



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:48 PM GMT

PDB ID : 4G73  
Title : Crystal structure of NDH with NADH and Quinone  
Authors : Li, W.; Feng, Y.; Ge, J.; Yang, M.  
Deposited on : 2012-07-19  
Resolution : 2.52 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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
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) : trunk26865

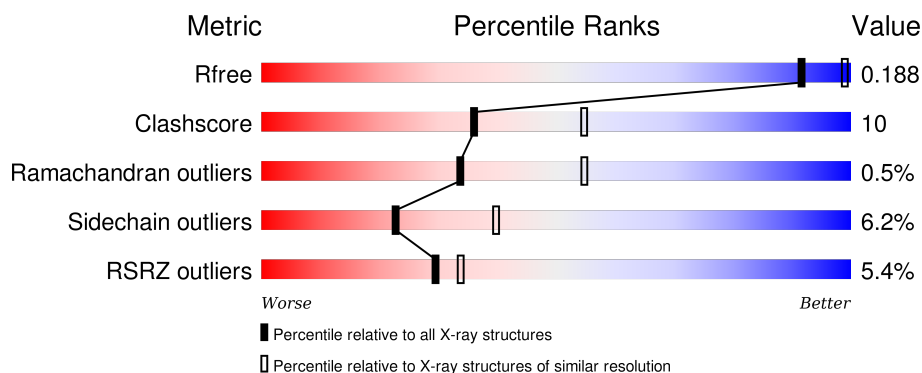
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.52 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	4241 (2.54-2.50)
Clashscore	102246	4968 (2.54-2.50)
Ramachandran outliers	100387	4873 (2.54-2.50)
Sidechain outliers	100360	4875 (2.54-2.50)
RSRZ outliers	91569	4253 (2.54-2.50)

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	502	<div> <div>4%</div> <div>77%</div> <div>15%</div> <div>6%</div> </div>
1	B	502	<div> <div>6%</div> <div>77%</div> <div>16%</div> <div>6%</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
2	UQ5	A	601	-	-	-	X
2	UQ5	A	602	-	X	-	-
2	UQ5	A	604	-	-	X	X
2	UQ5	B	605	-	-	X	X
3	MG	A	609	-	-	-	X
3	MG	B	608	-	-	-	X
3	MG	B	610	-	-	-	X
5	NAI	B	603	-	-	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8170 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.

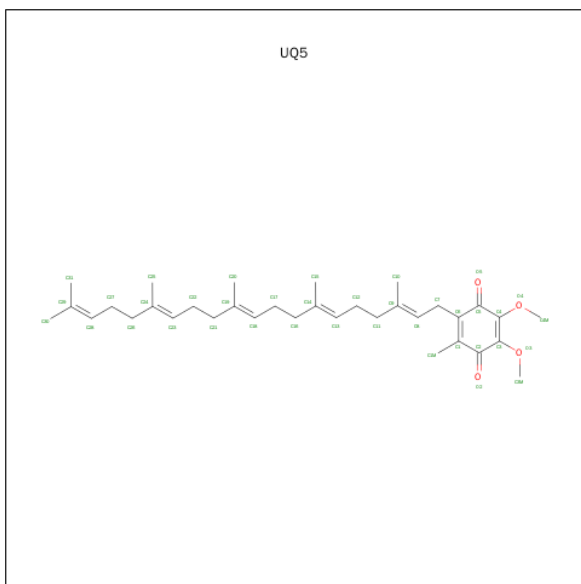
- Molecule 1 is a protein called Rotenone-insensitive NADH-ubiquinone oxidoreductase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	472	Total	C	N	O	S	0	0	0
			3738	2415	635	683	5			
1	B	472	Total	C	N	O	S	0	0	0
			3738	2415	635	683	5			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	12	MET	-	EXPRESSION TAG	UNP P32340
A	13	ARG	-	EXPRESSION TAG	UNP P32340
A	14	GLY	-	EXPRESSION TAG	UNP P32340
A	15	SER	-	EXPRESSION TAG	UNP P32340
A	16	HIS	-	EXPRESSION TAG	UNP P32340
A	17	HIS	-	EXPRESSION TAG	UNP P32340
A	18	HIS	-	EXPRESSION TAG	UNP P32340
A	19	HIS	-	EXPRESSION TAG	UNP P32340
A	20	HIS	-	EXPRESSION TAG	UNP P32340
A	21	HIS	-	EXPRESSION TAG	UNP P32340
A	22	GLY	-	EXPRESSION TAG	UNP P32340
A	23	SER	-	EXPRESSION TAG	UNP P32340
B	12	MET	-	EXPRESSION TAG	UNP P32340
B	13	ARG	-	EXPRESSION TAG	UNP P32340
B	14	GLY	-	EXPRESSION TAG	UNP P32340
B	15	SER	-	EXPRESSION TAG	UNP P32340
B	16	HIS	-	EXPRESSION TAG	UNP P32340
B	17	HIS	-	EXPRESSION TAG	UNP P32340
B	18	HIS	-	EXPRESSION TAG	UNP P32340
B	19	HIS	-	EXPRESSION TAG	UNP P32340
B	20	HIS	-	EXPRESSION TAG	UNP P32340
B	21	HIS	-	EXPRESSION TAG	UNP P32340
B	22	GLY	-	EXPRESSION TAG	UNP P32340
B	23	SER	-	EXPRESSION TAG	UNP P32340

- Molecule 2 is 2,3-DIMETHOXY-5-METHYL-6-(3,11,15,19-TETRAMETHYL-EICOSA-2,6,10,14,18-PENTAENYL)-[1,4]BENZOQUINONE (three-letter code: UQ5) (formula:  $C_{34}H_{50}O_4$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C 20 20	0	0
2	A	1	Total C 10 10	0	0
2	A	1	Total C O 24 20 4	0	0
2	B	1	Total C 20 20	0	0
2	B	1	Total C O 28 24 4	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

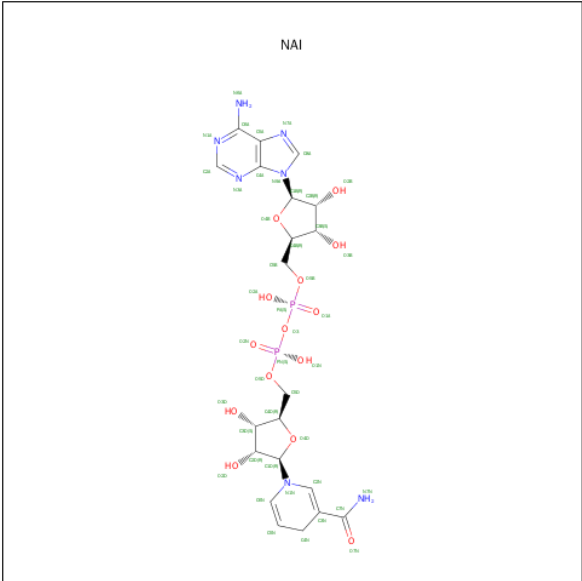
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	6	Total Mg 6 6	0	0
3	A	5	Total Mg 5 5	0	0

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:  $C_{27}H_{33}N_9O_{15}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
4	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 5 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (three-letter code: NAI) (formula: C<sub>21</sub>H<sub>29</sub>N<sub>7</sub>O<sub>14</sub>P<sub>2</sub>).



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

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

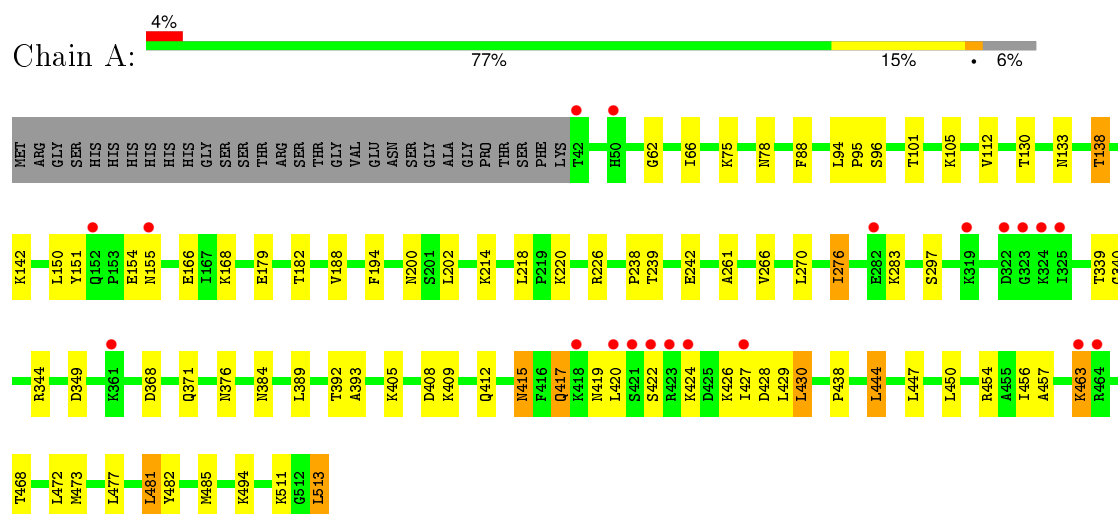
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	198	Total	O	0	0
			198	198		
6	B	189	Total	O	0	0
			189	189		

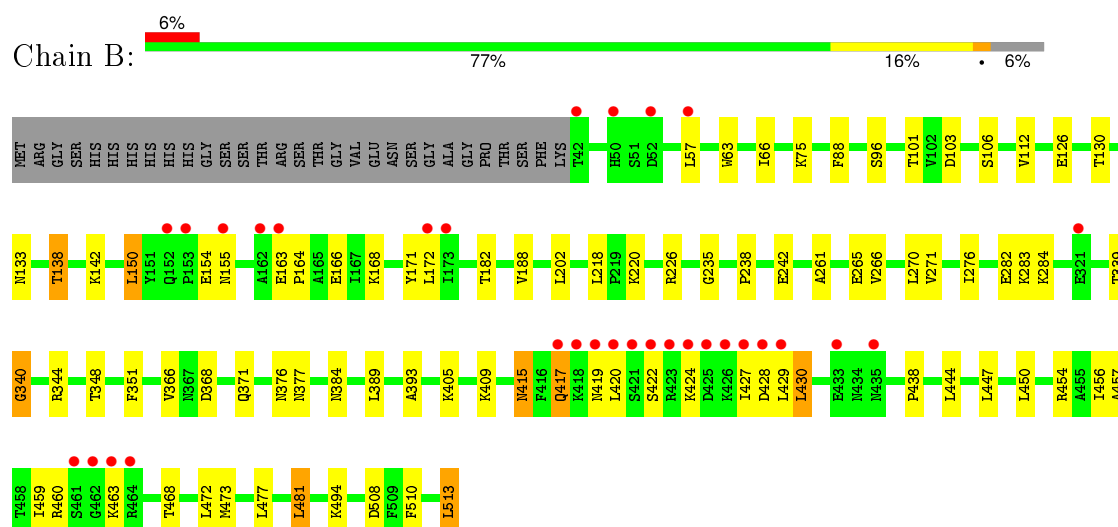
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Rotenone-insensitive NADH-ubiquinone oxidoreductase, mitochondrial



- Molecule 1: Rotenone-insensitive NADH-ubiquinone oxidoreductase, mitochondrial





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	127.41Å 230.19Å 112.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.36 – 2.52 29.62 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.0 (38.36-2.52) 99.7 (29.62-2.45)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.11 (at 2.45Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, $R_{free}$	0.189 , 0.217 0.189 , 0.188	Depositor DCC
$R_{free}$ test set	2855 reflections (5.36%)	DCC
Wilson B-factor (Å <sup>2</sup> )	36.9	Xtriage
Anisotropy	0.219	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 48.2	EDS
Estimated twinning fraction	0.018 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.027 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	1 of 60968 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	8170	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.22% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, NAI, UQ5, FAD

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	0.55	0/3824	0.73	9/5177 (0.2%)
1	B	0.56	0/3824	0.80	9/5177 (0.2%)
All	All	0.56	0/7648	0.76	18/10354 (0.2%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	454	ARG	NE-CZ-NH2	-17.13	111.73	120.30
1	B	454	ARG	NE-CZ-NH1	15.86	128.23	120.30
1	B	344	ARG	NE-CZ-NH2	-14.22	113.19	120.30
1	B	344	ARG	NE-CZ-NH1	13.07	126.83	120.30
1	A	226	ARG	NE-CZ-NH2	11.74	126.17	120.30
1	A	226	ARG	NE-CZ-NH1	-11.56	114.52	120.30
1	B	226	ARG	NE-CZ-NH1	11.46	126.03	120.30
1	A	454	ARG	NE-CZ-NH1	-11.45	114.58	120.30
1	B	226	ARG	NE-CZ-NH2	-11.44	114.58	120.30
1	A	344	ARG	NE-CZ-NH1	-11.14	114.73	120.30
1	A	454	ARG	NE-CZ-NH2	11.01	125.80	120.30
1	A	344	ARG	NE-CZ-NH2	10.32	125.46	120.30
1	B	454	ARG	CD-NE-CZ	7.53	134.14	123.60
1	B	344	ARG	CD-NE-CZ	6.47	132.66	123.60
1	B	226	ARG	CD-NE-CZ	6.18	132.26	123.60
1	A	226	ARG	CD-NE-CZ	6.04	132.05	123.60
1	A	454	ARG	CD-NE-CZ	5.72	131.60	123.60
1	A	344	ARG	CD-NE-CZ	5.27	130.98	123.60

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3738	0	3803	64	0
1	B	3738	0	3803	61	0
2	A	54	0	62	27	0
2	B	48	0	56	23	0
3	A	5	0	0	0	0
3	B	6	0	0	0	0
4	A	53	0	31	5	0
4	B	53	0	31	4	0
5	A	44	0	27	4	0
5	B	44	0	27	6	0
6	A	198	0	0	9	0
6	B	189	0	0	15	0
All	All	8170	0	7840	150	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 10.

All (150) 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)
2:B:605:UQ5:C1M	2:B:605:UQ5:H112	1.41	1.47
2:B:605:UQ5:C11	2:B:605:UQ5:H1M1	1.52	1.37
1:A:393:ALA:CB	2:A:604:UQ5:H4M1	1.65	1.25
1:A:393:ALA:CB	2:A:604:UQ5:C4M	2.24	1.15
1:A:393:ALA:HB1	2:A:604:UQ5:H4M1	1.25	1.11
1:A:393:ALA:HB3	2:A:604:UQ5:C4M	1.84	1.07
2:A:604:UQ5:H121	2:A:604:UQ5:C1M	1.87	1.04
1:A:393:ALA:HB1	2:A:604:UQ5:C4M	1.91	0.97
2:A:604:UQ5:H1M1	2:A:604:UQ5:H121	1.46	0.97
1:A:214:LYS:NZ	6:A:795:HOH:O	2.01	0.92
1:B:368:ASP:HB3	1:B:438:PRO:HB3	1.57	0.87
1:A:368:ASP:HB3	1:A:438:PRO:HB3	1.57	0.86
2:B:605:UQ5:H1M1	2:B:605:UQ5:C8	2.05	0.84
2:B:605:UQ5:C1M	2:B:605:UQ5:C11	2.30	0.84
2:B:605:UQ5:H1M1	2:B:605:UQ5:C9	2.09	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:238:PRO:HB2	5:B:603:NAI:H42N	1.59	0.83
1:A:393:ALA:HB3	2:A:604:UQ5:H4M2	1.60	0.83
2:A:604:UQ5:C8	2:A:604:UQ5:H1M1	2.09	0.82
2:B:605:UQ5:H112	2:B:605:UQ5:H1M1	0.84	0.82
1:A:238:PRO:HB2	5:A:610:NAI:H42N	1.60	0.81
1:A:393:ALA:CB	2:A:604:UQ5:H4M2	2.09	0.81
1:B:459:ILE:HD12	2:B:605:UQ5:H101	1.64	0.80
1:A:78:ASN:ND2	6:A:845:HOH:O	2.07	0.79
1:A:349:ASP:OD2	6:A:805:HOH:O	1.99	0.79
2:A:604:UQ5:H1M1	2:A:604:UQ5:C12	2.15	0.77
1:A:463:LYS:NZ	6:A:893:HOH:O	2.18	0.76
1:B:393:ALA:CB	2:B:605:UQ5:H4M1	2.16	0.75
2:B:605:UQ5:H1M2	2:B:605:UQ5:H112	1.65	0.75
1:B:377:ASN:OD1	6:B:702:HOH:O	2.05	0.75
5:B:603:NAI:H4N	6:B:889:HOH:O	1.87	0.75
1:B:138:THR:HG21	1:B:166:GLU:OE1	1.87	0.74
1:A:485:MET:HE1	2:A:604:UQ5:H72	1.69	0.74
1:B:510:PHE:O	6:B:859:HOH:O	2.05	0.74
1:B:366:VAL:O	6:B:810:HOH:O	2.07	0.73
1:B:271:VAL:HG12	5:B:603:NAI:H2A	1.71	0.72
1:B:282:GLU:OE1	6:B:718:HOH:O	2.07	0.72
2:B:605:UQ5:O2	2:B:605:UQ5:H3M2	1.89	0.71
1:A:138:THR:HG21	1:A:166:GLU:OE1	1.90	0.71
2:B:605:UQ5:C4M	2:B:605:UQ5:O3	2.38	0.71
1:B:513:LEU:O	6:B:859:HOH:O	2.08	0.70
1:B:265:GLU:OE2	6:B:767:HOH:O	2.09	0.70
2:A:604:UQ5:H3M2	2:A:604:UQ5:O2	1.92	0.69
1:A:393:ALA:HB3	2:A:604:UQ5:H4M1	1.47	0.67
1:A:151:TYR:OH	6:A:825:HOH:O	2.12	0.66
1:B:126:GLU:OE1	6:B:728:HOH:O	2.13	0.66
1:A:485:MET:CE	2:A:604:UQ5:H72	2.26	0.65
2:A:604:UQ5:H1M1	2:A:604:UQ5:C11	2.26	0.65
1:B:282:GLU:OE2	6:B:779:HOH:O	2.14	0.64
1:A:485:MET:SD	2:A:604:UQ5:C6	2.85	0.64
1:B:133:ASN:HB2	1:B:138:THR:HG22	1.78	0.64
1:B:238:PRO:HB2	5:B:603:NAI:C4N	2.28	0.63
1:B:393:ALA:HB1	2:B:605:UQ5:H4M1	1.80	0.63
1:A:133:ASN:HB2	1:A:138:THR:HG22	1.82	0.61
1:A:371:GLN:HE21	1:A:376:ASN:HA	1.65	0.61
1:A:238:PRO:HB2	5:A:610:NAI:C4N	2.30	0.61
1:B:393:ALA:HB3	2:B:605:UQ5:H4M1	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:605:UQ5:O2	2:B:605:UQ5:C3M	2.50	0.60
1:A:447:LEU:CD1	2:A:604:UQ5:H103	2.31	0.59
1:B:130:THR:HG21	1:B:142:LYS:HG3	1.83	0.59
1:B:371:GLN:HE21	1:B:376:ASN:HA	1.68	0.58
1:B:282:GLU:CD	6:B:779:HOH:O	2.42	0.58
1:A:405:LYS:HE2	1:A:409:LYS:HE3	1.85	0.58
1:A:130:THR:HG21	1:A:142:LYS:HG3	1.85	0.57
1:B:419:ASN:ND2	1:B:424:LYS:O	2.28	0.57
1:A:419:ASN:ND2	1:A:424:LYS:O	2.28	0.57
1:B:508:ASP:OD1	6:B:788:HOH:O	2.16	0.57
1:B:63:TRP:HB2	4:B:606:FAD:H5'2	1.85	0.56
1:B:393:ALA:CB	2:B:605:UQ5:C4M	2.84	0.56
1:B:393:ALA:HB3	2:B:605:UQ5:C4M	2.36	0.55
1:A:482:TYR:OH	5:A:610:NAI:N7N	2.39	0.55
1:A:393:ALA:C	2:A:604:UQ5:H4M2	2.27	0.55
2:B:605:UQ5:H4M2	2:B:605:UQ5:O3	2.06	0.54
1:A:485:MET:SD	2:A:604:UQ5:C5	2.97	0.53
1:A:349:ASP:OD1	6:A:863:HOH:O	2.19	0.53
1:B:405:LYS:HE2	1:B:409:LYS:HE3	1.91	0.53
1:A:447:LEU:HD12	2:A:604:UQ5:H103	1.91	0.52
2:A:604:UQ5:C8	2:A:604:UQ5:C1M	2.86	0.52
1:B:508:ASP:OD2	6:B:821:HOH:O	2.19	0.52
4:B:606:FAD:PA	4:B:606:FAD:O4'	2.68	0.52
2:A:604:UQ5:C3M	2:A:604:UQ5:O2	2.57	0.51
1:A:415:ASN:OD1	1:A:415:ASN:N	2.43	0.51
1:A:392:THR:HB	4:A:605:FAD:O2	2.10	0.50
1:B:182:THR:HG22	1:B:188:VAL:HG22	1.93	0.50
1:B:419:ASN:HA	1:B:422:SER:HB3	1.93	0.50
1:B:284:LYS:NZ	6:B:779:HOH:O	2.25	0.50
1:A:412:GLN:NE2	6:A:775:HOH:O	2.44	0.50
1:B:415:ASN:OD1	1:B:415:ASN:N	2.44	0.50
2:B:605:UQ5:H4M3	2:B:605:UQ5:O3	2.12	0.50
1:B:477:LEU:HG	1:B:481:LEU:HD22	1.95	0.49
4:B:606:FAD:O2A	4:B:606:FAD:O4'	2.30	0.49
1:A:220:LYS:HE3	1:A:261:ALA:HB1	1.95	0.49
1:B:150:LEU:O	6:B:861:HOH:O	2.19	0.49
1:B:457:ALA:O	1:B:468:THR:HA	2.12	0.49
2:A:604:UQ5:H4M3	2:A:604:UQ5:O3	2.12	0.48
1:B:419:ASN:OD1	1:B:427:ILE:HG13	2.13	0.48
1:B:66:ILE:HG13	1:B:88:PHE:HB2	1.95	0.48
1:A:447:LEU:HD11	2:A:604:UQ5:H103	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:477:LEU:HG	1:A:481:LEU:HD22	1.95	0.48
2:A:604:UQ5:H121	2:A:604:UQ5:H1M2	1.87	0.48
1:A:419:ASN:OD1	1:A:427:ILE:HG13	2.14	0.48
1:A:419:ASN:HA	1:A:422:SER:HB3	1.95	0.47
1:A:96:SER:HB3	1:A:101:THR:HB	1.96	0.47
1:A:62:GLY:HA3	4:A:605:FAD:O5B	2.15	0.47
1:A:426:LYS:HB3	1:A:426:LYS:HE2	1.77	0.47
2:B:605:UQ5:C1M	2:B:605:UQ5:C8	2.83	0.47
1:B:63:TRP:CH2	2:B:605:UQ5:H4M3	2.50	0.47
1:B:447:LEU:HD11	2:B:605:UQ5:H103	1.97	0.46
1:B:459:ILE:CD1	2:B:605:UQ5:H101	2.41	0.46
1:B:96:SER:HB3	1:B:101:THR:HB	1.97	0.46
1:A:75:LYS:HE3	1:A:75:LYS:HB2	1.79	0.45
1:A:450:LEU:HD11	1:A:456:ILE:HG23	1.99	0.45
1:B:417:GLN:HA	1:B:420:LEU:HD22	1.99	0.45
1:B:220:LYS:HE3	1:B:261:ALA:HB1	1.98	0.45
1:B:63:TRP:CD1	4:B:606:FAD:H4'	2.52	0.45
1:B:450:LEU:HD11	1:B:456:ILE:HG23	1.99	0.45
1:B:348:THR:O	1:B:351:PHE:HB2	2.17	0.44
1:B:460:ARG:HD3	6:B:840:HOH:O	2.16	0.44
1:A:95:PRO:HD3	4:A:605:FAD:HM72	2.00	0.44
1:B:57:LEU:O	1:B:172:LEU:HD12	2.17	0.44
1:A:444:LEU:HB3	2:A:604:UQ5:H122	1.99	0.44
1:B:171:TYR:CZ	1:B:420:LEU:HD11	2.53	0.43
1:A:179:GLU:OE1	6:A:771:HOH:O	2.21	0.43
1:B:238:PRO:O	1:B:242:GLU:HG3	2.18	0.43
1:B:481:LEU:HA	1:B:481:LEU:HD12	1.85	0.43
1:A:66:ILE:HG13	1:A:88:PHE:HB2	2.00	0.43
1:B:513:LEU:HA	1:B:513:LEU:HD13	1.71	0.43
1:A:405:LYS:O	1:A:408:ASP:HB2	2.19	0.42
1:A:481:LEU:HA	1:A:481:LEU:HD12	1.80	0.42
1:A:239:THR:OG1	5:A:610:NAI:H5N	2.19	0.42
1:A:276:ILE:H	1:A:276:ILE:HD13	1.83	0.42
1:A:417:GLN:HA	1:A:420:LEU:HD22	2.02	0.42
1:A:194:PHE:O	1:A:200:ASN:HB3	2.19	0.42
1:A:511:LYS:NZ	6:A:866:HOH:O	1.96	0.42
1:A:238:PRO:O	1:A:242:GLU:HG3	2.21	0.41
1:A:105:LYS:NZ	1:B:103:ASP:OD1	2.44	0.41
1:A:133:ASN:OD1	1:A:138:THR:HG22	2.21	0.41
1:A:94:LEU:HB3	4:A:605:FAD:HM72	2.02	0.41
1:B:163:GLU:HA	1:B:164:PRO:HD3	1.90	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:340:GLY:HA2	5:B:603:NAI:O2N	2.20	0.41
4:A:605:FAD:H1'1	4:A:605:FAD:H9	1.84	0.41
1:B:75:LYS:HB2	1:B:75:LYS:HE3	1.78	0.41
1:A:513:LEU:HA	1:A:513:LEU:HD13	1.71	0.41
1:B:430:LEU:HA	1:B:430:LEU:HD12	1.91	0.40
1:B:235:GLY:HA2	5:B:603:NAI:O4B	2.20	0.40
1:A:182:THR:HG22	1:A:188:VAL:HG22	2.02	0.40
1:B:393:ALA:HB1	2:B:605:UQ5:C4M	2.46	0.40
1:A:430:LEU:HD12	1:A:430:LEU:HA	1.92	0.40
1:B:106:SER:HA	1:B:494:LYS:HD3	2.02	0.40
1:A:494:LYS:HE3	1:B:508:ASP:OD1	2.21	0.40
1:A:457:ALA:O	1:A:468:THR:HA	2.22	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	470/502 (94%)	453 (96%)	14 (3%)	3 (1%)	30	49
1	B	470/502 (94%)	454 (97%)	14 (3%)	2 (0%)	39	60
All	All	940/1004 (94%)	907 (96%)	28 (3%)	5 (0%)	34	55

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	340	GLY
1	B	340	GLY
1	A	154	GLU
1	A	297	SER
1	B	154	GLU

### 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	403/427 (94%)	378 (94%)	25 (6%)	23	40
1	B	403/427 (94%)	378 (94%)	25 (6%)	23	40
All	All	806/854 (94%)	756 (94%)	50 (6%)	23	40

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	112	VAL
1	A	138	THR
1	A	150	LEU
1	A	155	ASN
1	A	168	LYS
1	A	202	LEU
1	A	218	LEU
1	A	266	VAL
1	A	270	LEU
1	A	276	ILE
1	A	283	LYS
1	A	339	THR
1	A	384	ASN
1	A	389	LEU
1	A	415	ASN
1	A	417	GLN
1	A	428	ASP
1	A	429	LEU
1	A	430	LEU
1	A	444	LEU
1	A	463	LYS
1	A	472	LEU
1	A	473	MET
1	A	481	LEU
1	A	513	LEU
1	B	112	VAL
1	B	138	THR

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Mol	Chain	Res	Type
1	B	150	LEU
1	B	155	ASN
1	B	168	LYS
1	B	202	LEU
1	B	218	LEU
1	B	266	VAL
1	B	270	LEU
1	B	276	ILE
1	B	283	LYS
1	B	339	THR
1	B	384	ASN
1	B	389	LEU
1	B	415	ASN
1	B	417	GLN
1	B	428	ASP
1	B	429	LEU
1	B	430	LEU
1	B	444	LEU
1	B	463	LYS
1	B	472	LEU
1	B	473	MET
1	B	481	LEU
1	B	513	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	377	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

Of 20 ligands modelled in this entry, 11 are monoatomic - leaving 9 for Mogul analysis.

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	UQ5	A	601	-	19,19,38	4.96	6 (31%)	22,22,49	3.63	18 (81%)
2	UQ5	A	602	-	9,9,38	4.09	2 (22%)	10,10,49	3.46	10 (100%)
2	UQ5	A	604	-	24,24,38	3.72	9 (37%)	29,32,49	2.88	10 (34%)
4	FAD	A	605	3	48,58,58	1.30	5 (10%)	54,89,89	2.13	10 (18%)
5	NAI	A	610	-	38,48,48	1.76	7 (18%)	48,73,73	2.05	6 (12%)
5	NAI	B	603	-	38,48,48	1.88	10 (26%)	48,73,73	2.35	14 (29%)
2	UQ5	B	604	-	19,19,38	4.89	4 (21%)	22,22,49	3.46	16 (72%)
2	UQ5	B	605	-	28,28,38	3.74	10 (35%)	34,37,49	2.75	13 (38%)
4	FAD	B	606	3	48,58,58	1.31	5 (10%)	54,89,89	2.13	10 (18%)

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	UQ5	A	601	-	-	0/20/20/57	0/0/0/1
2	UQ5	A	602	-	-	1/8/8/57	0/0/0/1
2	UQ5	A	604	-	-	0/17/41/57	0/1/1/1
4	FAD	A	605	3	-	0/30/50/50	0/6/6/6
5	NAI	A	610	-	-	0/25/72/72	0/5/5/5
5	NAI	B	603	-	-	0/25/72/72	0/5/5/5
2	UQ5	B	604	-	-	0/20/20/57	0/0/0/1
2	UQ5	B	605	-	-	0/21/45/57	0/1/1/1
4	FAD	B	606	3	-	0/30/50/50	0/6/6/6

All (58) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	603	NAI	C4N-C5N	-3.37	1.41	1.49
5	A	610	NAI	C4N-C5N	-3.26	1.42	1.49
2	B	605	UQ5	C3-C2	-2.95	1.40	1.48
2	A	604	UQ5	C3-C2	-2.93	1.40	1.48
2	B	605	UQ5	C4-C5	-2.69	1.41	1.48
2	A	604	UQ5	C4-C5	-2.65	1.41	1.48
2	B	605	UQ5	O4-C4M	-2.61	1.39	1.45
2	A	604	UQ5	O4-C4M	-2.59	1.39	1.45
5	B	603	NAI	PA-O5B	-2.32	1.48	1.59
2	B	605	UQ5	C6-C5	-2.09	1.40	1.46
5	B	603	NAI	O2B-C2B	-2.07	1.38	1.43
2	A	604	UQ5	C6-C5	-2.04	1.40	1.46
2	A	601	UQ5	C16-C14	2.03	1.55	1.51
2	A	601	UQ5	C12-C13	2.10	1.56	1.50
5	A	610	NAI	C2A-N1A	2.17	1.38	1.33
5	B	603	NAI	C5A-C4A	2.32	1.45	1.40
5	B	603	NAI	C2A-N3A	2.58	1.36	1.32
2	A	604	UQ5	C6-C1	2.68	1.41	1.35
2	B	605	UQ5	C6-C1	2.71	1.41	1.35
5	B	603	NAI	C2A-N1A	2.79	1.39	1.33
5	A	610	NAI	C2N-C3N	2.94	1.41	1.34
4	B	606	FAD	C5A-C4A	2.95	1.47	1.40
4	A	605	FAD	C5A-C4A	2.95	1.47	1.40
4	A	605	FAD	C8-C7	2.96	1.48	1.41
4	B	606	FAD	C8-C7	2.99	1.49	1.41
5	A	610	NAI	C6A-N6A	3.01	1.44	1.34
5	B	603	NAI	C2N-C3N	3.04	1.42	1.34
5	B	603	NAI	C6A-N6A	3.11	1.44	1.34
5	A	610	NAI	C2A-N3A	3.20	1.37	1.32
4	A	605	FAD	C9A-C5X	3.26	1.49	1.42
4	B	606	FAD	C9A-C5X	3.31	1.49	1.42
4	A	605	FAD	C4-C4X	3.57	1.48	1.41
4	B	606	FAD	C4-C4X	3.58	1.48	1.41
5	A	610	NAI	C6N-C5N	3.70	1.40	1.33
4	A	605	FAD	C4X-C10	3.86	1.48	1.41
4	B	606	FAD	C4X-C10	3.86	1.48	1.41
5	B	603	NAI	C6N-C5N	4.09	1.41	1.33
5	A	610	NAI	C7N-N7N	4.63	1.46	1.33
5	B	603	NAI	C7N-N7N	4.90	1.47	1.33
2	A	604	UQ5	O2-C2	6.26	1.37	1.23
2	B	605	UQ5	O2-C2	6.31	1.37	1.23
2	B	605	UQ5	O5-C5	6.33	1.37	1.23
2	A	604	UQ5	O5-C5	6.37	1.38	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	604	UQ5	C18-C19	7.20	1.54	1.32
2	B	605	UQ5	C18-C19	7.32	1.54	1.32
2	A	601	UQ5	C18-C19	7.35	1.54	1.32
2	A	602	UQ5	C13-C14	7.35	1.54	1.32
2	A	602	UQ5	C8-C9	9.41	1.53	1.32
2	B	605	UQ5	C8-C9	9.73	1.52	1.33
2	A	604	UQ5	C8-C9	9.76	1.52	1.33
2	A	604	UQ5	C13-C14	10.48	1.53	1.33
2	B	605	UQ5	C13-C14	10.51	1.53	1.33
2	A	601	UQ5	C8-C9	11.00	1.54	1.33
2	B	604	UQ5	C6-C1	11.18	1.54	1.33
2	B	604	UQ5	C8-C9	11.21	1.55	1.33
2	A	601	UQ5	C6-C1	11.33	1.55	1.33
2	B	604	UQ5	C13-C14	11.47	1.55	1.33
2	A	601	UQ5	C13-C14	11.89	1.56	1.33

All (107) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	603	NAI	N3A-C2A-N1A	-12.33	119.45	128.89
5	A	610	NAI	N3A-C2A-N1A	-10.74	120.67	128.89
2	A	604	UQ5	C10-C9-C8	-7.18	109.41	123.50
2	B	605	UQ5	C10-C9-C8	-7.16	109.45	123.50
2	A	602	UQ5	C10-C9-C8	-7.05	110.20	123.83
4	A	605	FAD	N3A-C2A-N1A	-7.03	123.51	128.89
4	B	606	FAD	N3A-C2A-N1A	-6.96	123.57	128.89
2	B	604	UQ5	C1M-C1-C6	-6.92	109.92	123.50
2	A	601	UQ5	C10-C9-C8	-6.80	110.14	123.50
2	A	601	UQ5	C1M-C1-C6	-6.34	111.05	123.50
2	A	604	UQ5	C15-C14-C13	-6.27	111.19	123.50
2	B	605	UQ5	C15-C14-C13	-6.24	111.26	123.50
2	A	604	UQ5	C15-C14-C16	-5.85	109.06	115.68
4	A	605	FAD	C4-C4X-C10	-5.76	116.26	119.94
2	B	604	UQ5	C10-C9-C8	-5.64	112.43	123.50
2	A	601	UQ5	C15-C14-C13	-5.53	112.65	123.50
4	B	606	FAD	C4-C4X-C10	-5.43	116.46	119.94
2	A	601	UQ5	C1M-C1-C2	-5.30	109.68	115.68
2	A	604	UQ5	C7-C8-C9	-5.28	117.75	126.70
2	B	605	UQ5	C7-C8-C9	-5.26	117.78	126.70
2	B	604	UQ5	C15-C14-C13	-4.96	113.76	123.50
2	B	604	UQ5	C10-C9-C11	-4.78	108.11	115.41
2	B	605	UQ5	C12-C13-C14	-4.62	117.71	127.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	604	UQ5	C12-C13-C14	-4.61	117.73	127.76
2	A	601	UQ5	C10-C9-C11	-4.45	108.61	115.41
2	B	605	UQ5	C16-C14-C13	-4.37	112.77	121.05
2	B	604	UQ5	C11-C9-C8	-4.21	113.06	121.05
2	B	605	UQ5	C15-C14-C16	-4.19	109.00	115.41
2	A	604	UQ5	C11-C9-C8	-4.14	113.20	121.05
2	B	604	UQ5	C1M-C1-C2	-4.07	111.07	115.68
4	B	606	FAD	C4X-C4-N3	-3.95	118.19	123.59
2	B	605	UQ5	C11-C9-C8	-3.93	113.59	121.05
2	B	604	UQ5	C16-C14-C13	-3.88	113.69	121.05
4	A	605	FAD	C4X-C4-N3	-3.87	118.30	123.59
4	B	606	FAD	P-O3P-PA	-3.77	122.14	132.73
2	A	602	UQ5	C10-C9-C11	-3.73	109.71	115.41
2	A	601	UQ5	C11-C9-C8	-3.71	114.01	121.05
4	A	605	FAD	P-O3P-PA	-3.67	122.42	132.73
2	A	604	UQ5	C16-C14-C13	-3.61	112.69	120.74
2	A	604	UQ5	C10-C9-C11	-3.46	110.13	115.41
2	B	604	UQ5	C20-C19-C18	-3.44	111.54	122.61
2	B	605	UQ5	C10-C9-C11	-3.44	110.16	115.41
2	B	604	UQ5	C12-C13-C14	-3.39	120.39	127.76
2	A	601	UQ5	C20-C19-C18	-3.31	111.96	122.61
4	A	605	FAD	C4A-C5A-N7A	-3.26	106.48	109.48
5	A	610	NAI	C3N-C2N-N1N	-3.24	118.50	123.14
4	B	606	FAD	C4A-C5A-N7A	-3.23	106.51	109.48
5	B	603	NAI	PN-O3-PA	-3.19	123.78	132.73
2	A	601	UQ5	C12-C13-C14	-3.06	121.11	127.76
2	A	602	UQ5	C15-C14-C13	-3.04	112.84	122.61
2	A	601	UQ5	C2-C1-C6	-3.02	114.00	120.74
2	A	601	UQ5	C16-C14-C13	-3.01	115.34	121.05
2	B	604	UQ5	C21-C19-C18	-2.98	113.03	122.61
2	B	604	UQ5	C7-C6-C1	-2.95	117.94	126.80
2	A	602	UQ5	C16-C14-C13	-2.87	113.38	122.61
2	A	602	UQ5	C11-C9-C8	-2.77	113.42	120.12
2	A	601	UQ5	C21-C19-C18	-2.77	113.69	122.61
2	B	605	UQ5	C20-C19-C18	-2.73	113.82	122.61
2	B	605	UQ5	C21-C19-C18	-2.73	113.83	122.61
2	B	604	UQ5	C17-C18-C19	-2.71	117.30	127.73
2	B	605	UQ5	C17-C18-C19	-2.59	117.75	127.73
5	A	610	NAI	C4A-C5A-N7A	-2.55	107.13	109.48
2	A	601	UQ5	C7-C6-C1	-2.52	119.24	126.80
2	A	602	UQ5	C12-C13-C14	-2.50	118.11	127.73
5	B	603	NAI	O2A-PA-O5B	-2.49	95.92	108.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	606	FAD	C4X-C10-N10	-2.42	119.10	120.52
2	A	602	UQ5	C7-C8-C9	-2.30	120.68	126.57
4	A	605	FAD	C4X-C10-N10	-2.28	119.17	120.52
5	A	610	NAI	PN-O3-PA	-2.27	126.36	132.73
2	A	601	UQ5	C7-C8-C9	-2.25	120.03	126.80
5	B	603	NAI	C3N-C2N-N1N	-2.23	119.94	123.14
5	B	603	NAI	O2A-PA-O3	-2.20	95.12	105.09
5	B	603	NAI	C4A-C5A-N7A	-2.16	107.49	109.48
5	B	603	NAI	C4D-O4D-C1D	-2.14	104.80	109.52
2	A	602	UQ5	C16-C14-C15	-2.10	109.48	114.64
2	B	604	UQ5	C2-C1-C6	-2.07	116.12	120.74
2	A	601	UQ5	C21-C19-C20	-2.07	109.56	114.64
5	B	603	NAI	O4B-C4B-C3B	2.04	109.26	105.15
2	A	604	UQ5	C8-C7-C6	2.05	117.80	111.64
5	B	603	NAI	O3-PN-O5D	2.05	108.38	102.94
5	B	603	NAI	O4B-C1B-N9A	2.17	112.65	108.10
2	B	604	UQ5	C11-C12-C13	2.18	117.40	111.69
5	A	610	NAI	O4B-C4B-C3B	2.19	109.56	105.15
2	A	602	UQ5	C12-C11-C9	2.21	119.92	112.71
2	B	605	UQ5	C16-C17-C18	2.31	117.75	111.69
2	A	604	UQ5	C11-C12-C13	2.35	117.85	111.69
5	B	603	NAI	O5B-PA-O1A	2.36	118.76	109.62
5	B	603	NAI	C2A-N1A-C6A	2.40	123.06	118.77
4	A	605	FAD	C5X-C9A-N10	2.41	119.45	117.62
2	B	605	UQ5	C11-C12-C13	2.43	118.06	111.69
2	A	601	UQ5	C12-C11-C9	2.45	120.69	112.71
4	B	606	FAD	C5X-C9A-N10	2.51	119.53	117.62
2	B	604	UQ5	C8-C7-C6	2.53	119.79	112.22
2	B	604	UQ5	C12-C11-C9	2.55	121.03	112.71
4	B	606	FAD	C4-C4X-N5	2.63	121.91	118.72
5	B	603	NAI	C5N-C4N-C3N	2.67	119.88	112.52
2	A	601	UQ5	C8-C7-C6	2.68	120.23	112.22
5	A	610	NAI	C5N-C4N-C3N	2.89	120.48	112.52
4	A	605	FAD	C4-C4X-N5	2.93	122.27	118.72
5	B	603	NAI	O4D-C1D-N1N	3.00	114.40	108.07
2	A	602	UQ5	C11-C12-C13	3.17	119.98	111.69
2	A	601	UQ5	C16-C17-C18	3.18	120.01	111.69
2	A	601	UQ5	C11-C12-C13	3.71	121.40	111.69
4	A	605	FAD	C4X-N5-C5X	3.71	121.03	116.76
4	B	606	FAD	C4X-N5-C5X	3.72	121.04	116.76
4	B	606	FAD	C4-N3-C2	7.94	122.11	115.25
4	A	605	FAD	C4-N3-C2	7.94	122.11	115.25

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	602	UQ5	C7-C8-C9-C10

There are no ring outliers.

6 monomers are involved in 69 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	604	UQ5	27	0
4	A	605	FAD	5	0
5	A	610	NAI	4	0
5	B	603	NAI	6	0
2	B	605	UQ5	23	0
4	B	606	FAD	4	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	472/502 (94%)	-0.10	20 (4%)	40	45	23, 37, 80, 117	0
1	B	472/502 (94%)	-0.01	31 (6%)	22	24	23, 38, 81, 119	0
All	All	944/1004 (94%)	-0.05	51 (5%)	29	34	23, 37, 81, 119	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	422	SER	6.8
1	A	152	GLN	6.8
1	B	423	ARG	6.6
1	B	152	GLN	6.3
1	A	421	SER	6.1
1	B	421	SER	5.4
1	B	427	ILE	5.0
1	A	50	HIS	5.0
1	B	425	ASP	4.9
1	B	420	LEU	4.5
1	B	429	LEU	4.3
1	B	424	LYS	4.2
1	A	423	ARG	4.2
1	A	324	LYS	4.0
1	A	422	SER	3.9
1	B	50	HIS	3.8
1	B	418	LYS	3.5
1	B	426	LYS	3.3
1	A	325	ILE	3.3
1	A	463	LYS	3.2
1	B	321	GLU	3.1
1	B	52	ASP	3.1
1	A	322	ASP	3.1
1	B	42	THR	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	42	THR	3.0
1	A	323	GLY	2.9
1	B	153	PRO	2.8
1	B	155	ASN	2.8
1	B	428	ASP	2.8
1	B	463	LYS	2.7
1	B	163	GLU	2.7
1	B	162	ALA	2.7
1	A	424	LYS	2.7
1	A	420	LEU	2.6
1	B	417	GLN	2.6
1	A	155	ASN	2.6
1	A	464	ARG	2.6
1	A	282	GLU	2.5
1	B	461	SER	2.5
1	B	173	ILE	2.4
1	B	462	GLY	2.4
1	B	464	ARG	2.4
1	A	418	LYS	2.3
1	A	427	ILE	2.3
1	B	57	LEU	2.2
1	B	419	ASN	2.2
1	B	433	GLU	2.1
1	B	435	ASN	2.1
1	B	172	LEU	2.1
1	A	361	LYS	2.1
1	A	319	LYS	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	UQ5	A	601	20/38	0.84	0.34	13.67	51,65,79,81	0
2	UQ5	A	604	24/38	0.69	0.43	7.49	49,67,84,87	0
3	MG	A	609	1/1	0.92	0.18	5.40	51,51,51,51	0
2	UQ5	B	605	28/38	0.77	0.39	4.12	51,62,84,88	0
5	NAI	B	603	44/44	0.84	0.25	3.54	30,63,102,126	0
3	MG	B	610	1/1	0.88	0.22	2.64	51,51,51,51	0
3	MG	B	608	1/1	0.94	0.27	2.36	45,45,45,45	0
3	MG	A	603	1/1	0.74	0.25	1.08	70,70,70,70	0
5	NAI	A	610	44/44	0.86	0.17	0.98	21,65,85,102	0
4	FAD	B	606	53/53	0.96	0.17	0.50	20,30,37,40	0
4	FAD	A	605	53/53	0.97	0.14	0.11	15,28,34,37	0
3	MG	B	609	1/1	0.96	0.15	-0.51	46,46,46,46	0
3	MG	B	601	1/1	0.97	0.09	-0.87	23,23,23,23	0
3	MG	A	607	1/1	0.98	0.12	-1.29	42,42,42,42	0
3	MG	A	606	1/1	0.98	0.09	-1.44	38,38,38,38	0
3	MG	B	607	1/1	0.93	0.12	-1.51	42,42,42,42	0
3	MG	A	608	1/1	0.97	0.09	-1.81	41,41,41,41	0
2	UQ5	B	604	20/38	0.80	0.37	-	48,73,86,93	0
3	MG	B	602	1/1	0.88	0.09	-	60,60,60,60	0
2	UQ5	A	602	10/38	0.79	0.19	-	44,65,75,82	0

## 6.5 Other polymers

There are no such residues in this entry.