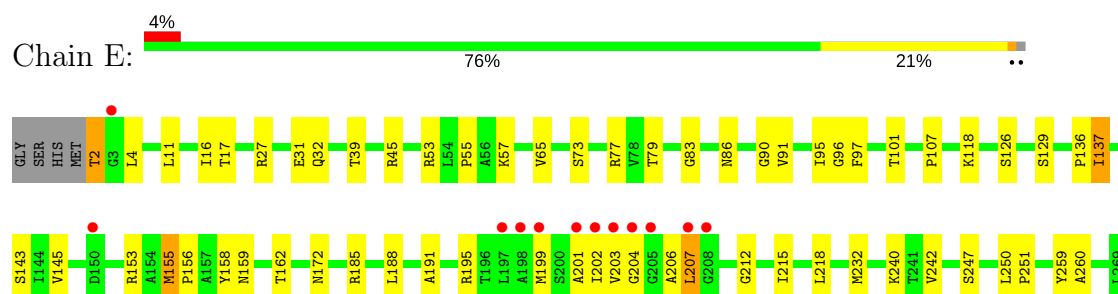
- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



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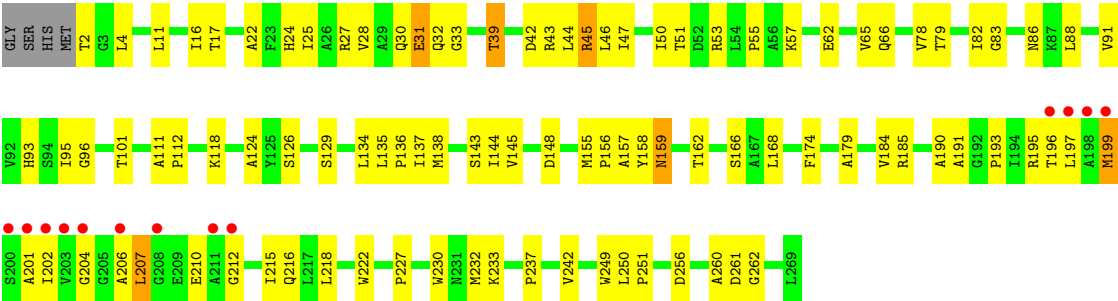


- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



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





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	103.58Å 103.58Å 121.39Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.00 – 3.20 44.85 – 3.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (45.00-3.20) 100.0 (44.85-3.20)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.82 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.177 , 0.235 0.174 , 0.234	Depositor DCC
$R_{free}$ test set	1230 reflections (5.40%)	DCC
Wilson B-factor (Å <sup>2</sup> )	111.8	Xtriage
Anisotropy	0.083	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 63.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.027 for -h,-k,l 0.059 for h,-h-k,-l 0.039 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8271	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	117.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.18% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 3KY, 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	0.39	1/2034 (0.0%)	0.59	0/2761
1	C	0.37	1/2034 (0.0%)	0.60	0/2761
1	E	0.38	1/2034 (0.0%)	0.57	0/2761
1	G	0.37	0/2034	0.57	0/2761
All	All	0.38	3/8136 (0.0%)	0.58	0/11044

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	2	THR	CB-CG2	5.92	1.71	1.52
1	C	2	THR	CB-CG2	5.26	1.69	1.52
1	E	2	THR	CB-CG2	5.20	1.69	1.52

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	1996	0	2013	48	0
1	C	1996	0	2013	40	0
1	E	1996	0	2013	37	0
1	G	1996	0	2013	61	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	44	0	26	2	0
2	C	44	0	26	3	0
2	E	44	0	26	3	0
2	G	44	0	26	3	0
3	A	23	0	25	4	0
3	C	23	0	24	5	0
3	E	23	0	25	9	0
3	G	23	0	25	11	0
4	A	3	0	0	3	0
4	C	3	0	0	0	0
4	E	8	0	0	4	0
4	G	5	0	0	4	0
All	All	8271	0	8255	184	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 11.

All (184) 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:G:195:ARG:HB2	4:G:403:HOH:O	1.60	1.01
2:G:301:NAD:H2D	3:G:302:3KY:H1	1.55	0.86
1:G:215:ILE:HD13	3:G:302:3KY:H21	1.61	0.83
1:E:215:ILE:HD13	3:E:302:3KY:H21	1.60	0.82
1:E:77:ARG:HB2	4:E:407:HOH:O	1.80	0.80
1:E:16:ILE:HG23	1:E:17:THR:HG23	1.66	0.76
1:G:210:GLU:HA	4:G:405:HOH:O	1.86	0.74
2:G:301:NAD:H2D	3:G:302:3KY:C18	2.20	0.71
1:G:96:GLY:O	3:G:302:3KY:H3	1.91	0.71
1:A:46:LEU:HA	4:A:401:HOH:O	1.89	0.70
1:E:96:GLY:O	3:E:302:3KY:H3	1.92	0.70
1:G:45:ARG:NH2	1:G:45:ARG:HB2	2.08	0.68
1:E:2:THR:HG1	1:G:2:THR:N	1.93	0.67
1:G:4:LEU:H	1:G:32:GLN:HE21	1.42	0.66
1:A:5:LEU:HB2	4:A:403:HOH:O	1.96	0.66
3:C:302:3KY:H11	3:C:302:3KY:H17	1.77	0.66
1:A:83:GLY:O	1:A:86:ASN:HB2	1.96	0.66
1:A:2:THR:N	1:C:2:THR:HG1	1.95	0.65
1:E:65:VAL:HG22	2:E:301:NAD:N1A	2.12	0.65
1:C:96:GLY:O	3:C:302:3KY:H3	1.97	0.64
1:G:233:LYS:HG3	4:G:403:HOH:O	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:44:LEU:HD21	1:C:62:GLU:HG3	1.79	0.63
1:G:148:ASP:O	1:G:190:ALA:HA	1.98	0.62
1:A:202:ILE:HG12	1:A:207:LEU:HB2	1.80	0.62
1:G:212:GLY:O	1:G:216:GLN:HB2	2.01	0.61
1:E:191:ALA:HA	1:E:260:ALA:O	2.02	0.59
1:A:105:ILE:HG23	1:A:211:ALA:HB2	1.85	0.59
1:G:126:SER:HA	1:G:129:SER:HB2	1.85	0.59
1:G:65:VAL:HG22	2:G:301:NAD:N1A	2.18	0.58
3:A:302:3KY:H11	3:A:302:3KY:H17	1.85	0.57
1:A:149:PHE:CE2	3:A:302:3KY:H11	2.38	0.57
1:A:221:GLY:O	1:A:225:ARG:HG3	2.05	0.57
1:G:16:ILE:HG23	1:G:17:THR:HG23	1.86	0.56
1:C:99:PRO:HG2	1:C:115:ASP:HB3	1.88	0.56
1:E:136:PRO:HD2	4:E:406:HOH:O	2.05	0.56
1:C:91:VAL:HB	1:C:144:ILE:HG23	1.88	0.56
1:C:83:GLY:O	1:C:86:ASN:HB2	2.07	0.55
1:A:241:THR:HG23	1:C:250:LEU:HD23	1.88	0.55
2:C:301:NAD:H2D	3:C:302:3KY:C18	2.36	0.54
3:E:302:3KY:H17	3:E:302:3KY:H11	1.87	0.54
1:G:202:ILE:HG23	1:G:207:LEU:HB3	1.88	0.54
1:C:97:PHE:HE2	1:C:118:LYS:HD3	1.72	0.54
1:C:4:LEU:HB2	1:C:32:GLN:HE21	1.73	0.54
1:G:78:VAL:HG11	1:G:88:LEU:HD11	1.89	0.54
1:A:44:LEU:HD21	1:A:62:GLU:HG3	1.90	0.53
1:E:137:ILE:HG13	1:E:137:ILE:O	2.08	0.53
3:G:302:3KY:H11	3:G:302:3KY:H17	1.88	0.53
1:G:30:GLN:O	1:G:33:GLY:N	2.34	0.53
1:E:259:TYR:HB2	1:G:256:ASP:OD2	2.08	0.53
1:G:157:ALA:O	1:G:159:ASN:N	2.41	0.53
1:G:134:LEU:O	1:G:138:MET:HG3	2.10	0.53
2:C:301:NAD:H2D	3:C:302:3KY:H1	1.92	0.52
1:G:27:ARG:HG2	1:G:31:GLU:OE2	2.09	0.52
1:G:46:LEU:O	1:G:50:ILE:HG12	2.08	0.52
1:E:195:ARG:HH21	1:E:203:VAL:HG21	1.73	0.52
1:E:27:ARG:HG2	1:E:31:GLU:OE2	2.10	0.52
1:A:65:VAL:HG22	2:A:301:NAD:N1A	2.24	0.52
1:G:17:THR:HG21	1:G:197:LEU:HD22	1.89	0.52
1:A:53:ARG:N	1:A:53:ARG:HD2	2.25	0.52
1:E:202:ILE:HG12	1:E:207:LEU:HB2	1.90	0.52
1:C:149:PHE:H	1:C:165:LYS:HD2	1.74	0.51
1:A:215:ILE:HD13	3:A:302:3KY:H21	1.90	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:301:NAD:H2D	3:E:302:3KY:C18	2.41	0.51
1:C:46:LEU:O	1:C:50:ILE:HG12	2.11	0.51
1:C:159:ASN:HA	1:G:174:PHE:CE2	2.45	0.51
1:E:4:LEU:H	1:E:32:GLN:HB3	1.75	0.51
1:G:193:PRO:HB3	3:G:302:3KY:C5	2.41	0.51
1:G:22:ALA:HA	1:G:25:ILE:HD12	1.93	0.51
1:A:250:LEU:HD23	1:C:241:THR:HG23	1.92	0.51
1:A:135:LEU:N	1:A:136:PRO:HD2	2.26	0.51
1:A:3:GLY:HA3	1:A:32:GLN:O	2.11	0.51
1:A:2:THR:HG1	1:C:2:THR:N	2.09	0.50
1:E:83:GLY:O	1:E:86:ASN:HB2	2.12	0.50
1:G:124:ALA:HA	1:G:168:LEU:HD13	1.93	0.50
1:A:174:PHE:O	1:A:177:ARG:HB2	2.11	0.50
1:A:87:LYS:HE2	1:A:137:ILE:HA	1.93	0.50
1:C:87:LYS:HE2	1:C:137:ILE:HA	1.94	0.50
1:A:227:PRO:O	1:C:180:GLY:HA3	2.12	0.49
1:G:91:VAL:HB	1:G:144:ILE:HG23	1.94	0.49
1:C:202:ILE:HG12	1:C:207:LEU:HB2	1.93	0.49
1:E:218:LEU:HD12	3:E:302:3KY:H22	1.95	0.49
1:E:73:SER:O	1:E:77:ARG:HG3	2.12	0.49
1:G:191:ALA:HA	1:G:260:ALA:O	2.12	0.49
1:C:65:VAL:HG22	2:C:301:NAD:N1A	2.28	0.49
1:C:249:TRP:C	1:C:251:PRO:HD3	2.33	0.49
1:G:218:LEU:HD12	3:G:302:3KY:H22	1.94	0.49
1:A:45:ARG:HB2	1:A:45:ARG:HH21	1.77	0.48
1:G:179:ALA:HB1	1:G:184:VAL:HB	1.94	0.48
1:G:78:VAL:O	1:G:82:ILE:HG12	2.13	0.48
1:C:16:ILE:HG23	1:C:17:THR:HG23	1.95	0.48
1:G:45:ARG:CZ	1:G:45:ARG:HB2	2.43	0.48
1:A:227:PRO:HD2	1:A:262:GLY:O	2.13	0.48
2:E:301:NAD:H2D	3:E:302:3KY:H1	1.96	0.47
1:G:201:ALA:O	1:G:206:ALA:HB3	2.15	0.47
1:C:113:TYR:CE2	1:C:117:SER:HB2	2.49	0.47
1:G:145:VAL:HG11	1:G:242:VAL:HG13	1.97	0.47
1:E:212:GLY:C	4:E:402:HOH:O	2.53	0.47
3:G:302:3KY:C11	3:G:302:3KY:N	2.78	0.47
1:A:237:PRO:HB3	1:C:251:PRO:HG2	1.96	0.47
1:E:136:PRO:CD	4:E:406:HOH:O	2.61	0.47
1:E:201:ALA:O	1:E:206:ALA:HB3	2.14	0.47
1:A:16:ILE:HG23	1:A:17:THR:HG23	1.96	0.46
1:G:11:LEU:HD21	1:G:39:THR:HG22	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:45:ARG:HH21	1:G:45:ARG:HB2	1.79	0.46
1:G:249:TRP:C	1:G:251:PRO:HD3	2.36	0.46
1:A:215:ILE:HD13	3:A:302:3KY:H23	1.96	0.46
1:G:227:PRO:HD2	1:G:262:GLY:O	2.15	0.46
1:E:251:PRO:HG2	1:G:237:PRO:HB3	1.98	0.46
1:C:153:ARG:CZ	1:E:153:ARG:CZ	2.94	0.46
1:G:135:LEU:N	1:G:136:PRO:HD2	2.31	0.46
1:A:155:MET:HB2	1:A:156:PRO:HD2	1.99	0.45
1:A:27:ARG:HG2	1:A:31:GLU:OE2	2.16	0.45
1:A:148:ASP:O	1:A:190:ALA:HA	2.15	0.45
1:G:193:PRO:HB3	3:G:302:3KY:H25	1.99	0.45
1:A:45:ARG:H	1:A:45:ARG:HG3	1.57	0.45
1:C:47:ILE:O	1:C:51:THR:HG23	2.16	0.45
1:E:172:ASN:HB2	1:E:188:LEU:HD11	1.99	0.45
1:E:90:GLY:HA2	1:E:143:SER:O	2.17	0.45
1:A:66:GLN:HG2	1:A:121:HIS:CE1	2.51	0.45
1:G:199:MET:SD	4:G:404:HOH:O	2.61	0.45
1:G:202:ILE:HG12	1:G:207:LEU:HB2	1.99	0.45
1:A:150:ASP:HA	1:A:151:PRO:HD3	1.88	0.44
1:A:58:ALA:HB1	1:A:59:PRO:HD2	2.00	0.44
1:C:135:LEU:N	1:C:136:PRO:HD2	2.32	0.44
1:E:11:LEU:HD12	1:E:91:VAL:HG13	2.00	0.44
1:C:185:ARG:NH1	1:C:250:LEU:O	2.50	0.44
1:A:251:PRO:HG2	1:C:237:PRO:HB3	1.98	0.44
1:E:158:TYR:HB2	3:E:302:3KY:H18	2.00	0.44
1:E:97:PHE:HE2	1:E:118:LYS:HD3	1.82	0.44
1:C:44:LEU:HD21	1:C:62:GLU:CG	2.48	0.44
1:A:33:GLY:C	4:A:403:HOH:O	2.56	0.44
1:C:269:LEU:HD22	1:E:218:LEU:HD23	1.99	0.44
1:C:148:ASP:O	1:C:190:ALA:HA	2.17	0.44
1:C:145:VAL:HG11	1:C:242:VAL:HG13	2.00	0.44
1:C:53:ARG:HD2	1:C:53:ARG:N	2.33	0.44
1:A:10:ILE:HD13	1:A:246:LEU:HD13	2.00	0.43
1:G:222:TRP:HE1	1:G:261:ASP:HB2	1.82	0.43
1:A:73:SER:O	1:A:77:ARG:HG3	2.18	0.43
1:G:193:PRO:HB3	3:G:302:3KY:H24	2.00	0.43
1:A:191:ALA:HA	1:A:260:ALA:O	2.18	0.43
1:C:191:ALA:HA	1:C:260:ALA:O	2.18	0.43
1:A:30:GLN:O	1:A:33:GLY:N	2.50	0.43
1:G:156:PRO:O	1:G:157:ALA:HB3	2.18	0.43
1:G:83:GLY:O	1:G:86:ASN:HB2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:96:GLY:O	3:G:302:3KY:C16	2.64	0.43
1:E:4:LEU:H	1:E:32:GLN:HE21	1.67	0.43
1:G:44:LEU:HD21	1:G:62:GLU:HG3	2.01	0.43
1:C:93:HIS:CE1	1:C:95:ILE:HB	2.54	0.43
1:E:240:LYS:HD2	1:G:251:PRO:HG3	2.01	0.43
1:G:193:PRO:HD2	1:G:230:TRP:NE1	2.34	0.43
1:G:4:LEU:N	1:G:32:GLN:HE21	2.13	0.43
1:C:86:ASN:HA	1:C:86:ASN:HD22	1.67	0.42
1:A:45:ARG:HB2	1:A:45:ARG:NH2	2.33	0.42
1:A:74:LEU:O	1:A:75:ALA:C	2.57	0.42
1:E:96:GLY:O	3:E:302:3KY:C16	2.62	0.42
1:C:97:PHE:O	1:C:119:GLY:HA2	2.18	0.42
1:E:4:LEU:HD22	1:G:249:TRP:CE2	2.54	0.42
1:E:126:SER:HA	1:E:129:SER:HB2	2.02	0.42
1:E:145:VAL:HG11	1:E:242:VAL:HG13	2.01	0.42
1:G:185:ARG:NH1	1:G:250:LEU:O	2.52	0.42
1:C:43:ARG:HB3	1:C:43:ARG:HH11	1.85	0.42
1:A:203:VAL:HG12	1:A:203:VAL:O	2.19	0.42
3:E:302:3KY:N	3:E:302:3KY:C11	2.82	0.42
1:C:227:PRO:HD2	1:C:262:GLY:O	2.19	0.42
1:G:111:ALA:HA	1:G:112:PRO:HD3	1.92	0.42
1:A:4:LEU:H	1:A:32:GLN:HB3	1.84	0.41
1:E:95:ILE:HG22	1:E:96:GLY:N	2.35	0.41
1:G:24:HIS:O	1:G:28:VAL:HG23	2.20	0.41
1:A:222:TRP:HE1	1:A:261:ASP:HB2	1.85	0.41
1:C:155:MET:HB2	1:C:156:PRO:HD2	2.02	0.41
1:G:143:SER:HB2	1:G:185:ARG:NH2	2.36	0.41
1:A:126:SER:HA	1:A:129:SER:HB2	2.03	0.41
3:C:302:3KY:C4	3:C:302:3KY:H17	2.46	0.41
1:E:185:ARG:NH1	1:E:250:LEU:O	2.54	0.41
1:G:93:HIS:CE1	1:G:95:ILE:HB	2.56	0.41
1:G:43:ARG:O	1:G:47:ILE:HG13	2.21	0.41
1:A:196:THR:HG21	2:A:301:NAD:O1N	2.21	0.40
1:A:202:ILE:CG2	1:A:215:ILE:HG13	2.51	0.40
1:A:195:ARG:HH21	1:A:203:VAL:HG21	1.86	0.40
1:E:155:MET:HB2	1:E:156:PRO:HD2	2.03	0.40
1:A:204:GLY:O	1:A:206:ALA:N	2.51	0.40
1:C:57:LYS:HD3	1:C:57:LYS:H	1.87	0.40
1:G:47:ILE:O	1:G:51:THR:HG23	2.21	0.40
1:G:66:GLN:HE22	1:G:118:LYS:HG3	1.87	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	266/272 (98%)	243 (91%)	18 (7%)	5 (2%)	9	46
1	C	266/272 (98%)	241 (91%)	22 (8%)	3 (1%)	17	58
1	E	266/272 (98%)	237 (89%)	24 (9%)	5 (2%)	9	46
1	G	266/272 (98%)	239 (90%)	19 (7%)	8 (3%)	5	32
All	All	1064/1088 (98%)	960 (90%)	83 (8%)	21 (2%)	9	44

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	159	ASN
1	G	204	GLY
1	C	203	VAL
1	G	158	TYR
1	G	159	ASN
1	G	196	THR
1	A	203	VAL
1	E	204	GLY
1	G	42	ASP
1	E	107	PRO
1	A	159	ASN
1	A	205	GLY
1	C	206	ALA
1	G	31	GLU
1	G	55	PRO
1	A	204	GLY
1	G	137	ILE
1	C	204	GLY
1	A	137	ILE
1	E	55	PRO
1	E	137	ILE

### 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	204/207 (99%)	194 (95%)	10 (5%)	29	68
1	C	204/207 (99%)	193 (95%)	11 (5%)	26	65
1	E	204/207 (99%)	192 (94%)	12 (6%)	23	62
1	G	204/207 (99%)	192 (94%)	12 (6%)	23	62
All	All	816/828 (99%)	771 (94%)	45 (6%)	25	64

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	ARG
1	A	57	LYS
1	A	79	THR
1	A	101	THR
1	A	155	MET
1	A	199	MET
1	A	220	GLU
1	A	232	MET
1	A	247	SER
1	A	269	LEU
1	C	4	LEU
1	C	39	THR
1	C	43	ARG
1	C	57	LYS
1	C	79	THR
1	C	101	THR
1	C	155	MET
1	C	199	MET
1	C	207	LEU
1	C	220	GLU
1	C	232	MET
1	E	39	THR
1	E	45	ARG
1	E	53	ARG

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Mol	Chain	Res	Type
1	E	57	LYS
1	E	79	THR
1	E	101	THR
1	E	155	MET
1	E	162	THR
1	E	199	MET
1	E	207	LEU
1	E	232	MET
1	E	247	SER
1	G	39	THR
1	G	45	ARG
1	G	53	ARG
1	G	57	LYS
1	G	79	THR
1	G	101	THR
1	G	155	MET
1	G	162	THR
1	G	166	SER
1	G	199	MET
1	G	207	LEU
1	G	232	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	86	ASN
1	C	32	GLN
1	C	86	ASN
1	C	187	ASN
1	C	265	HIS
1	E	32	GLN
1	E	86	ASN
1	E	100	GLN
1	G	32	GLN
1	G	66	GLN
1	G	86	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 5.5 Carbohydrates [i](#)

There are no 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	301	-	41,48,48	0.87	1 (2%)	43,73,73	1.49	5 (11%)
3	3KY	A	302	-	24,25,25	1.59	4 (16%)	31,36,36	2.08	5 (16%)
2	NAD	C	301	-	41,48,48	0.90	1 (2%)	43,73,73	1.55	5 (11%)
3	3KY	C	302	-	24,25,25	1.43	3 (12%)	31,36,36	2.10	4 (12%)
2	NAD	E	301	-	41,48,48	0.91	1 (2%)	43,73,73	1.55	4 (9%)
3	3KY	E	302	-	24,25,25	1.62	4 (16%)	31,36,36	2.15	5 (16%)
2	NAD	G	301	-	41,48,48	0.92	2 (4%)	43,73,73	1.50	5 (11%)
3	3KY	G	302	-	24,25,25	1.47	3 (12%)	31,36,36	2.15	6 (19%)

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	301	-	-	0/22/62/62	0/5/5/5
3	3KY	A	302	-	-	0/8/20/20	0/3/3/3
2	NAD	C	301	-	-	0/22/62/62	0/5/5/5
3	3KY	C	302	-	-	0/8/20/20	1/3/3/3
2	NAD	E	301	-	-	0/22/62/62	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	3KY	E	302	-	-	0/8/20/20	1/3/3/3
2	NAD	G	301	-	-	0/22/62/62	0/5/5/5
3	3KY	G	302	-	-	0/8/20/20	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	302	3KY	C12-C13	-2.42	1.46	1.49
3	A	302	3KY	C12-C13	-2.13	1.47	1.49
2	G	301	NAD	C2A-N3A	2.05	1.35	1.32
3	G	302	3KY	C2-N	2.71	1.39	1.34
3	C	302	3KY	C2-N	3.05	1.40	1.34
2	A	301	NAD	C5A-C4A	3.09	1.47	1.40
3	A	302	3KY	C2-N	3.23	1.40	1.34
2	C	301	NAD	C5A-C4A	3.24	1.47	1.40
3	E	302	3KY	C2-N	3.30	1.40	1.34
2	E	301	NAD	C5A-C4A	3.36	1.48	1.40
2	G	301	NAD	C5A-C4A	3.47	1.48	1.40
3	C	302	3KY	C19-N	3.51	1.39	1.33
3	C	302	3KY	C1-C	3.58	1.43	1.38
3	G	302	3KY	C19-N	3.65	1.39	1.33
3	A	302	3KY	C19-N	3.96	1.40	1.33
3	G	302	3KY	C1-C	4.17	1.44	1.38
3	E	302	3KY	C1-C	4.22	1.44	1.38
3	E	302	3KY	C19-N	4.27	1.40	1.33
3	A	302	3KY	C1-C	4.61	1.45	1.38

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	302	3KY	C12-C19-N	-8.94	115.54	124.12
3	C	302	3KY	C12-C19-N	-8.87	115.61	124.12
3	G	302	3KY	C12-C19-N	-8.71	115.76	124.12
3	A	302	3KY	C12-C19-N	-8.70	115.77	124.12
2	A	301	NAD	N3A-C2A-N1A	-6.34	123.33	128.86
2	C	301	NAD	N3A-C2A-N1A	-5.97	123.66	128.86
2	E	301	NAD	N3A-C2A-N1A	-5.90	123.72	128.86
2	G	301	NAD	N3A-C2A-N1A	-5.28	124.26	128.86
2	C	301	NAD	C4B-O4B-C1B	-4.08	105.43	109.77
2	G	301	NAD	C4B-O4B-C1B	-3.80	105.72	109.77
2	E	301	NAD	C4B-O4B-C1B	-3.56	105.98	109.77
2	A	301	NAD	C4B-O4B-C1B	-3.42	106.13	109.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	302	3KY	C1-C2-N	-3.03	119.63	122.91
3	C	302	3KY	C1-C2-N	-3.00	119.66	122.91
2	E	301	NAD	C4A-C5A-N7A	-2.95	106.56	109.41
3	A	302	3KY	C1-C2-N	-2.93	119.74	122.91
2	G	301	NAD	C4A-C5A-N7A	-2.84	106.66	109.41
3	E	302	3KY	C1-C2-N	-2.81	119.87	122.91
2	C	301	NAD	C4A-C5A-N7A	-2.65	106.85	109.41
2	A	301	NAD	C4A-C5A-N7A	-2.65	106.85	109.41
2	A	301	NAD	C1B-N9A-C4A	-2.49	122.33	126.64
2	C	301	NAD	C5N-C4N-C3N	-2.08	117.90	120.35
2	G	301	NAD	O7N-C7N-C3N	-2.05	117.22	119.62
3	G	302	3KY	C3-C2-C1	2.02	125.50	120.83
3	E	302	3KY	C11-C10-C7	2.02	116.65	112.94
2	A	301	NAD	C2A-N1A-C6A	2.09	122.42	118.77
3	C	302	3KY	C6-C5-C4	2.10	114.37	110.38
2	G	301	NAD	C3N-C7N-N7N	2.15	120.23	117.77
3	E	302	3KY	C6-C5-C4	2.17	114.50	110.38
3	A	302	3KY	C5-C6-C7	2.24	117.06	112.94
3	G	302	3KY	C11-C10-C7	2.27	117.11	112.94
3	A	302	3KY	C11-C10-C7	2.46	117.47	112.94
3	G	302	3KY	C5-C6-C7	2.53	117.60	112.94
2	E	301	NAD	C3N-C7N-N7N	2.66	120.82	117.77
2	C	301	NAD	C2N-C3N-C4N	2.85	121.52	118.26
3	C	302	3KY	C19-N-C2	4.84	123.77	117.26
3	A	302	3KY	C19-N-C2	5.12	124.15	117.26
3	E	302	3KY	C19-N-C2	5.17	124.21	117.26
3	G	302	3KY	C19-N-C2	5.34	124.44	117.26

There are no chirality outliers.

There are no torsion outliers.

All (2) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	302	3KY	C10-C11-C4-C5-C6-C7
3	C	302	3KY	C10-C11-C4-C5-C6-C7

8 monomers are involved in 34 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	NAD	2	0
3	A	302	3KY	4	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	301	NAD	3	0
3	C	302	3KY	5	0
2	E	301	NAD	3	0
3	E	302	3KY	9	0
2	G	301	NAD	3	0
3	G	302	3KY	11	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	268/272 (98%)	-0.37	1 (0%)	92 89	74, 102, 142, 215	0
1	C	268/272 (98%)	-0.43	1 (0%)	92 89	76, 100, 133, 213	0
1	E	268/272 (98%)	-0.08	12 (4%)	34 21	85, 126, 211, 299	0
1	G	268/272 (98%)	0.05	13 (4%)	30 18	88, 117, 235, 314	0
All	All	1072/1088 (98%)	-0.21	27 (2%)	58 43	74, 111, 176, 314	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	198	ALA	11.1
1	G	199	MET	10.5
1	G	200	SER	8.9
1	G	204	GLY	8.1
1	G	203	VAL	7.0
1	G	196	THR	6.0
1	G	197	LEU	6.0
1	A	207	LEU	5.2
1	G	201	ALA	4.8
1	E	199	MET	4.5
1	G	208	GLY	3.7
1	E	202	ILE	3.6
1	C	207	LEU	3.5
1	G	202	ILE	3.2
1	E	150	ASP	3.1
1	E	204	GLY	2.9
1	E	203	VAL	2.8
1	E	201	ALA	2.7
1	E	198	ALA	2.7
1	G	206	ALA	2.6
1	E	3	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	G	212	GLY	2.4
1	E	207	LEU	2.3
1	E	197	LEU	2.2
1	E	208	GLY	2.2
1	E	205	GLY	2.2
1	G	211	ALA	2.1

## 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	3KY	A	302	23/23	0.97	0.26	2.06	45,49,55,57	0
3	3KY	C	302	23/23	0.96	0.22	1.31	43,48,53,56	0
3	3KY	G	302	23/23	0.95	0.28	0.04	47,50,54,57	0
3	3KY	E	302	23/23	0.95	0.22	-0.25	47,51,57,59	0
2	NAD	E	301	44/44	0.94	0.19	-0.35	97,114,132,138	0
2	NAD	C	301	44/44	0.96	0.18	-0.42	88,96,106,116	0
2	NAD	G	301	44/44	0.96	0.14	-0.58	89,104,116,125	0
2	NAD	A	301	44/44	0.96	0.16	-0.70	85,102,121,122	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.