



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 12:57 PM GMT

PDB ID : 1T6Y  
Title : Crystal structure of ADP, AMP, and FMN bound TM379  
Authors : Shin, D.H.; Wang, W.; Kim, R.; Yokota, H.; Kim, S.-H.; Berkeley Structural Genomics Center (BSGC)  
Deposited on : 2004-05-07  
Resolution : 2.80 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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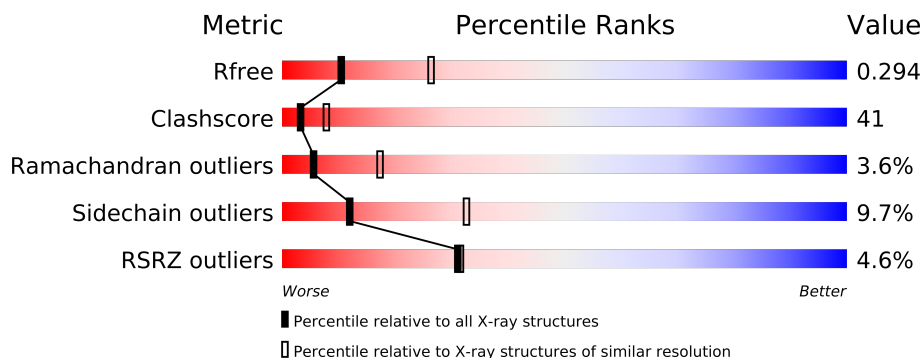
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.80 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))
$R_{free}$	66092	1799 (2.80-2.80)
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (2.80-2.80)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	293	
1	B	293	

## 2 Entry composition i

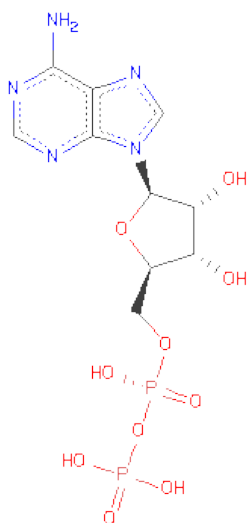
There are 5 unique types of molecules in this entry. The entry contains 4560 atoms, of which 0 are hydrogen and 0 are deuterium.

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 riboflavin kinase/FMN adenylyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	0	0	0
			2176	1400	378	392	6			
1	B	270	Total	C	N	O	S	0	0	0
			2189	1413	373	397	6			

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).



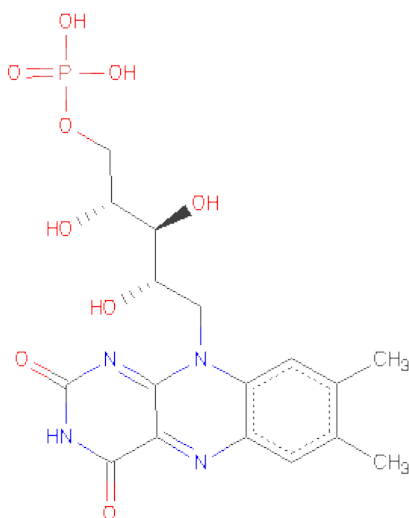
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 3 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula:  $C_{10}H_{14}N_5O_7P$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
3	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 4 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula:  $C_{17}H_{21}N_4O_9P$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
4	B	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 5 is water.

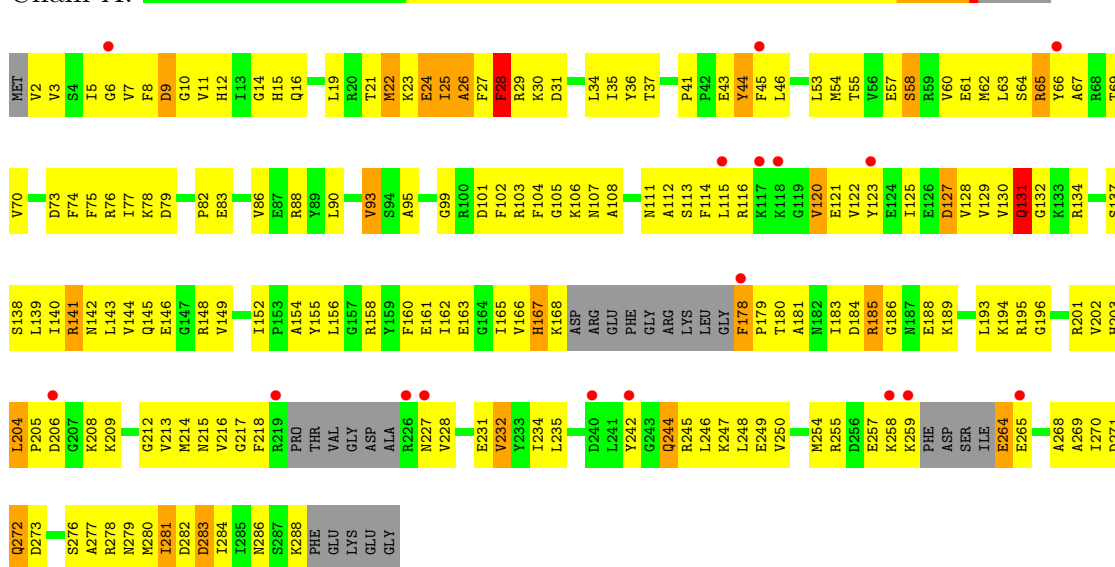
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	18	Total 18	O 18	0	0
5	B	15	Total 15	O 15	0	0

### 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.

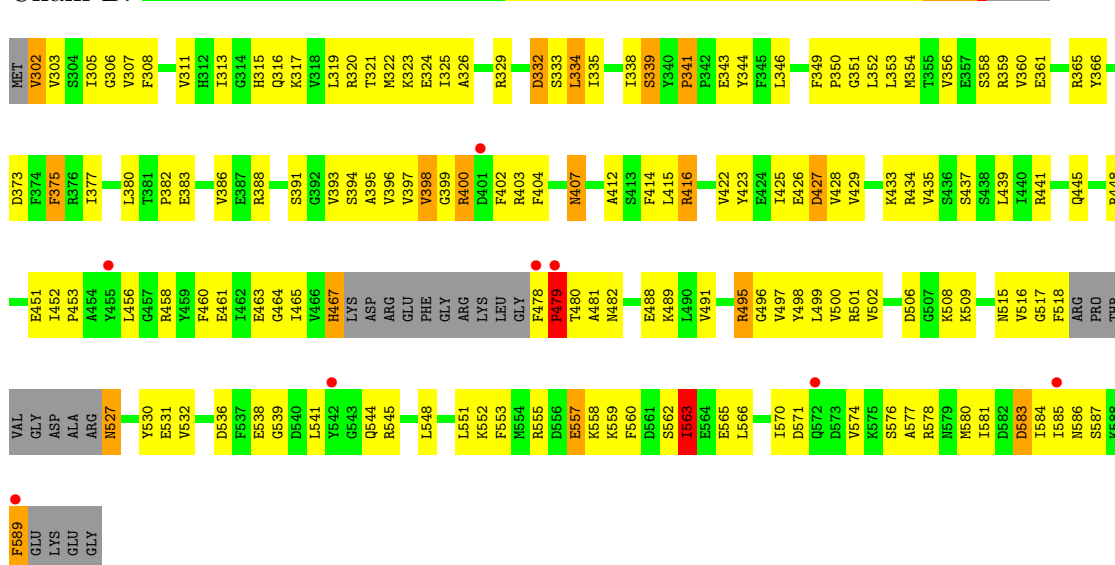
- Molecule 1: riboflavin kinase/FMN adenylyltransferase

Chain A:



- Molecule 1: riboflavin kinase/FMN adenylyltransferase

Chain B:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	67.80Å 82.66Å 66.72Å 90.00° 116.42° 90.00°	Depositor
Resolution (Å)	19.92 – 2.80 29.87 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.1 (19.92-2.80) 99.1 (29.87-2.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.85 (at 2.80Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.231 , 0.298 0.230 , 0.294	Depositor DCC
$R_{free}$ test set	1636 reflections (10.10%)	DCC
Wilson B-factor (Å <sup>2</sup> )	76.4	Xtriage
Anisotropy	0.231	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 49.2	EDS
Estimated twinning fraction	0.024 for l,-k,h	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 16237 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4560	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.25% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FMN, AMP, ADP

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	0/2213	0.81	1/2968 (0.0%)
1	B	0.52	0/2229	0.75	0/2994
All	All	0.53	0/4442	0.78	1/5962 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	9	ASP	CB-CG-OD2	5.18	122.96	118.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2176	0	2221	204	0
1	B	2189	0	2221	167	0
2	A	27	0	12	1	0
2	B	27	0	12	1	0
3	A	23	0	12	4	0
3	B	23	0	12	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	31	0	19	1	0
4	B	31	0	19	6	0
5	A	18	0	0	2	1
5	B	15	0	0	2	1
All	All	4560	0	4528	368	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 41.

All (368) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:214:MET:HB2	1:A:232:VAL:HG23	1.37	1.06
1:A:103:ARG:HD3	1:A:108:ALA:HB1	1.34	1.04
1:A:3:VAL:HG12	1:A:34:LEU:HB3	1.38	1.03
1:A:212:GLY:HA2	1:A:235:LEU:HD13	1.49	0.93
1:B:495:ARG:H	1:B:495:ARG:HH11	1.12	0.90
1:A:21:THR:HG21	1:A:125:ILE:HD11	1.53	0.89
1:A:270:ILE:HD12	1:A:271:ASP:N	1.88	0.88
1:A:280:MET:O	1:A:283:ASP:HB2	1.74	0.87
1:A:3:VAL:CG1	1:A:34:LEU:HB3	2.04	0.87
1:A:161:GLU:OE1	1:A:247:LYS:HE3	1.80	0.81
1:A:269:ALA:HA	1:A:272:GLN:NE2	1.95	0.81
1:A:166:VAL:HG12	1:A:167:HIS:H	1.45	0.81
1:A:140:ILE:O	1:A:144:VAL:HG23	1.80	0.81
1:A:255:ARG:HG3	1:A:255:ARG:HH11	1.46	0.80
1:B:388:ARG:HD3	5:B:629:HOH:O	1.81	0.80
1:A:166:VAL:HG23	1:A:244:GLN:O	1.81	0.80
1:B:321:THR:HG21	1:B:425:ILE:HD11	1.64	0.80
1:A:75:PHE:HA	1:A:78:LYS:HE3	1.65	0.79
1:B:465:ILE:HD11	1:B:545:ARG:HG2	1.64	0.79
1:A:88:ARG:HB2	1:A:88:ARG:HH11	1.47	0.79
1:A:158:ARG:HG2	1:A:158:ARG:HH11	1.47	0.78
1:A:41:PRO:HB2	1:A:43:GLU:OE2	1.84	0.77
1:B:517:GLY:O	1:B:518:PHE:HB2	1.84	0.77
1:B:463:GLU:HG2	1:B:464:GLY:N	2.00	0.77
1:A:203:HIS:ND1	1:A:209:LYS:HG2	2.00	0.76
1:A:9:ASP:OD1	1:A:141:ARG:NH1	2.18	0.76
1:B:382:PRO:O	1:B:386:VAL:HG23	1.85	0.75
1:A:178:PHE:N	1:A:179:PRO:HA	2.02	0.75
1:B:465:ILE:CD1	1:B:545:ARG:HG2	2.16	0.74
1:A:139:LEU:HD12	1:B:506:ASP:OD1	1.86	0.74
1:B:501:ARG:HB2	1:B:551:LEU:HD11	1.70	0.73

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:166:VAL:HB	1:A:242:TYR:O	1.89	0.73
1:A:129:VAL:HG12	1:A:134:ARG:HA	1.71	0.72
1:B:354:MET:HE1	1:B:359:ARG:HA	1.71	0.72
1:A:202:VAL:HG22	1:A:248:LEU:HD22	1.72	0.71
1:B:495:ARG:NH1	1:B:495:ARG:H	1.88	0.71
1:A:178:PHE:N	5:A:625:HOH:O	2.24	0.71
1:B:501:ARG:HH11	1:B:501:ARG:HG2	1.55	0.71
1:A:127:ASP:HA	3:A:295:AMP:C2	2.27	0.70
1:A:103:ARG:CD	1:A:108:ALA:HB1	2.16	0.70
1:A:29:ARG:C	1:A:30:LYS:HD2	2.12	0.70
1:A:202:VAL:HG22	1:A:248:LEU:CD2	2.23	0.69
1:B:570:ILE:O	1:B:574:VAL:HG13	1.93	0.69
1:B:577:ALA:O	1:B:581:ILE:HG13	1.92	0.69
1:B:315:HIS:O	1:B:319:LEU:HG	1.93	0.69
1:B:307:VAL:O	1:B:308:PHE:HB2	1.93	0.68
1:B:437:SER:O	1:B:441:ARG:HG3	1.93	0.68
1:B:495:ARG:N	1:B:495:ARG:HH11	1.87	0.68
1:B:407:ASN:N	1:B:407:ASN:HD22	1.92	0.68
1:A:12:HIS:O	1:A:15:HIS:HB2	1.92	0.68
1:B:356:VAL:O	1:B:360:VAL:HG23	1.94	0.68
1:A:255:ARG:HG3	1:A:255:ARG:NH1	2.08	0.67
1:A:166:VAL:HG12	1:A:167:HIS:N	2.08	0.67
1:A:8:PHE:HA	1:A:15:HIS:CE1	2.30	0.67
1:A:88:ARG:HB2	1:A:88:ARG:NH1	2.09	0.66
1:B:358:SER:O	1:B:361:GLU:HB3	1.95	0.66
1:A:167:HIS:O	1:A:168:LYS:HB3	1.96	0.65
1:A:83:GLU:CD	1:A:83:GLU:H	1.99	0.65
1:B:325:ILE:HD13	1:B:423:TYR:CD2	2.31	0.65
1:B:302:VAL:CG1	1:B:326:ALA:HB2	2.27	0.65
1:A:158:ARG:HG2	1:A:158:ARG:NH1	2.10	0.64
1:A:60:VAL:HG12	1:A:60:VAL:O	1.96	0.64
1:A:34:LEU:HD11	1:A:70:VAL:HG23	1.80	0.64
1:A:130:VAL:HG21	1:A:154:ALA:O	1.97	0.64
1:A:11:VAL:HA	1:A:15:HIS:ND1	2.12	0.64
1:A:178:PHE:CZ	1:A:270:ILE:HD13	2.33	0.64
1:B:316:GLN:HG2	1:B:366:TYR:OH	1.96	0.64
1:A:195:ARG:HD2	1:A:257:GLU:HG2	1.81	0.63
1:A:178:PHE:O	1:A:178:PHE:HD2	1.81	0.63
1:B:329:ARG:HH12	1:B:394:SER:HB2	1.63	0.63
1:B:383:GLU:H	1:B:383:GLU:CD	2.02	0.63
1:A:149:VAL:HG11	1:A:250:VAL:CG1	2.29	0.63
1:A:19:LEU:O	1:A:22:MET:HB3	1.99	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:481:ALA:O	1:B:531:GLU:HA	1.99	0.62
1:B:377:ILE:HA	1:B:380:LEU:HG	1.81	0.62
1:A:12:HIS:O	1:A:16:GLN:HG3	1.98	0.62
1:A:258:LYS:HG2	1:A:259:LYS:N	2.13	0.62
1:A:112:ALA:HB1	1:A:122:VAL:HG11	1.80	0.62
1:A:127:ASP:N	1:A:127:ASP:OD2	2.32	0.61
1:B:316:GLN:HB3	1:B:320:ARG:HH12	1.65	0.61
1:A:245:ARG:C	1:A:246:LEU:HD12	2.20	0.61
1:A:45:PHE:HE2	1:A:74:PHE:CD2	2.18	0.61
1:A:28:PHE:HD1	1:A:28:PHE:O	1.83	0.61
1:B:343:GLU:H	1:B:343:GLU:CD	2.04	0.61
1:A:45:PHE:O	1:A:46:LEU:HD23	2.00	0.61
1:A:21:THR:CG2	1:A:125:ILE:HD11	2.28	0.61
1:A:3:VAL:HG12	1:A:34:LEU:CB	2.23	0.60
1:B:352:LEU:O	1:B:491:VAL:HG12	2.00	0.60
1:A:178:PHE:HZ	1:A:270:ILE:HD13	1.67	0.60
1:B:354:MET:CE	1:B:359:ARG:HA	2.31	0.60
1:A:37:THR:HB	1:A:63:LEU:HD13	1.84	0.60
1:B:373:ASP:OD1	1:B:375:PHE:N	2.35	0.60
1:A:130:VAL:HB	1:A:155:TYR:CE1	2.36	0.60
1:B:580:MET:O	1:B:583:ASP:HB2	2.01	0.60
1:B:555:ARG:NH2	4:B:596:FMN:N5	2.50	0.59
1:A:269:ALA:HA	1:A:272:GLN:HE22	1.65	0.59
1:A:83:GLU:HG3	1:A:114:PHE:HZ	1.67	0.59
1:A:58:SER:O	1:A:61:GLU:HB3	2.02	0.59
1:B:552:LYS:HD3	1:B:580:MET:HE1	1.85	0.59
1:A:130:VAL:HG12	1:A:131:GLN:HG3	1.85	0.59
1:B:495:ARG:N	1:B:495:ARG:HD3	2.16	0.59
1:B:488:GLU:HG3	1:B:489:LYS:N	2.18	0.59
1:B:365:ARG:HG3	1:B:365:ARG:HH11	1.67	0.59
1:A:43:GLU:CD	1:A:43:GLU:H	2.06	0.59
1:A:99:GLY:HA3	3:A:295:AMP:O2'	2.03	0.59
1:B:325:ILE:HD12	1:B:397:VAL:HG22	1.84	0.59
1:B:498:TYR:CE1	1:B:553:PHE:HD1	2.21	0.58
1:A:23:LYS:O	1:A:26:ALA:HB3	2.03	0.58
1:A:134:ARG:NH2	1:B:539:GLY:HA3	2.18	0.58
1:B:311:VAL:HA	1:B:315:HIS:ND1	2.18	0.58
1:B:501:ARG:NH1	1:B:501:ARG:HG2	2.16	0.58
1:A:53:LEU:H	1:A:53:LEU:HD12	1.68	0.58
1:B:576:SER:O	1:B:580:MET:HG3	2.03	0.58
1:A:37:THR:HG23	1:A:37:THR:O	2.02	0.58
1:B:562:SER:HB3	1:B:565:GLU:CD	2.25	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:536:ASP:OD2	1:B:578:ARG:NH2	2.35	0.58
1:A:53:LEU:HD11	1:A:141:ARG:HD3	1.86	0.57
1:B:563:ILE:HD12	1:B:563:ILE:O	2.04	0.57
1:A:103:ARG:HH11	1:A:108:ALA:HB3	1.69	0.57
1:B:365:ARG:HG3	1:B:365:ARG:NH1	2.19	0.57
1:A:24:GLU:O	1:A:26:ALA:N	2.38	0.56
1:B:383:GLU:HG3	1:B:414:PHE:HZ	1.70	0.56
1:B:517:GLY:O	1:B:518:PHE:CB	2.52	0.56
1:B:527:ASN:HA	5:B:616:HOH:O	2.04	0.56
1:A:19:LEU:HD12	1:A:66:TYR:CD2	2.41	0.56
1:B:448:ARG:HB3	1:B:451:GLU:OE1	2.06	0.56
1:A:73:ASP:OD2	1:A:75:PHE:HB2	2.06	0.55
1:B:429:VAL:HA	1:B:433:LYS:O	2.07	0.55
1:A:129:VAL:HG23	1:A:129:VAL:O	2.05	0.55
1:B:480:THR:HG21	4:B:596:FMN:HM81	1.89	0.55
1:B:555:ARG:NH1	4:B:596:FMN:O4	2.38	0.55
1:B:488:GLU:HG3	1:B:489:LYS:H	1.71	0.55
1:B:398:VAL:HG11	1:B:402:PHE:CD2	2.41	0.55
1:A:64:SER:O	1:A:66:TYR:N	2.40	0.55
1:A:22:MET:HE1	1:A:67:ALA:HB2	1.87	0.55
1:B:502:VAL:HG22	1:B:548:LEU:HD22	1.88	0.55
1:A:166:VAL:CG1	1:A:167:HIS:H	2.19	0.54
1:B:302:VAL:N	1:B:332:ASP:O	2.40	0.54
1:A:178:PHE:N	1:A:179:PRO:CA	2.70	0.54
1:B:325:ILE:HD13	1:B:423:TYR:HD2	1.72	0.54
1:B:427:ASP:HB2	1:B:434:ARG:HH21	1.71	0.54
1:A:203:HIS:CE1	1:A:209:LYS:HG2	2.42	0.54
1:A:35:ILE:CD1	1:A:67:ALA:HB3	2.38	0.54
1:A:212:GLY:HA2	1:A:235:LEU:CD1	2.31	0.54
1:B:354:MET:HE2	1:B:359:ARG:N	2.21	0.54
1:A:205:PRO:O	1:A:206:ASP:HB2	2.08	0.54
1:A:254:MET:HG2	1:A:280:MET:HE1	1.90	0.54
1:A:88:ARG:CB	1:A:88:ARG:HH11	2.18	0.54
1:B:441:ARG:O	1:B:445:GLN:HG3	2.08	0.53
1:A:142:ASN:O	1:A:145:GLN:N	2.42	0.53
1:B:399:GLY:HA2	1:B:425:ILE:H	1.72	0.53
1:A:43:GLU:O	1:A:45:PHE:N	2.41	0.53
1:A:216:VAL:HG12	1:A:228:VAL:CG1	2.39	0.53
1:A:138:SER:HB2	1:B:508:LYS:HD2	1.89	0.53
1:B:558:LYS:HG2	1:B:559:LYS:N	2.23	0.53
1:A:116:ARG:HA	1:A:120:VAL:O	2.09	0.53
1:A:62:MET:O	1:A:63:LEU:C	2.47	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:272:GLN:O	1:A:276:SER:HB2	2.09	0.53
1:A:29:ARG:NH1	1:A:29:ARG:HG2	2.24	0.53
1:B:463:GLU:HG2	1:B:464:GLY:H	1.74	0.53
1:A:149:VAL:HG11	1:A:250:VAL:HG12	1.91	0.53
1:A:53:LEU:HD12	1:A:53:LEU:N	2.24	0.52
1:B:329:ARG:NH1	1:B:394:SER:HB2	2.24	0.52
1:A:265:GLU:HG2	1:A:268:ALA:HB3	1.91	0.52
1:A:106:LYS:O	1:A:107:ASN:HB2	2.10	0.52
1:B:496:GLY:HA2	1:B:557:GLU:HB2	1.90	0.52
1:B:495:ARG:CD	1:B:495:ARG:H	2.22	0.51
1:A:254:MET:HG2	1:A:280:MET:CE	2.40	0.51
1:A:103:ARG:HH11	1:A:108:ALA:CB	2.23	0.51
1:B:344:TYR:HA	1:B:349:PHE:CD2	2.45	0.51
1:B:495:ARG:H	1:B:495:ARG:HD3	1.76	0.51
1:B:353:LEU:HD23	1:B:491:VAL:HG11	1.92	0.51
1:B:344:TYR:OH	1:B:489:LYS:HG2	2.10	0.51
1:A:138:SER:CB	1:B:508:LYS:HD2	2.41	0.51
1:B:322:MET:CE	1:B:333:SER:HB3	2.40	0.51
1:B:495:ARG:NH1	1:B:495:ARG:HG2	2.25	0.51
1:A:216:VAL:HG12	1:A:228:VAL:HG13	1.93	0.51
1:B:467:HIS:CG	1:B:467:HIS:O	2.63	0.51
1:B:303:VAL:HG22	1:B:334:LEU:HB3	1.93	0.51
1:A:286:ASN:O	1:A:288:LYS:N	2.43	0.50
1:B:563:ILE:O	1:B:566:LEU:HB3	2.12	0.50
1:B:325:ILE:O	1:B:325:ILE:HG22	2.11	0.50
1:A:257:GLU:OE2	1:A:257:GLU:HA	2.10	0.50
1:A:90:LEU:O	1:A:93:VAL:HG12	2.11	0.50
1:A:202:VAL:O	1:A:209:LYS:HA	2.11	0.50
1:B:571:ASP:O	1:B:574:VAL:HG22	2.11	0.50
1:A:19:LEU:O	1:A:22:MET:CB	2.58	0.50
1:B:343:GLU:HA	1:B:346:LEU:HD12	1.94	0.50
1:A:134:ARG:NH2	1:B:538:GLU:O	2.44	0.50
1:A:64:SER:C	1:A:66:TYR:H	2.15	0.50
1:A:95:ALA:HA	1:A:121:GLU:O	2.11	0.50
1:A:24:GLU:C	1:A:26:ALA:N	2.65	0.50
1:B:552:LYS:HD3	1:B:580:MET:CE	2.41	0.50
1:B:398:VAL:HG21	1:B:402:PHE:HD2	1.77	0.50
1:A:27:PHE:C	1:A:29:ARG:H	2.16	0.50
1:A:14:GLY:HA2	1:A:128:VAL:HG21	1.93	0.49
1:B:532:VAL:HG11	1:B:548:LEU:HD11	1.94	0.49
1:A:204:LEU:HB2	1:A:208:LYS:HB2	1.94	0.49
1:A:149:VAL:HG11	1:A:250:VAL:HG11	1.94	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:460:PHE:CD1	1:B:461:GLU:N	2.81	0.49
1:A:37:THR:HB	1:A:63:LEU:CD1	2.41	0.49
1:B:396:VAL:HG12	1:B:396:VAL:O	2.11	0.49
1:A:28:PHE:CD1	1:A:28:PHE:O	2.64	0.49
1:A:103:ARG:HD3	1:A:108:ALA:CB	2.23	0.49
1:A:111:ASN:OD1	1:A:113:SER:HB2	2.13	0.49
1:A:288:LYS:HB2	5:A:630:HOH:O	2.12	0.49
1:A:7:VAL:O	1:A:8:PHE:HB2	2.13	0.49
1:B:404:PHE:CZ	1:B:415:LEU:HD21	2.48	0.49
1:B:349:PHE:CZ	1:B:352:LEU:HD23	2.48	0.48
1:A:204:LEU:HD22	1:A:204:LEU:N	2.28	0.48
1:B:516:VAL:HG22	1:B:530:TYR:CD1	2.48	0.48
1:A:167:HIS:HB3	1:A:168:LYS:H	1.47	0.48
1:A:73:ASP:HB3	1:A:76:ARG:HD3	1.95	0.48
1:A:268:ALA:O	1:A:272:GLN:OE1	2.31	0.48
1:B:320:ARG:HG3	1:B:320:ARG:HH11	1.79	0.48
1:A:195:ARG:CD	1:A:217:GLY:HA2	2.44	0.48
1:A:264:GLU:O	1:A:264:GLU:OE2	2.31	0.48
1:A:24:GLU:C	1:A:26:ALA:H	2.16	0.48
1:B:407:ASN:ND2	1:B:407:ASN:N	2.61	0.48
1:A:232:VAL:CG1	1:A:234:ILE:HG13	2.44	0.48
1:B:560:PHE:CD1	1:B:566:LEU:HA	2.49	0.48
1:B:349:PHE:HZ	1:B:352:LEU:HD23	1.79	0.48
1:B:322:MET:HE1	1:B:333:SER:HB3	1.95	0.47
1:A:34:LEU:HD11	1:A:70:VAL:CG2	2.44	0.47
1:A:113:SER:O	1:A:116:ARG:HB2	2.13	0.47
1:B:341:PRO:HB2	1:B:343:GLU:OE2	2.14	0.47
1:B:584:ILE:O	1:B:584:ILE:HG22	2.13	0.47
1:B:302:VAL:HG13	1:B:326:ALA:HB2	1.95	0.47
1:B:502:VAL:HG13	1:B:548:LEU:CD2	2.45	0.47
1:A:158:ARG:CG	1:A:158:ARG:NH1	2.75	0.47
1:A:43:GLU:O	1:A:44:TYR:C	2.53	0.47
1:B:415:LEU:HB2	1:B:422:VAL:CG2	2.44	0.47
1:B:571:ASP:HA	1:B:574:VAL:HG22	1.96	0.47
1:B:316:GLN:HB3	1:B:320:ARG:NH1	2.28	0.47
1:A:112:ALA:CB	1:A:122:VAL:HG11	2.44	0.47
1:B:416:ARG:HG3	1:B:416:ARG:HH11	1.79	0.47
3:A:295:AMP:H8	3:A:295:AMP:H5'1	1.79	0.47
1:A:116:ARG:NH2	1:A:122:VAL:O	2.48	0.47
1:B:322:MET:HA	1:B:397:VAL:HG21	1.96	0.47
1:A:265:GLU:O	1:A:268:ALA:HB3	2.14	0.46
1:B:465:ILE:HD13	1:B:545:ARG:HG2	1.95	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:3:VAL:HG11	1:A:34:LEU:HD23	1.98	0.46
1:A:284:ILE:H	1:A:284:ILE:HG13	1.58	0.46
1:A:216:VAL:CG1	1:A:228:VAL:CG1	2.94	0.46
1:B:415:LEU:HB2	1:B:422:VAL:HG21	1.97	0.46
1:B:497:VAL:HG22	1:B:515:ASN:HD22	1.80	0.46
1:A:201:ARG:HA	1:A:201:ARG:HD2	1.71	0.46
1:A:270:ILE:HD12	1:A:270:ILE:C	2.36	0.46
1:A:25:ILE:HG22	1:A:29:ARG:HD2	1.98	0.46
1:A:269:ALA:O	1:A:272:GLN:HB2	2.16	0.45
1:B:427:ASP:CB	1:B:434:ARG:HE	2.28	0.45
1:A:105:GLY:O	1:A:106:LYS:C	2.54	0.45
1:A:258:LYS:CG	1:A:259:LYS:N	2.80	0.45
1:B:495:ARG:HA	1:B:516:VAL:HG12	1.97	0.45
1:A:29:ARG:HH21	1:A:121:GLU:CD	2.20	0.45
1:B:429:VAL:HG12	1:B:433:LYS:H	1.82	0.45
1:A:27:PHE:O	1:A:29:ARG:N	2.49	0.45
1:A:103:ARG:O	1:A:104:PHE:HB3	2.15	0.45
1:A:60:VAL:CG1	1:A:60:VAL:O	2.65	0.45
1:B:495:ARG:HG2	1:B:495:ARG:HH11	1.81	0.45
1:A:137:SER:O	1:A:141:ARG:HB2	2.17	0.45
1:A:195:ARG:HD3	1:A:218:PHE:HD1	1.82	0.45
1:B:324:GLU:C	1:B:326:ALA:N	2.68	0.45
1:B:427:ASP:HA	3:B:595:AMP:N1	2.32	0.45
1:B:338:ILE:O	1:B:339:SER:C	2.55	0.45
1:B:391:SER:O	1:B:393:VAL:HG23	2.17	0.45
1:B:302:VAL:HB	1:B:395:ALA:HB3	1.99	0.44
1:B:412:ALA:HB1	1:B:422:VAL:HG11	1.98	0.44
1:B:313:ILE:HD11	1:B:435:VAL:HG21	2.00	0.44
1:B:558:LYS:CG	1:B:559:LYS:N	2.80	0.44
1:B:502:VAL:HG13	1:B:548:LEU:HD21	1.99	0.44
1:B:303:VAL:O	1:B:396:VAL:HA	2.17	0.44
1:A:184:ASP:C	1:A:186:GLY:H	2.21	0.44
1:A:2:VAL:HA	1:A:95:ALA:O	2.18	0.44
1:A:193:LEU:O	1:A:194:LYS:C	2.55	0.44
1:B:463:GLU:CG	1:B:464:GLY:N	2.77	0.44
1:B:566:LEU:O	1:B:570:ILE:HG13	2.17	0.44
1:B:437:SER:O	1:B:441:ARG:CG	2.62	0.44
1:B:400:ARG:HH11	1:B:400:ARG:HB3	1.83	0.44
1:A:9:ASP:OD1	1:A:53:LEU:HD12	2.19	0.43
1:A:273:ASP:O	1:A:276:SER:HB3	2.17	0.43
1:B:317:LYS:HE3	1:B:425:ILE:HG23	1.99	0.43
1:B:586:ASN:HA	1:B:589:PHE:HE1	1.83	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:456:LEU:HD12	1:B:456:LEU:O	2.17	0.43
1:A:281:ILE:O	1:A:283:ASP:N	2.51	0.43
1:A:167:HIS:O	1:A:168:LYS:CB	2.66	0.43
1:A:10:GLY:HA2	1:A:54:MET:SD	2.58	0.43
1:B:319:LEU:HD22	1:B:335:ILE:HD13	2.01	0.43
1:A:165:ILE:HG22	1:A:166:VAL:N	2.33	0.43
1:A:144:VAL:O	1:A:193:LEU:HD23	2.19	0.43
1:B:416:ARG:NH1	1:B:416:ARG:HG3	2.33	0.43
1:B:325:ILE:O	1:B:329:ARG:HG3	2.18	0.43
1:A:115:LEU:O	1:A:120:VAL:HG23	2.19	0.43
1:B:496:GLY:HA2	1:B:557:GLU:H	1.83	0.43
1:A:286:ASN:O	1:A:288:LYS:HG3	2.18	0.43
1:B:305:ILE:HG22	1:B:306:GLY:N	2.33	0.43
1:A:146:GLU:O	1:A:148:ARG:HG3	2.19	0.43
1:A:45:PHE:CE2	1:A:74:PHE:CD2	3.05	0.43
4:A:296:FMN:H1'1	4:A:296:FMN:H9	1.79	0.42
1:B:541:LEU:HA	1:B:544:GLN:OE1	2.19	0.42
1:A:195:ARG:CD	1:A:218:PHE:HD1	2.32	0.42
1:A:217:GLY:O	1:A:228:VAL:HA	2.19	0.42
1:A:245:ARG:O	1:A:246:LEU:HD12	2.18	0.42
1:B:428:VAL:HG23	3:B:595:AMP:N1	2.35	0.42
1:A:8:PHE:CE1	1:A:19:LEU:HD21	2.53	0.42
1:B:478:PHE:HB3	1:B:479:PRO:HD2	2.01	0.42
1:A:127:ASP:HA	3:A:295:AMP:N1	2.34	0.42
1:B:480:THR:HB	1:B:531:GLU:HB3	2.02	0.42
1:A:55:THR:OG1	1:A:58:SER:HB2	2.20	0.42
1:A:152:ILE:HD13	1:A:160:PHE:HB2	2.01	0.42
1:A:26:ALA:O	1:A:27:PHE:C	2.58	0.42
1:B:574:VAL:O	1:B:577:ALA:HB3	2.19	0.42
1:B:320:ARG:HG3	1:B:320:ARG:NH1	2.34	0.42
1:B:531:GLU:OE1	4:B:596:FMN:H4'	2.19	0.42
1:A:201:ARG:HB3	1:A:249:GLU:HB2	2.01	0.42
1:A:181:ALA:O	1:A:231:GLU:HA	2.19	0.42
1:A:180:THR:O	2:A:294:ADP:O1B	2.38	0.42
1:A:78:LYS:NZ	1:B:589:PHE:CZ	2.88	0.42
1:A:82:PRO:O	1:A:86:VAL:HG23	2.20	0.42
1:A:25:ILE:O	1:A:25:ILE:HG22	2.19	0.42
1:B:308:PHE:HA	1:B:315:HIS:CE1	2.55	0.42
1:A:205:PRO:O	1:A:206:ASP:CB	2.66	0.41
1:A:79:ASP:HA	1:A:106:LYS:HE3	2.01	0.41
1:A:105:GLY:O	1:A:108:ALA:N	2.53	0.41
1:A:137:SER:HB3	1:A:141:ARG:NH1	2.34	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:25:ILE:O	1:A:29:ARG:HB2	2.20	0.41
1:B:350:PRO:O	1:B:491:VAL:HA	2.20	0.41
1:A:232:VAL:HG13	1:A:234:ILE:HG13	2.02	0.41
1:A:5:ILE:HG22	1:A:6:GLY:N	2.36	0.41
1:A:195:ARG:HD3	1:A:217:GLY:HA2	2.02	0.41
1:B:502:VAL:O	1:B:509:LYS:HG2	2.20	0.41
1:B:479:PRO:HA	2:B:594:ADP:O2A	2.21	0.41
1:B:349:PHE:CE2	1:B:351:GLY:HA2	2.56	0.41
1:B:499:LEU:HD12	1:B:500:VAL:N	2.36	0.41
1:B:412:ALA:CB	1:B:422:VAL:HG11	2.50	0.41
1:A:64:SER:C	1:A:66:TYR:N	2.73	0.41
1:A:11:VAL:CA	1:A:15:HIS:ND1	2.83	0.41
1:A:142:ASN:O	1:A:143:LEU:C	2.59	0.41
1:B:585:ILE:C	1:B:587:SER:H	2.23	0.41
1:B:516:VAL:HG22	1:B:530:TYR:CE1	2.56	0.41
1:A:45:PHE:O	1:A:46:LEU:CD2	2.68	0.41
1:B:560:PHE:HD1	1:B:566:LEU:HD13	1.86	0.41
1:A:6:GLY:O	1:A:37:THR:HA	2.21	0.41
4:B:596:FMN:H9	4:B:596:FMN:H1'1	1.96	0.41
1:A:196:GLY:O	1:A:215:ASN:HA	2.21	0.41
1:B:452:ILE:O	1:B:453:PRO:C	2.59	0.41
1:A:218:PHE:HA	1:A:227:ASN:O	2.21	0.41
1:A:277:ALA:O	1:A:278:ARG:C	2.60	0.41
1:B:495:ARG:CD	1:B:495:ARG:N	2.80	0.40
1:A:144:VAL:O	1:A:193:LEU:HA	2.21	0.40
1:A:25:ILE:HD12	1:A:123:TYR:CD1	2.55	0.40
1:B:562:SER:O	1:B:563:ILE:C	2.60	0.40
1:A:8:PHE:HB3	1:A:63:LEU:HD21	2.03	0.40
1:B:480:THR:CG2	4:B:596:FMN:HM81	2.50	0.40
1:A:276:SER:O	1:A:279:ASN:HB2	2.21	0.40
1:A:156:LEU:HG	1:A:158:ARG:NH1	2.36	0.40
1:A:29:ARG:NH2	1:A:121:GLU:OE2	2.54	0.40
1:B:453:PRO:HA	1:B:458:ARG:O	2.21	0.40
1:B:527:ASN:C	1:B:527:ASN:HD22	2.24	0.40
1:B:308:PHE:CE1	1:B:319:LEU:HD21	2.56	0.40
1:B:311:VAL:O	1:B:311:VAL:HG12	2.21	0.40
1:B:358:SER:O	1:B:361:GLU:CB	2.67	0.40
1:B:400:ARG:NH1	1:B:400:ARG:HB3	2.36	0.40
1:A:167:HIS:CD2	1:A:242:TYR:HD1	2.39	0.40
1:B:402:PHE:CD1	1:B:403:ARG:N	2.89	0.40
1:B:500:VAL:HG21	1:B:548:LEU:HD13	2.02	0.40
1:A:143:LEU:HG	1:A:143:LEU:H	1.74	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:162:ILE:HD12	1:A:185:ARG:NE	2.36	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:A:602:HOH:O	5:B:629:HOH:O[2_656]	2.10	0.10

## 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	260/293 (89%)	210 (81%)	39 (15%)	11 (4%)	4	13
1	B	264/293 (90%)	222 (84%)	34 (13%)	8 (3%)	7	22
All	All	524/586 (89%)	432 (82%)	73 (14%)	19 (4%)	5	17

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	PHE
1	A	44	TYR
1	A	65	ARG
1	B	398	VAL
1	A	25	ILE
1	A	26	ALA
1	A	102	PHE
1	A	131	GLN
1	A	185	ARG
1	A	282	ASP
1	B	375	PHE
1	B	479	PRO
1	B	341	PRO
1	B	557	GLU
1	A	281	ILE

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Mol	Chain	Res	Type
1	B	323	LYS
1	B	339	SER
1	B	563	ILE
1	A	132	GLY

### 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	237/257 (92%)	208 (88%)	29 (12%)	7	20
1	B	239/257 (93%)	222 (93%)	17 (7%)	21	51
All	All	476/514 (93%)	430 (90%)	46 (10%)	12	32

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	MET
1	A	24	GLU
1	A	28	PHE
1	A	31	ASP
1	A	36	TYR
1	A	57	GLU
1	A	58	SER
1	A	65	ARG
1	A	69	THR
1	A	77	ILE
1	A	93	VAL
1	A	101	ASP
1	A	120	VAL
1	A	127	ASP
1	A	131	GLN
1	A	141	ARG
1	A	163	GLU
1	A	167	HIS
1	A	178	PHE
1	A	183	ILE
1	A	188	GLU

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Mol	Chain	Res	Type
1	A	189	LYS
1	A	204	LEU
1	A	213	VAL
1	A	232	VAL
1	A	244	GLN
1	A	264	GLU
1	A	272	GLN
1	A	283	ASP
1	B	302	VAL
1	B	332	ASP
1	B	334	LEU
1	B	400	ARG
1	B	407	ASN
1	B	416	ARG
1	B	426	GLU
1	B	427	ASP
1	B	439	LEU
1	B	467	HIS
1	B	479	PRO
1	B	482	ASN
1	B	495	ARG
1	B	527	ASN
1	B	563	ILE
1	B	583	ASP
1	B	589	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	131	GLN
1	A	167	HIS
1	A	244	GLN
1	A	272	GLN
1	A	279	ASN
1	B	407	ASN
1	B	442	ASN
1	B	515	ASN
1	B	527	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

6 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	ADP	A	294	-	29,29,29	1.56	7 (24%)	45,45,45	2.46	10 (22%)
3	AMP	A	295	-	25,25,25	1.88	5 (20%)	38,38,38	1.98	10 (26%)
4	FMN	A	296	-	33,33,33	2.57	10 (30%)	46,50,50	3.30	17 (36%)
2	ADP	B	594	-	29,29,29	1.44	5 (17%)	45,45,45	2.31	5 (11%)
3	AMP	B	595	-	25,25,25	2.35	5 (20%)	38,38,38	1.84	9 (23%)
4	FMN	B	596	-	33,33,33	2.62	12 (36%)	46,50,50	3.22	17 (36%)

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	ADP	A	294	-	-	0/16/32/32	0/1/3/3
3	AMP	A	295	-	-	0/10/26/26	0/1/3/3
4	FMN	A	296	-	-	0/18/18/18	0/0/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	B	594	-	-	0/16/32/32	0/1/3/3
3	AMP	B	595	-	-	0/10/26/26	0/1/3/3
4	FMN	B	596	-	-	0/18/18/18	0/0/3/3

All (44) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	596	FMN	C4A-C10	8.02	1.54	1.40
4	A	296	FMN	C4A-C10	7.88	1.54	1.40
3	B	595	AMP	C2'-C1'	7.30	1.63	1.53
4	A	296	FMN	C2-N3	5.93	1.48	1.37
4	B	596	FMN	C2-N3	5.54	1.48	1.37
3	B	595	AMP	C3'-C4'	5.25	1.67	1.53
3	A	295	AMP	C2'-C1'	4.92	1.60	1.53
4	B	596	FMN	C4-N3	4.39	1.44	1.37
4	B	596	FMN	C5A-N5	4.38	1.42	1.35
4	B	596	FMN	C9-C9A	4.12	1.48	1.40
4	A	296	FMN	C9-C9A	4.04	1.48	1.40
3	B	595	AMP	C8-N9	3.96	1.42	1.36
4	A	296	FMN	C5A-N5	3.92	1.41	1.35
3	A	295	AMP	C3'-C4'	3.90	1.63	1.53
4	B	596	FMN	C5'-C4'	3.86	1.57	1.51
4	A	296	FMN	C4A-N5	-3.83	1.28	1.36
4	B	596	FMN	C9A-C5A	3.79	1.50	1.42
3	B	595	AMP	O3'-C3'	3.76	1.52	1.43
4	A	296	FMN	C4-N3	3.57	1.43	1.37
4	A	296	FMN	C5'-C4'	3.55	1.57	1.51
3	A	295	AMP	O3'-C3'	3.54	1.51	1.43
2	B	594	ADP	C4-N9	-3.47	1.32	1.37
3	A	295	AMP	C8-N9	3.46	1.41	1.36
2	A	294	ADP	C4-N9	-3.40	1.32	1.37
4	A	296	FMN	C9A-C5A	3.11	1.48	1.42
4	A	296	FMN	C6-C7	2.92	1.45	1.37
4	B	596	FMN	C6-C7	2.90	1.45	1.37
2	B	594	ADP	O4'-C1'	2.90	1.45	1.41
2	A	294	ADP	C2-N3	2.82	1.37	1.32
2	A	294	ADP	O4'-C1'	2.74	1.45	1.41
2	A	294	ADP	C5-N7	-2.58	1.30	1.40
4	B	596	FMN	C4A-N5	-2.54	1.30	1.36
3	B	595	AMP	O4'-C1'	2.54	1.45	1.41
2	B	594	ADP	C2-N3	2.50	1.37	1.32
2	B	594	ADP	C5-N7	-2.36	1.31	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	296	FMN	C8-C7	2.33	1.47	1.40
2	A	294	ADP	C4-N3	2.30	1.39	1.35
4	B	596	FMN	C4'-C3'	2.22	1.58	1.53
3	A	295	AMP	C4-N9	-2.21	1.34	1.37
4	B	596	FMN	C8-C7	2.12	1.47	1.40
2	A	294	ADP	C5'-C4'	2.04	1.58	1.51
2	A	294	ADP	PA-O2A	-2.03	1.46	1.55
2	B	594	ADP	C4-N3	2.02	1.38	1.35
4	B	596	FMN	O3'-C3'	2.02	1.47	1.43

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	594	ADP	N3-C2-N1	-12.57	118.20	128.71
2	A	294	ADP	N3-C2-N1	-12.39	118.36	128.71
4	B	596	FMN	C1'-N10-C9A	9.85	128.46	118.87
4	A	296	FMN	C2-N1-C10	9.33	124.38	114.98
4	B	596	FMN	C2-N1-C10	9.28	124.33	114.98
4	A	296	FMN	C2'-C1'-N10	8.84	124.17	112.45
4	A	296	FMN	C5A-C9A-N10	8.16	124.84	116.80
4	B	596	FMN	C5A-C9A-N10	7.92	124.59	116.80
4	A	296	FMN	C1'-N10-C9A	7.10	125.78	118.87
4	A	296	FMN	C9A-N10-C10	-6.95	114.95	121.77
4	B	596	FMN	C9A-N10-C10	-6.56	115.33	121.77
3	A	295	AMP	C4-C5-N7	5.56	114.28	109.52
2	A	294	ADP	N3-C4-N9	5.22	134.86	125.43
4	B	596	FMN	C2'-C1'-N10	5.21	119.36	112.45
3	B	595	AMP	C4-C5-N7	5.07	113.86	109.52
4	A	296	FMN	C4A-C10-N1	-5.03	117.70	122.73
3	A	295	AMP	N3-C2-N1	-4.86	124.65	128.71
4	A	296	FMN	C7M-C7-C6	-4.85	108.69	120.38
2	B	594	ADP	N3-C4-N9	4.83	134.15	125.43
3	B	595	AMP	N3-C2-N1	-4.62	124.85	128.71
4	B	596	FMN	C7M-C7-C6	-4.60	109.30	120.38
4	A	296	FMN	C7M-C7-C8	4.42	130.95	120.74
4	B	596	FMN	C4A-C10-N1	-4.41	118.32	122.73
2	A	294	ADP	O4'-C1'-N9	4.31	112.45	108.44
4	B	596	FMN	C7M-C7-C8	4.25	130.55	120.74
4	B	596	FMN	C9A-C5A-N5	-3.57	116.89	122.37
4	A	296	FMN	C4A-N5-C5A	3.37	120.48	116.69
3	A	295	AMP	C8-N9-C4	-3.32	104.36	106.90
4	B	596	FMN	C5'-C4'-C3'	3.30	118.28	112.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	595	AMP	C8-N9-C4	-3.29	104.39	106.90
4	A	296	FMN	C9A-C5A-N5	-3.28	117.33	122.37
3	A	295	AMP	O4'-C4'-C5'	-3.28	97.64	109.36
3	A	295	AMP	N6-C6-N1	3.22	125.69	119.36
4	A	296	FMN	C1'-C2'-C3'	-3.17	100.75	109.82
4	B	596	FMN	C4A-N5-C5A	3.16	120.24	116.69
4	A	296	FMN	C5'-C4'-C3'	3.13	117.96	112.06
3	A	295	AMP	O3'-C3'-C4'	-3.11	101.91	111.08
4	A	296	FMN	C4'-C3'-C2'	3.10	120.25	113.25
2	B	594	ADP	C2-N3-C4	3.08	122.78	114.01
4	B	596	FMN	C4'-C3'-C2'	3.06	120.16	113.25
4	B	596	FMN	C6-C5A-N5	3.04	122.52	118.97
3	B	595	AMP	C3'-C2'-C1'	3.04	105.66	100.91
2	A	294	ADP	C2-N3-C4	2.99	122.53	114.01
2	A	294	ADP	C5-C4-N3	-2.94	119.30	125.70
3	B	595	AMP	N6-C6-N1	2.86	124.98	119.36
4	A	296	FMN	C4-C4A-C10	2.84	121.54	116.95
3	A	295	AMP	C3'-C2'-C1'	2.81	105.30	100.91
4	A	296	FMN	C6-C5A-N5	2.80	122.23	118.97
3	B	595	AMP	O4'-C4'-C5'	-2.80	99.37	109.36
2	B	594	ADP	C5-C4-N3	-2.80	119.61	125.70
3	B	595	AMP	C2-N1-C6	2.76	123.75	118.77
2	A	294	ADP	C1'-N9-C4	2.66	131.23	126.64
4	B	596	FMN	C1'-C2'-C3'	-2.60	102.39	109.82
2	A	294	ADP	C8-N9-C1'	-2.60	121.26	126.38
3	A	295	AMP	C2-N1-C6	2.60	123.46	118.77
2	B	594	ADP	C4-C5-N7	-2.54	107.35	109.52
3	B	595	AMP	O3'-C3'-C4'	-2.38	104.08	111.08
4	B	596	FMN	C4-C4A-C10	2.30	120.67	116.95
2	A	294	ADP	C2'-C3'-C4'	2.26	107.16	102.65
4	A	296	FMN	O3'-C3'-C4'	-2.24	103.06	108.74
3	A	295	AMP	C2'-C1'-N9	-2.22	107.56	113.27
2	A	294	ADP	C4-C5-N7	-2.21	107.62	109.52
4	B	596	FMN	O2'-C2'-C3'	2.20	114.52	109.05
4	A	296	FMN	O2'-C2'-C3'	2.13	114.36	109.05
3	B	595	AMP	O4'-C4'-C3'	2.11	109.45	105.17
2	A	294	ADP	C4'-O4'-C1'	2.10	112.03	109.75
4	B	596	FMN	C1'-N10-C10	-2.08	116.21	119.17
3	A	295	AMP	P-O5'-C5'	-2.05	112.25	118.19

There are no chirality outliers.

There are no torsion outliers.



There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	268/293 (91%)	0.17	17 (6%)	19 18	31, 77, 119, 144	0
1	B	270/293 (92%)	0.01	8 (2%)	48 49	42, 74, 116, 141	0
All	All	538/586 (91%)	0.09	25 (4%)	31 31	31, 75, 117, 144	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	478	PHE	4.9
1	A	226	ARG	4.1
1	A	227	ASN	4.1
1	A	258	LYS	3.9
1	A	178	PHE	3.5
1	A	265	GLU	3.4
1	A	219	ARG	2.9
1	B	572	GLN	2.9
1	B	401	ASP	2.8
1	A	66	TYR	2.7
1	A	45	PHE	2.7
1	A	123	TYR	2.6
1	B	589	PHE	2.6
1	A	242	TYR	2.5
1	A	6	GLY	2.4
1	A	117	LYS	2.4
1	B	585	ILE	2.4
1	A	206	ASP	2.4
1	B	455	TYR	2.3
1	A	240	ASP	2.2
1	A	115	LEU	2.2
1	A	118	LYS	2.1
1	B	479	PRO	2.1
1	B	542	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	259	LYS	2.1

## 6.2 Non-standard residues in protein, DNA, RNA chains

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

## 6.3 Carbohydrates

There are no carbohydrates in this entry.

## 6.4 Ligands

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	AMP	B	595	23/23	0.29	1.59	76,87,93,93	0
4	FMN	B	596	31/31	0.26	0.69	59,66,127,128	0
3	AMP	A	295	23/23	0.26	0.58	72,75,86,86	0
4	FMN	A	296	31/31	0.26	0.16	77,85,123,123	0
2	ADP	A	294	27/27	0.22	-0.27	118,126,148,148	0
2	ADP	B	594	27/27	0.17	-0.64	124,130,138,138	0

## 6.5 Other polymers

There are no such residues in this entry.