



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2018 – 10:02 am 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 X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : trunk31020
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk31020

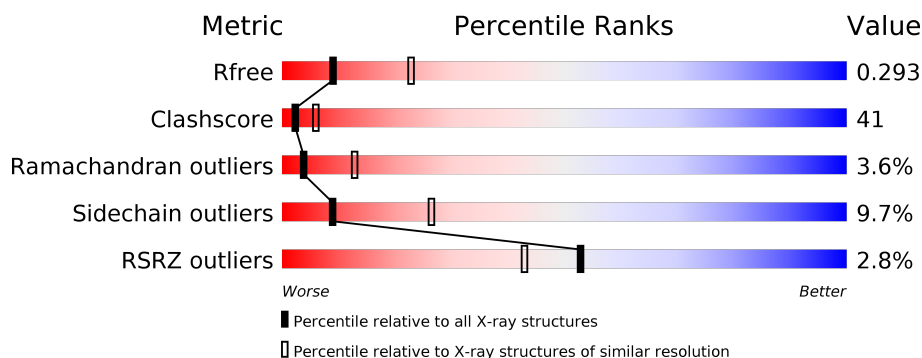
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	111664	2792 (2.80-2.80)
Clashscore	122126	3209 (2.80-2.80)
Ramachandran outliers	120053	3158 (2.80-2.80)
Sidechain outliers	120020	3160 (2.80-2.80)
RSRZ outliers	108989	2726 (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 ≥ 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	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 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 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		

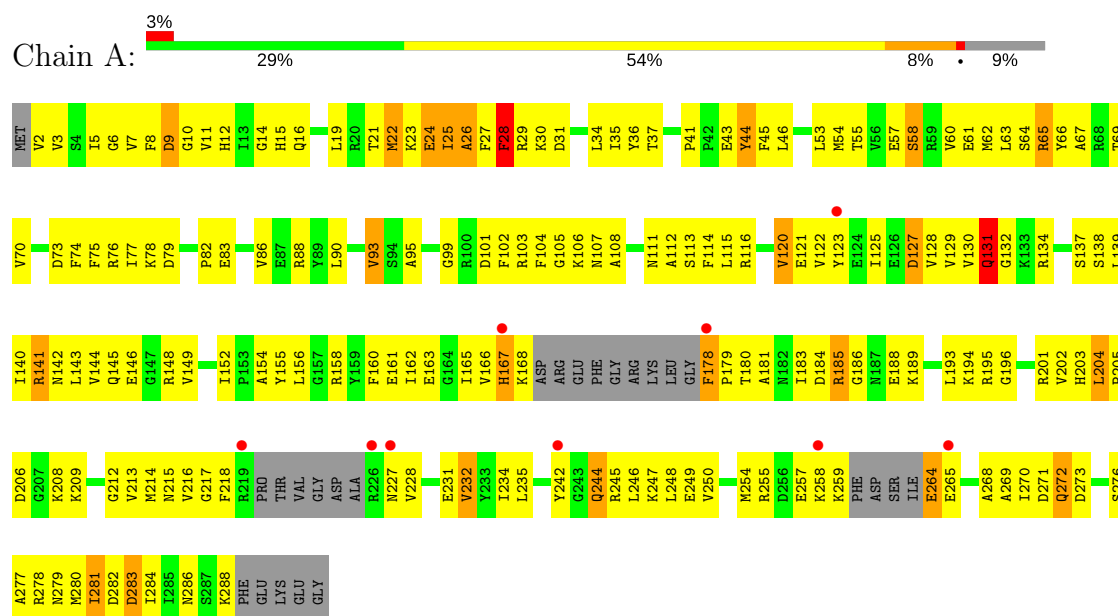
- Molecule 5 is water.

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

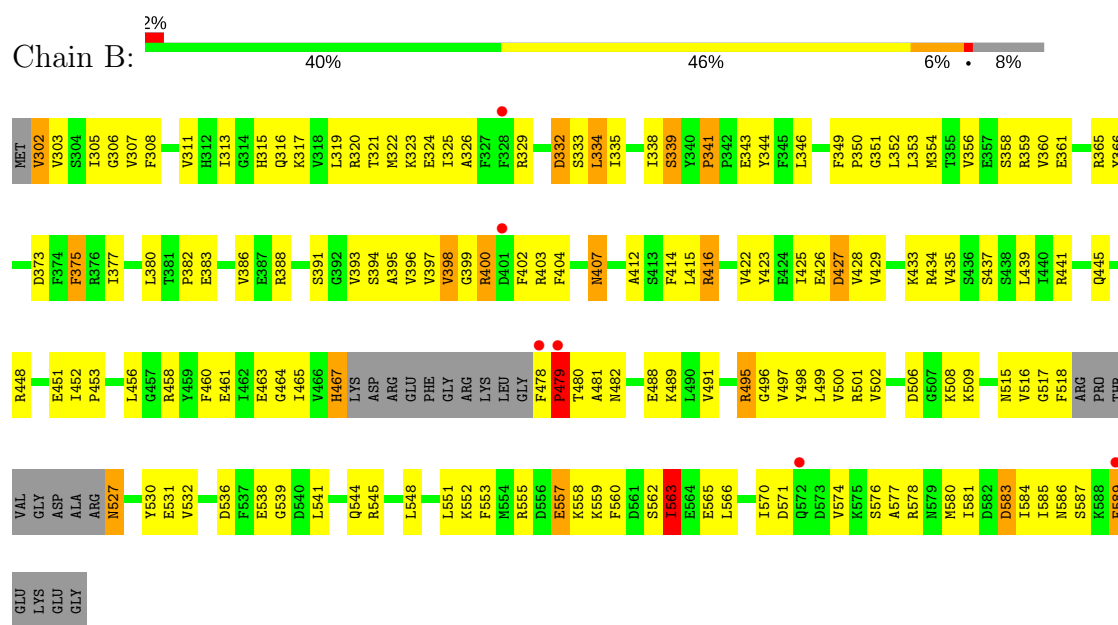
3 Residue-property plots

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



- Molecule 1: riboflavin kinase/FMN adenylyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	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
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 2.80Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.231 , 0.298 0.229 , 0.293	Depositor DCC
R_{free} test set	1639 reflections (10.09%)	wwPDB-VP
Wilson B-factor (Å ²)	76.4	Xtriage
Anisotropy	0.231	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 67.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.024 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4560	wwPDB-VP
Average B, all atoms (Å ²)	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.*

¹Intensities estimated from amplitudes.

²Theoretical 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: 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 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	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
4	A	31	0	19	1	0
4	B	31	0	19	6	0

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

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 41.

All (368) 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: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:A:139:LEU:HD12	1:B:506:ASP:OD1	1.86	0.74
1:B:465:ILE:CD1	1:B:545:ARG:HG2	2.16	0.74
1:B:501:ARG:HB2	1:B:551:LEU:HD11	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (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:HD22	1:B:407:ASN:N	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:8:PHE:HA	1:A:15:HIS:CE1	2.30	0.67
1:A:166:VAL:HG12	1:A:167:HIS:N	2.08	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:130:VAL:HG21	1:A:154:ALA:O	1.97	0.64
1:A:34:LEU:HD11	1:A:70:VAL:HG23	1.80	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:LEU:O	1:A:22:MET:HB3	1.99	0.62
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:112:ALA:HB1	1:A:122:VAL:HG11	1.80	0.62
1:A:258:LYS:HG2	1:A:259:LYS:N	2.13	0.62
1:A:127:ASP:N	1:A:127:ASP:OD2	2.32	0.61
1:A:245:ARG:C	1:A:246:LEU:HD12	2.20	0.61
1:B:316:GLN:HB3	1:B:320:ARG:HH12	1.65	0.61
1:A:28:PHE:HD1	1:A:28:PHE:O	1.83	0.61
1:A:45:PHE:HE2	1:A:74:PHE:CD2	2.18	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:58:SER:O	1:A:61:GLU:HB3	2.02	0.59
1:A:83:GLU:HG3	1:A:114:PHE:HZ	1.67	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:488:GLU:HG3	1:B:489:LYS:N	2.18	0.59
1:B:495:ARG:N	1:B:495:ARG:HD3	2.16	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:B:311:VAL:HA	1:B:315:HIS:ND1	2.18	0.58
1:A:134:ARG:NH2	1:B:539:GLY:HA3	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:488:GLU:HG3	1:B:489:LYS:H	1.71	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: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:A:142:ASN:O	1:A:145:GLN:N	2.42	0.53
1:B:441:ARG:O	1:B:445:GLN:HG3	2.08	0.53
1:B:399:GLY:HA2	1:B:425:ILE:H	1.72	0.53
1:A:138:SER:HB2	1:B:508:LYS:HD2	1.89	0.53
1:A:216:VAL:HG12	1:A:228:VAL:CG1	2.39	0.53
1:A:43:GLU:O	1:A:45:PHE:N	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:A:149:VAL:HG11	1:A:250:VAL:HG12	1.91	0.53
1:B:463:GLU:HG2	1:B:464:GLY:H	1.74	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:A:254:MET:HG2	1:A:280:MET:CE	2.40	0.51
1:B:495:ARG:CD	1:B:495:ARG:H	2.22	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:353:LEU:HD23	1:B:491:VAL:HG11	1.92	0.51
1:B:495:ARG:H	1:B:495:ARG:HD3	1.76	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: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:495:ARG:NH1	1:B:495:ARG:HG2	2.25	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: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:B:325:ILE:O	1:B:325:ILE:HG22	2.11	0.50
1:B:563:ILE:O	1:B:566:LEU:HB3	2.12	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:398:VAL:HG21	1:B:402:PHE:HD2	1.77	0.50
1:B:552:LYS:HD3	1:B:580:MET:CE	2.41	0.50
1:A:27:PHE:C	1:A:29:ARG:H	2.16	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:GLY:HA2	1:A:128:VAL:HG21	1.93	0.49
1:A:204:LEU:HB2	1:A:208:LYS:HB2	1.94	0.49
1:B:532:VAL:HG11	1:B:548:LEU:HD11	1.94	0.49
1:A:149:VAL:HG11	1:A:250:VAL:HG11	1.94	0.49
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:A:204:LEU:HD22	1:A:204:LEU:N	2.28	0.48
1:B:349:PHE:CZ	1:B:352:LEU:HD23	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:B:516:VAL:HG22	1:B:530:TYR:CD1	2.48	0.48
1:A:268:ALA:O	1:A:272:GLN:OE1	2.31	0.48
1:A:195:ARG:CD	1:A:217:GLY:HA2	2.44	0.48
1:B:320:ARG:HG3	1:B:320:ARG:HH11	1.79	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:349:PHE:HZ	1:B:352:LEU:HD23	1.79	0.48
1:B:560:PHE:CD1	1:B:566:LEU:HA	2.49	0.48
1:B:322:MET:HE1	1:B:333:SER:HB3	1.95	0.47
1:A:113:SER:O	1:A:116:ARG:HB2	2.13	0.47
1:A:34:LEU:HD11	1:A:70:VAL:CG2	2.44	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:316:GLN:HB3	1:B:320:ARG:NH1	2.28	0.47
1:B:571:ASP:HA	1:B:574:VAL:HG22	1.96	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:ARG:NH2	1:A:122:VAL:O	2.48	0.47
3:A:295:AMP:H8	3:A:295:AMP:H5'1	1.79	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
1:A:284:ILE:H	1:A:284:ILE:HG13	1.58	0.46
1:A:3:VAL:HG11	1:A:34:LEU:HD23	1.98	0.46
1:A:201:ARG:HA	1:A:201:ARG:HD2	1.71	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: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: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:B:495:ARG:HA	1:B:516:VAL:HG12	1.97	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: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:495:ARG:HG2	1:B:495:ARG:HH11	1.81	0.45
1:B:324:GLU:C	1:B:326:ALA:N	2.68	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:427:ASP:HA	3:B:595:AMP:N1	2.32	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:A:184:ASP:C	1:A:186:GLY:H	2.21	0.44
1:B:303:VAL:O	1:B:396:VAL:HA	2.17	0.44
1:B:502:VAL:HG13	1:B:548:LEU:HD21	1.99	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:437:SER:O	1:B:441:ARG:CG	2.62	0.44
1:B:463:GLU:CG	1:B:464:GLY:N	2.77	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:566:LEU:O	1:B:570:ILE:HG13	2.17	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:456:LEU:HD12	1:B:456:LEU:O	2.17	0.43
1:B:586:ASN:HA	1:B:589:PHE:HE1	1.83	0.43
1:A:167:HIS:O	1:A:168:LYS:CB	2.66	0.43
1:A:281:ILE:O	1:A:283:ASP:N	2.51	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:A:115:LEU:O	1:A:120:VAL:HG23	2.19	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:B:325:ILE:O	1:B:329:ARG:HG3	2.18	0.43
1:B:496:GLY:HA2	1:B:557:GLU:H	1.83	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:A:55:THR:OG1	1:A:58:SER:HB2	2.20	0.42
1:B:480:THR:HB	1:B:531:GLU:HB3	2.02	0.42
1:A:152:ILE:HD13	1:A:160:PHE:HB2	2.01	0.42
1:A:180:THR:O	2:A:294:ADP:O1B	2.38	0.42
1:A:181:ALA:O	1:A:231:GLU:HA	2.19	0.42
1:A:201:ARG:HB3	1:A:249:GLU:HB2	2.01	0.42
1:A:26:ALA:O	1:A:27:PHE:C	2.58	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:B:574:VAL:O	1:B:577:ALA:HB3	2.19	0.42
1:A:82:PRO:O	1:A:86:VAL:HG23	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:LYS:NZ	1:B:589:PHE:CZ	2.88	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
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:195:ARG:HD3	1:A:217:GLY:HA2	2.02	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: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:412:ALA:CB	1:B:422:VAL:HG11	2.50	0.41
1:B:499:LEU:HD12	1:B:500:VAL:N	2.36	0.41
1:A:64:SER:C	1:A:66:TYR:N	2.73	0.41
1:A:142:ASN:O	1:A:143:LEU:C	2.59	0.41
1:A:11:VAL:CA	1:A:15:HIS:ND1	2.83	0.41
1:B:585:ILE:C	1:B:587:SER:H	2.23	0.41
1:A:196:GLY:O	1:A:215:ASN:HA	2.21	0.41
1:A:45:PHE:O	1:A:46:LEU:CD2	2.68	0.41
1:A:6:GLY:O	1:A:37:THR:HA	2.21	0.41
1:B:452:ILE:O	1:B:453:PRO:C	2.59	0.41
1:B:560:PHE:HD1	1:B:566:LEU:HD13	1.86	0.41
1:B:516:VAL:HG22	1:B:530:TYR:CE1	2.56	0.41
4:B:596:FMN:H9	4:B:596:FMN:H1'1	1.96	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: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:A:8:PHE:HB3	1:A:63:LEU:HD21	2.03	0.40
1:B:495:ARG:CD	1:B:495:ARG:N	2.80	0.40
1:B:562:SER:O	1:B:563:ILE:C	2.60	0.40
1:B:480:THR:CG2	4:B:596:FMN:HM81	2.50	0.40
1:A:156:LEU:HG	1:A:158:ARG:NH1	2.36	0.40
1:A:276:SER:O	1:A:279:ASN:HB2	2.21	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:308:PHE:CE1	1:B:319:LEU:HD21	2.56	0.40
1:B:311:VAL:HG12	1:B:311:VAL:O	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:143:LEU:HG	1:A:143:LEU:H	1.74	0.40
1:A:162:ILE:HD12	1:A:185:ARG:NE	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

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	Interatomic distance (Å)	Clash overlap (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%)	3	10
1	B	264/293 (90%)	222 (84%)	34 (13%)	8 (3%)	5	16
All	All	524/586 (89%)	432 (82%)	73 (14%)	19 (4%)	4	13

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

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Mol	Chain	Res	Type
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
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%)	5	16
1	B	239/257 (93%)	222 (93%)	17 (7%)	16	42
All	All	476/514 (93%)	430 (90%)	46 (10%)	9	26

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

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Mol	Chain	Res	Type
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
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 [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	-	25,29,29	1.57	6 (24%)	25,45,45	2.71	4 (16%)
3	AMP	A	295	-	22,25,25	1.74	4 (18%)	23,38,38	2.21	8 (34%)
4	FMN	A	296	-	31,33,33	2.77	10 (32%)	39,50,50	3.25	15 (38%)
2	ADP	B	594	-	25,29,29	1.44	4 (16%)	25,45,45	2.65	2 (8%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	AMP	B	595	-	22,25,25	2.08	5 (22%)	23,38,38	2.03	7 (30%)
4	FMN	B	596	-	31,33,33	2.85	11 (35%)	39,50,50	3.43	14 (35%)

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	-	-	0/12/32/32	0/3/3/3
3	AMP	A	295	-	-	0/6/26/26	0/3/3/3
4	FMN	A	296	-	-	0/16/18/18	0/3/3/3
2	ADP	B	594	-	-	0/12/32/32	0/3/3/3
3	AMP	B	595	-	-	0/6/26/26	0/3/3/3
4	FMN	B	596	-	-	0/16/18/18	0/3/3/3

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	296	FMN	C4A-N5	-3.62	1.28	1.33
2	A	294	ADP	C5-N7	-2.60	1.30	1.39
2	B	594	ADP	C5-N7	-2.35	1.31	1.39
4	B	596	FMN	C1'-N10	-2.28	1.46	1.48
3	B	595	AMP	C2'-C3'	2.03	1.58	1.53
2	A	294	ADP	C2-N1	2.07	1.37	1.33
2	A	294	ADP	C5'-C4'	2.08	1.58	1.51
2	B	594	ADP	C4-N3	2.17	1.38	1.35
3	A	295	AMP	O4'-C1'	2.27	1.44	1.41
4	B	596	FMN	C4'-C3'	2.37	1.58	1.53
4	B	596	FMN	C8-C7	2.41	1.47	1.40
2	A	294	ADP	C4-N3	2.47	1.39	1.35
4	A	296	FMN	C8-C7	2.65	1.47	1.40
3	B	595	AMP	O4'-C1'	2.87	1.45	1.41
2	B	594	ADP	C2-N3	3.06	1.37	1.32
4	B	596	FMN	C6-C7	3.08	1.45	1.37
2	A	294	ADP	O4'-C1'	3.09	1.45	1.41
4	A	296	FMN	C6-C7	3.10	1.45	1.37
4	A	296	FMN	C9A-C5A	3.11	1.48	1.42
2	B	594	ADP	O4'-C1'	3.26	1.45	1.41
2	A	294	ADP	C2-N3	3.45	1.37	1.32
3	A	295	AMP	O3'-C3'	3.47	1.51	1.43
3	B	595	AMP	O3'-C3'	3.68	1.52	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	296	FMN	C9-C9A	3.79	1.48	1.40
4	B	596	FMN	C9A-C5A	3.79	1.50	1.42
4	A	296	FMN	C5'-C4'	3.82	1.57	1.51
4	B	596	FMN	C9-C9A	3.86	1.48	1.40
4	A	296	FMN	C5A-N5	3.92	1.41	1.35
3	A	295	AMP	C8-N9	4.01	1.41	1.36
4	B	596	FMN	C5'-C4'	4.15	1.57	1.51
3	A	295	AMP	C3'-C4'	4.17	1.63	1.53
4	B	596	FMN	C5A-N5	4.39	1.42	1.35
3	B	595	AMP	C8-N9	4.62	1.42	1.36
4	B	596	FMN	C2-N3	5.00	1.48	1.38
4	A	296	FMN	C2-N3	5.39	1.48	1.38
3	B	595	AMP	C3'-C4'	5.60	1.67	1.53
4	A	296	FMN	C4-N3	5.66	1.43	1.33
4	B	596	FMN	C4-N3	6.42	1.44	1.33
4	A	296	FMN	C4A-C10	8.28	1.54	1.41
4	B	596	FMN	C4A-C10	8.43	1.54	1.41

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	594	ADP	N3-C2-N1	-12.46	118.20	128.86
2	A	294	ADP	N3-C2-N1	-12.28	118.36	128.86
4	A	296	FMN	C4A-C4-N3	-5.35	115.87	123.47
4	B	596	FMN	C4A-C4-N3	-5.17	116.11	123.47
4	A	296	FMN	C9A-N10-C10	-5.12	114.95	121.77
3	A	295	AMP	N3-C2-N1	-4.93	124.65	128.86
4	B	596	FMN	C9A-N10-C10	-4.83	115.33	121.77
4	A	296	FMN	C7M-C7-C6	-4.74	108.69	120.36
3	B	595	AMP	N3-C2-N1	-4.69	124.85	128.86
4	B	596	FMN	C7M-C7-C6	-4.49	109.30	120.36
4	B	596	FMN	C9A-C5A-N5	-3.82	116.89	122.32
3	A	295	AMP	O4'-C4'-C5'	-3.53	97.64	109.39
4	A	296	FMN	C9A-C5A-N5	-3.52	117.33	122.32
4	A	296	FMN	C1'-C2'-C3'	-3.17	100.75	109.82
3	A	295	AMP	O3'-C3'-C4'	-3.16	101.91	111.06
3	B	595	AMP	O4'-C4'-C5'	-3.01	99.37	109.39
4	B	596	FMN	C1'-C2'-C3'	-2.60	102.39	109.82
3	B	595	AMP	O3'-C3'-C4'	-2.41	104.08	111.06
4	A	296	FMN	O3'-C3'-C4'	-2.34	103.06	108.82
4	B	596	FMN	C1'-N10-C10	-2.25	116.21	118.46
3	A	295	AMP	P-O5'-C5'	-2.19	112.25	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	594	ADP	C4-C5-N7	-2.14	107.35	109.41
3	A	295	AMP	O4'-C4'-C3'	2.00	109.11	105.15
2	A	294	ADP	C4'-O4'-C1'	2.11	112.03	109.83
4	A	296	FMN	C4-C4A-C10	2.13	121.54	119.95
4	A	296	FMN	O2'-C2'-C3'	2.13	114.36	109.10
3	B	595	AMP	O4'-C4'-C3'	2.17	109.45	105.15
4	B	596	FMN	O2'-C2'-C3'	2.19	114.52	109.10
2	A	294	ADP	C2'-C3'-C4'	2.37	107.16	102.62
2	A	294	ADP	C1'-N9-C4	2.66	131.23	126.64
3	A	295	AMP	C2-N1-C6	2.77	123.46	118.75
4	A	296	FMN	C6-C5A-N5	2.86	122.23	118.95
3	B	595	AMP	C2-N1-C6	2.94	123.75	118.75
3	B	595	AMP	N6-C6-N1	3.09	124.98	118.57
4	B	596	FMN	C6-C5A-N5	3.11	122.52	118.95
4	B	596	FMN	C4'-C3'-C2'	3.20	120.16	113.40
4	A	296	FMN	C4'-C3'-C2'	3.25	120.25	113.40
4	B	596	FMN	C4A-N5-C5A	3.32	120.24	116.76
3	A	295	AMP	N6-C6-N1	3.43	125.69	118.57
4	A	296	FMN	C4A-N5-C5A	3.55	120.48	116.76
3	B	595	AMP	C4-C5-N7	4.61	113.86	109.41
4	B	596	FMN	C7M-C7-C8	4.71	130.55	120.72
4	A	296	FMN	C7M-C7-C8	4.90	130.95	120.72
3	A	295	AMP	C4-C5-N7	5.04	114.28	109.41
4	A	296	FMN	C4-N3-C2	8.32	122.23	115.14
4	A	296	FMN	C1'-N10-C9A	8.41	125.78	118.31
4	B	596	FMN	C4-N3-C2	8.53	122.41	115.14
4	B	596	FMN	C5A-C9A-N10	9.02	124.59	117.71
4	A	296	FMN	C5A-C9A-N10	9.34	124.84	117.71
4	B	596	FMN	C1'-N10-C9A	11.42	128.46	118.31

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	294	ADP	1	0
3	A	295	AMP	4	0
4	A	296	FMN	1	0
2	B	594	ADP	1	0
3	B	595	AMP	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	596	FMN	6	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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, 95th 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 ²)	Q < 0.9
1	A	268/293 (91%)	0.04	9 (3%)	45 35	31, 77, 119, 144	0
1	B	270/293 (92%)	-0.04	6 (2%)	62 52	42, 74, 116, 141	0
All	All	538/586 (91%)	0.00	15 (2%)	53 43	31, 75, 117, 144	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	478	PHE	3.7
1	A	227	ASN	3.4
1	B	479	PRO	3.2
1	A	178	PHE	3.0
1	A	265	GLU	3.0
1	A	242	TYR	2.9
1	A	226	ARG	2.8
1	A	123	TYR	2.6
1	B	401	ASP	2.5
1	A	219	ARG	2.4
1	A	167	HIS	2.4
1	B	572	GLN	2.4
1	B	328	PHE	2.2
1	A	258	LYS	2.2
1	B	589	PHE	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. The B-factors column lists the minimum, median, 95th 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	B-factors(Å ²)	Q<0.9
2	ADP	A	294	27/27	0.80	0.26	118,126,148,148	0
2	ADP	B	594	27/27	0.85	0.20	124,130,138,138	0
4	FMN	A	296	31/31	0.86	0.24	77,85,123,123	0
4	FMN	B	596	31/31	0.87	0.24	59,66,127,128	0
3	AMP	B	595	23/23	0.87	0.29	76,87,93,93	0
3	AMP	A	295	23/23	0.91	0.25	72,75,86,86	0

6.5 Other polymers [i](#)

There are no such residues in this entry.