



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 06:45 PM GMT

PDB ID : 3K30
Title : Histamine dehydrogenase from Nocardioides simplex
Authors : Scott, E.E.; Reed, T.M.; Limburg, J.
Deposited on : 2009-09-30
Resolution : 2.70 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
<http://wwpdb.org/validation/2016/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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

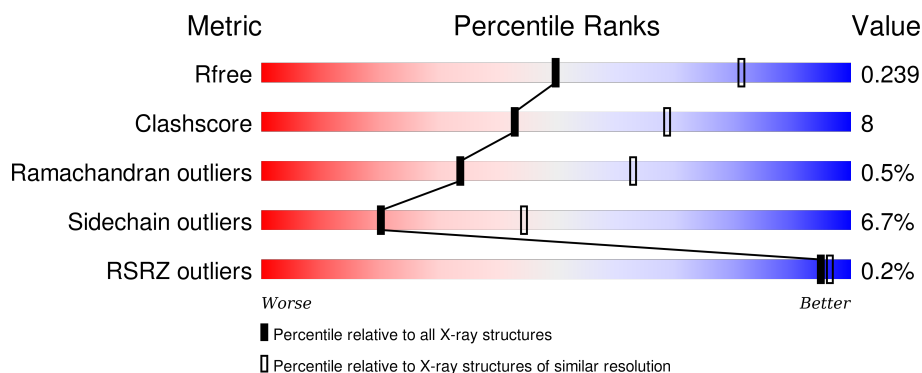
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.70 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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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	690	 76% 20% ..
1	B	690	 81% 17% ..

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 10684 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.

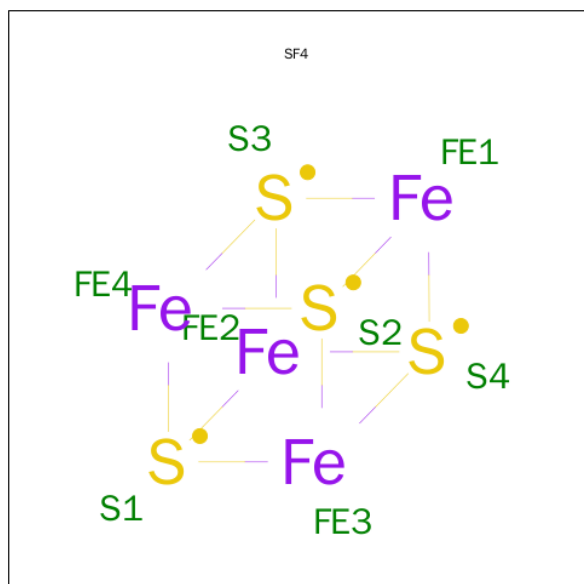
- Molecule 1 is a protein called Histamine dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	684	Total	C	N	O	S	0	0	0
			5272	3294	956	1001	21			
1	B	684	Total	C	N	O	S	0	0	0
			5272	3294	956	1001	21			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	615	ALA	THR	SEE REMARK 999	UNP Q6IWIJ5
B	615	ALA	THR	SEE REMARK 999	UNP Q6IWIJ5

- Molecule 2 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe₄S₄).



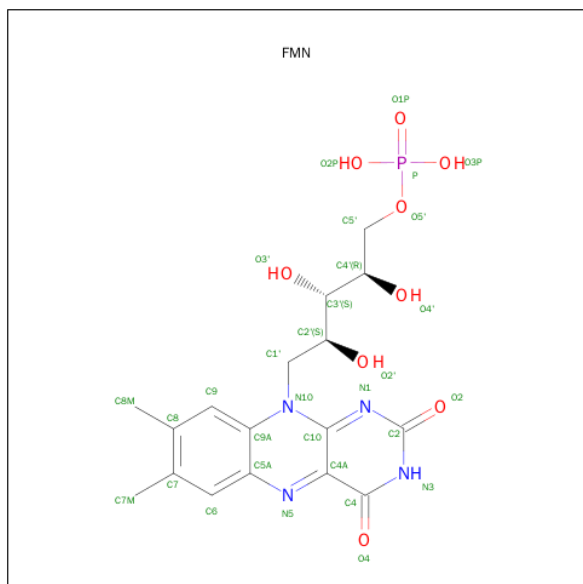
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			8	4	4		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 3 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



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

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $\text{C}_{10}\text{H}_{15}\text{N}_5\text{O}_{10}\text{P}_2$).



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

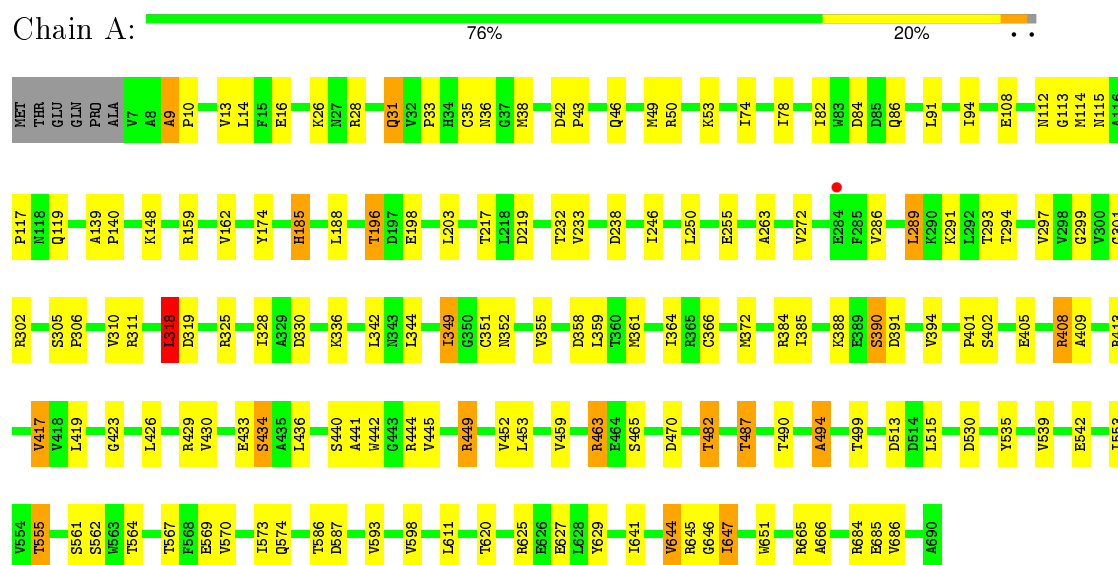
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	O	0	0
			2	2		
5	B	6	Total	O	0	0
			6	6		

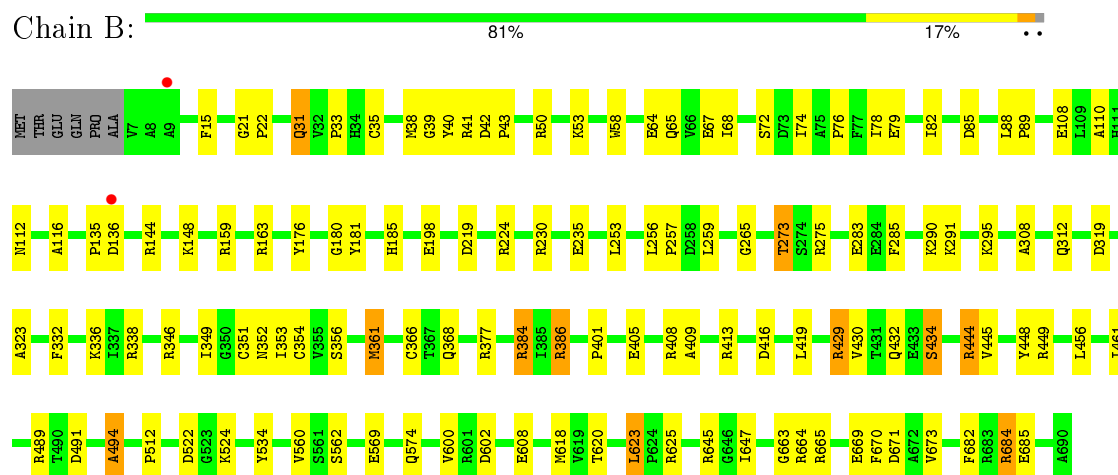
3 Residue-property plots

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: Histamine dehydrogenase



• Molecule 1: Histamine dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.14Å 107.03Å 153.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	84.43 – 2.70 84.43 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.6 (84.43-2.70) 99.6 (84.43-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.63 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.182 , 0.242 0.184 , 0.239	Depositor DCC
R_{free} test set	2336 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	29.2	Xtriage
Anisotropy	0.055	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 34.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 46249 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10684	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% 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, 0.375 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: FMN, SF4, 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	0.83	2/5379 (0.0%)	0.86	4/7306 (0.1%)
1	B	0.83	0/5379	0.85	3/7306 (0.0%)
All	All	0.83	2/10758 (0.0%)	0.86	7/14612 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	542	GLU	CG-CD	5.82	1.60	1.51
1	A	627	GLU	CG-CD	5.30	1.59	1.51

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	429	ARG	NE-CZ-NH2	-5.75	117.43	120.30
1	A	114	MET	CG-SD-CE	-5.73	91.03	100.20
1	A	318	LEU	CA-CB-CG	5.73	128.49	115.30
1	B	386	ARG	NE-CZ-NH2	-5.43	117.59	120.30
1	A	302	ARG	NE-CZ-NH2	5.34	122.97	120.30
1	A	645	ARG	NE-CZ-NH1	-5.24	117.68	120.30
1	B	664	ARG	NE-CZ-NH2	-5.16	117.72	120.30

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	5272	0	5177	96	0
1	B	5272	0	5176	79	0
2	A	8	0	0	1	0
2	B	8	0	0	1	0
3	A	31	0	19	4	0
3	B	31	0	18	5	0
4	A	27	0	12	2	0
4	B	27	0	12	2	0
5	A	2	0	0	0	0
5	B	6	0	0	0	0
All	All	10684	0	10414	172	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 8.

All (172) 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:35:CYS:SG	3:A:802:FMN:H6	1.13	1.70
1:A:35:CYS:SG	3:A:802:FMN:C6	2.08	1.41
1:A:351:CYS:O	1:A:352:ASN:HB2	1.64	0.98
1:B:384:ARG:HH11	1:B:384:ARG:HG3	1.30	0.97
1:A:352:ASN:ND2	1:A:567:THR:O	2.05	0.89
1:A:349:ILE:HD11	1:A:444:ARG:HB2	1.55	0.88
1:B:384:ARG:NH1	1:B:384:ARG:HG3	1.90	0.84
1:B:74:ILE:HD11	1:B:79:GLU:HG3	1.59	0.82
1:A:16:GLU:HB2	1:A:28:ARG:NH1	1.96	0.80
1:A:9:ALA:HB1	1:A:10:PRO:CD	2.13	0.78
1:A:349:ILE:CD1	1:A:444:ARG:HB2	2.13	0.77
1:A:482:THR:HG21	1:A:629:TYR:HB2	1.65	0.76
1:B:273:THR:CG2	1:B:275:ARG:HB3	2.16	0.76
1:B:290:LYS:NZ	1:B:319:ASP:OD2	2.20	0.74
1:A:16:GLU:HB2	1:A:28:ARG:HH11	1.49	0.74
1:B:136:ASP:OD1	1:B:181:TYR:OH	2.05	0.73
1:B:159:ARG:NH2	1:B:219:ASP:OD2	2.21	0.73
1:B:41:ARG:HH11	1:B:41:ARG:HG2	1.55	0.71
1:B:110:ALA:CB	1:B:176:TYR:HD2	2.02	0.71
1:B:429:ARG:HA	1:B:432:GLN:HE21	1.54	0.71
1:B:110:ALA:HB1	1:B:176:TYR:HD2	1.56	0.70
1:A:401:PRO:HD2	4:A:803:ADP:O3B	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:ALA:CB	1:A:10:PRO:CD	2.70	0.70
1:B:494:ALA:HB2	1:B:620:THR:HB	1.73	0.70
1:B:273:THR:HG21	1:B:569:GLU:OE2	1.92	0.69
1:A:84:ASP:HB3	1:A:86:GLN:H	1.58	0.68
1:A:647:ILE:HD11	1:A:666:ALA:HB2	1.75	0.67
1:A:233:VAL:HG22	1:A:246:ILE:HG12	1.75	0.67
1:A:31:GLN:HE22	1:A:50:ARG:HH11	1.42	0.66
1:A:405:GLU:OE2	1:A:408:ARG:NH1	2.27	0.66
1:A:9:ALA:HB1	1:A:10:PRO:HD2	1.78	0.66
1:A:9:ALA:CB	1:A:10:PRO:HD2	2.26	0.65
1:A:401:PRO:HG3	1:A:430:VAL:HG23	1.82	0.62
1:B:110:ALA:CB	1:B:176:TYR:CD2	2.83	0.62
1:A:441:ALA:O	1:A:444:ARG:HG3	1.99	0.62
1:B:43:PRO:HD2	1:B:685:GLU:OE1	2.00	0.61
1:B:15:PHE:CE2	1:B:338:ARG:HB2	2.36	0.61
1:B:135:PRO:O	1:B:136:ASP:OD2	2.19	0.61
1:A:487:THR:HG22	1:A:513:ASP:OD2	2.02	0.59
1:B:110:ALA:HB2	1:B:176:TYR:CD2	2.38	0.59
1:A:13:VAL:O	1:A:28:ARG:HD2	2.03	0.59
1:A:570:VAL:O	1:A:574:GLN:HG2	2.04	0.58
1:A:82:ILE:HG23	1:A:82:ILE:O	2.04	0.58
1:A:349:ILE:HG13	1:A:442:TRP:HA	1.86	0.57
1:B:82:ILE:O	1:B:82:ILE:HG23	2.04	0.57
1:B:35:CYS:HA	1:B:64:GLU:HB3	1.86	0.57
1:B:489:ARG:HG2	1:B:623:LEU:HD22	1.87	0.57
1:A:409:ALA:O	1:A:413:ARG:HG3	2.06	0.56
1:A:38:MET:HB2	1:A:42:ASP:HB2	1.88	0.56
1:A:36:ASN:HA	1:A:364:ILE:HG12	1.87	0.56
1:A:564:THR:HB	1:A:569:GLU:HB2	1.87	0.56
1:A:646:GLY:O	1:A:647:ILE:HD12	2.07	0.55
1:B:31:GLN:HE21	1:B:53:LYS:HD3	1.71	0.55
1:A:351:CYS:O	1:A:352:ASN:CB	2.45	0.55
1:A:49:MET:SD	1:A:372:MET:HE3	2.46	0.55
1:B:351:CYS:O	1:B:352:ASN:HB2	2.07	0.54
1:A:43:PRO:HD2	1:A:685:GLU:OE1	2.07	0.54
1:A:286:VAL:HA	1:A:289:LEU:HD22	1.88	0.54
1:B:31:GLN:HE21	1:B:53:LYS:CD	2.21	0.54
1:A:330:ASP:OD2	1:A:336:LYS:HE2	2.06	0.54
1:B:64:GLU:N	1:B:64:GLU:OE1	2.34	0.53
1:A:306:PRO:HG2	1:A:342:LEU:HD13	1.89	0.53
1:B:38:MET:HB3	1:B:42:ASP:HB2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:423:GLY:O	1:A:463:ARG:O	2.27	0.52
1:A:84:ASP:HB2	1:B:684:ARG:HH21	1.73	0.52
1:B:384:ARG:CG	1:B:384:ARG:HH11	2.11	0.52
1:A:9:ALA:HB1	1:A:10:PRO:HD3	1.90	0.52
1:A:82:ILE:HD12	1:A:91:LEU:HD13	1.91	0.52
1:B:429:ARG:HG2	4:B:803:ADP:O2A	2.09	0.52
1:B:349:ILE:HG22	1:B:444:ARG:HB2	1.92	0.52
1:B:332:PHE:O	1:B:336:LYS:HG3	2.09	0.51
1:A:417:VAL:HG13	1:A:459:VAL:HG22	1.92	0.51
1:A:232:THR:HA	1:A:263:ALA:O	2.11	0.51
1:A:586:THR:O	1:A:587:ASP:HB2	2.09	0.51
1:A:31:GLN:NE2	1:A:53:LYS:HZ2	2.08	0.51
1:A:31:GLN:HE21	1:A:53:LYS:NZ	2.09	0.50
1:A:366:CYS:HA	2:A:801:SF4:S4	2.52	0.50
1:B:489:ARG:NH1	1:B:491:ASP:OD2	2.44	0.50
1:A:445:VAL:O	1:A:449:ARG:HD2	2.12	0.50
1:B:419:LEU:HD23	1:B:461:ILE:HG12	1.93	0.50
1:A:31:GLN:HE21	1:A:53:LYS:HZ2	1.59	0.50
1:A:561:SER:O	1:A:564:THR:OG1	2.26	0.49
1:A:352:ASN:HB3	1:A:567:THR:HB	1.93	0.49
1:A:299:GLY:HA3	1:A:318:LEU:HD11	1.95	0.49
1:B:67:GLU:HG2	1:B:72:SER:O	2.12	0.49
1:B:273:THR:HG23	1:B:275:ARG:HB3	1.92	0.48
1:A:430:VAL:O	1:A:434:SER:OG	2.28	0.48
1:B:353:ILE:O	1:B:356:SER:HB3	2.12	0.48
1:B:670:PHE:O	1:B:671:ASP:HB2	2.12	0.48
1:B:351:CYS:SG	1:B:353:ILE:HB	2.53	0.48
1:A:433:GLU:OE1	1:A:535:TYR:OH	2.20	0.48
1:A:33:PRO:O	3:A:802:FMN:O2'	2.26	0.48
1:A:530:ASP:O	1:A:555:THR:HG23	2.13	0.48
1:A:325:ARG:HA	1:A:328:ILE:HD12	1.95	0.48
1:A:46:GLN:HG2	1:A:94:ILE:HG12	1.94	0.48
1:B:377:ARG:HB2	1:B:682:PHE:HB2	1.96	0.48
1:A:388:LYS:HD2	1:A:390:SER:O	2.14	0.48
1:A:35:CYS:H	3:A:802:FMN:C5A	2.27	0.47
1:B:112:ASN:HB2	1:B:116:ALA:HB2	1.96	0.47
1:B:409:ALA:O	1:B:413:ARG:HG3	2.15	0.47
1:B:253:LEU:HG	1:B:256:LEU:HD12	1.96	0.47
1:A:384:ARG:C	1:A:385:ILE:HG13	2.34	0.47
1:B:429:ARG:HA	1:B:432:GLN:NE2	2.28	0.47
1:B:405:GLU:OE2	1:B:408:ARG:NH1	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:ARG:NH2	1:A:219:ASP:OD2	2.48	0.47
1:A:355:VAL:O	1:A:358:ASP:N	2.48	0.46
1:B:40:TYR:CE2	1:B:78:ILE:HG13	2.50	0.46
1:B:346:ARG:HG2	1:B:368:GLN:HB3	1.96	0.46
1:B:366:CYS:HA	2:B:801:SF4:S4	2.56	0.46
1:B:346:ARG:HD3	1:B:448:TYR:CE2	2.50	0.46
1:B:110:ALA:HB1	1:B:176:TYR:CD2	2.44	0.46
1:B:401:PRO:HD2	4:B:803:ADP:O2A	2.15	0.45
1:B:33:PRO:O	3:B:802:FMN:O2'	2.28	0.45
1:B:430:VAL:O	1:B:434:SER:OG	2.25	0.45
1:A:453:LEU:HD22	1:A:459:VAL:HG11	1.99	0.45
1:A:641:ILE:HD11	1:A:644:VAL:HG22	1.99	0.45
1:B:600:VAL:O	1:B:608:GLU:HA	2.16	0.45
1:A:344:LEU:HD22	1:A:452:VAL:HG22	1.99	0.45
1:A:108:GLU:HA	1:A:174:TYR:HB2	1.98	0.45
1:B:41:ARG:HH11	1:B:41:ARG:CG	2.29	0.45
1:B:560:VAL:H	1:B:574:GLN:HE22	1.64	0.45
1:B:35:CYS:H	3:B:802:FMN:C5A	2.30	0.44
1:A:26:LYS:HE2	1:A:26:LYS:HB2	1.84	0.44
1:A:196:THR:CG2	1:A:196:THR:O	2.66	0.44
1:A:569:GLU:O	1:A:573:ILE:HG12	2.18	0.44
1:B:647:ILE:HD12	1:B:663:GLY:HA2	1.99	0.44
1:B:257:PRO:O	1:B:295:LYS:HE2	2.17	0.44
1:B:35:CYS:H	3:B:802:FMN:C6	2.32	0.43
1:B:384:ARG:NH1	1:B:384:ARG:CG	2.67	0.43
1:A:355:VAL:CG1	1:A:359:LEU:HD13	2.48	0.43
1:A:203:LEU:HA	1:A:203:LEU:HD23	1.72	0.43
1:B:88:LEU:N	1:B:89:PRO:CD	2.82	0.43
1:B:68:ILE:HA	1:B:82:ILE:HG22	2.00	0.43
1:B:230:ARG:NH1	3:B:802:FMN:O2	2.50	0.43
1:B:21:GLY:HA3	1:B:22:PRO:HD2	1.68	0.43
1:B:41:ARG:HG2	1:B:41:ARG:NH1	2.31	0.43
1:B:108:GLU:OE2	3:B:802:FMN:O2	2.36	0.43
1:A:139:ALA:HA	1:A:140:PRO:HD3	1.83	0.43
1:A:593:VAL:HG22	1:A:598:VAL:HG12	2.00	0.43
1:A:49:MET:HE3	1:A:50:ARG:CD	2.49	0.43
1:B:76:PRO:O	1:B:361:MET:HG2	2.19	0.43
1:A:112:ASN:O	1:A:115:ASN:HB3	2.19	0.43
1:A:84:ASP:HB2	1:B:684:ARG:NH2	2.33	0.42
1:A:31:GLN:NE2	1:A:53:LYS:NZ	2.67	0.42
1:B:180:GLY:HA2	1:B:185:HIS:CD2	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:647:ILE:CD1	1:A:666:ALA:HB2	2.46	0.42
1:B:512:PRO:HD3	1:B:618:MET:O	2.19	0.42
1:A:297:VAL:H	1:A:319:ASP:HB2	1.85	0.42
1:A:49:MET:SD	1:A:372:MET:CE	3.07	0.41
1:A:625:ARG:HG3	1:A:625:ARG:NH1	2.35	0.41
1:A:429:ARG:HG2	4:A:803:ADP:O1A	2.20	0.41
1:B:349:ILE:HG21	1:B:445:VAL:HG23	2.02	0.41
1:B:235:GLU:HA	1:B:265:GLY:HA3	2.01	0.41
1:A:651:TRP:O	1:B:144:ARG:HD3	2.20	0.41
1:A:436:LEU:HB3	1:A:539:VAL:CG1	2.51	0.41
1:A:494:ALA:HB2	1:A:620:THR:HB	2.02	0.41
1:A:185:HIS:CD2	1:A:185:HIS:C	2.93	0.41
1:A:349:ILE:HG22	1:A:351:CYS:HB3	2.03	0.41
1:A:31:GLN:HE22	1:A:50:ARG:NH1	2.15	0.41
1:A:306:PRO:O	1:A:310:VAL:HG23	2.21	0.41
1:B:308:ALA:O	1:B:312:GLN:HG2	2.21	0.41
1:A:162:VAL:HG21	1:A:217:THR:HG23	2.03	0.41
1:A:402:SER:HB2	1:A:647:ILE:HG23	2.02	0.41
1:B:31:GLN:HE22	1:B:50:ARG:HH11	1.68	0.41
1:A:113:GLY:C	1:A:115:ASN:H	2.24	0.41
1:A:148:LYS:HE3	1:A:198:GLU:OE2	2.21	0.41
1:B:58:TRP:CH2	1:B:323:ALA:HB3	2.56	0.40
1:A:641:ILE:CD1	1:A:644:VAL:HG22	2.51	0.40
1:B:65:GLN:HE22	1:B:112:ASN:HD21	1.69	0.40
1:B:645:ARG:NH2	1:B:669:GLU:OE1	2.51	0.40
1:A:117:PRO:HB2	1:A:119:GLN:HG3	2.04	0.40
1:B:283:GLU:C	1:B:285:PHE:N	2.73	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	682/690 (99%)	637 (93%)	40 (6%)	5 (1%)	26	55
1	B	682/690 (99%)	652 (96%)	28 (4%)	2 (0%)	46	75
All	All	1364/1380 (99%)	1289 (94%)	68 (5%)	7 (0%)	34	63

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	9	ALA
1	B	39	GLY
1	A	494	ALA
1	B	494	ALA
1	A	291	LYS
1	A	470	ASP
1	A	301	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	544/549 (99%)	499 (92%)	45 (8%)	14	31
1	B	544/549 (99%)	516 (95%)	28 (5%)	29	59
All	All	1088/1098 (99%)	1015 (93%)	73 (7%)	20	44

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	LEU
1	A	31	GLN
1	A	74	ILE
1	A	78	ILE
1	A	185	HIS
1	A	188	LEU
1	A	196	THR
1	A	238	ASP
1	A	250	LEU

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Mol	Chain	Res	Type
1	A	255	GLU
1	A	272	VAL
1	A	289	LEU
1	A	293	THR
1	A	294	THR
1	A	305	SER
1	A	311	ARG
1	A	318	LEU
1	A	349	ILE
1	A	361	MET
1	A	390	SER
1	A	391	ASP
1	A	394	VAL
1	A	408	ARG
1	A	417	VAL
1	A	419	LEU
1	A	426	LEU
1	A	434	SER
1	A	440	SER
1	A	449	ARG
1	A	463	ARG
1	A	465	SER
1	A	482	THR
1	A	487	THR
1	A	490	THR
1	A	499	THR
1	A	515	LEU
1	A	553	ILE
1	A	555	THR
1	A	562	SER
1	A	611	LEU
1	A	644	VAL
1	A	647	ILE
1	A	665	ARG
1	A	684	ARG
1	A	686	VAL
1	B	31	GLN
1	B	85	ASP
1	B	148	LYS
1	B	163	ARG
1	B	198	GLU
1	B	224	ARG

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Mol	Chain	Res	Type
1	B	259	LEU
1	B	273	THR
1	B	291	LYS
1	B	354	CYS
1	B	361	MET
1	B	384	ARG
1	B	386	ARG
1	B	416	ASP
1	B	434	SER
1	B	444	ARG
1	B	449	ARG
1	B	456	LEU
1	B	522	ASP
1	B	524	LYS
1	B	534	TYR
1	B	562	SER
1	B	602	ASP
1	B	623	LEU
1	B	625	ARG
1	B	665	ARG
1	B	673	VAL
1	B	684	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	B	31	GLN
1	B	112	ASN
1	B	432	GLN
1	B	574	GLN

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 [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	SF4	A	801	1	0,12,12	0.00	-	0,24,24	0.00	-
3	FMN	A	802	-	32,33,33	1.19	3 (9%)	34,50,50	2.36	10 (29%)
4	ADP	A	803	-	24,29,29	1.11	1 (4%)	23,45,45	1.94	5 (21%)
2	SF4	B	801	1	0,12,12	0.00	-	0,24,24	0.00	-
3	FMN	B	802	-	32,33,33	1.12	2 (6%)	34,50,50	2.32	11 (32%)
4	ADP	B	803	-	24,29,29	1.01	1 (4%)	23,45,45	1.89	3 (13%)

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	SF4	A	801	1	-	0/0/48/48	0/6/5/5
3	FMN	A	802	-	-	0/18/18/18	0/3/3/3
4	ADP	A	803	-	-	0/12/32/32	0/3/3/3
2	SF4	B	801	1	-	0/0/48/48	0/6/5/5
3	FMN	B	802	-	-	0/18/18/18	0/3/3/3
4	ADP	B	803	-	-	0/12/32/32	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	802	FMN	C2-N1	-2.30	1.33	1.38
4	B	803	ADP	C5-C4	2.27	1.45	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	802	FMN	C4A-N5	2.41	1.37	1.33
3	B	802	FMN	C9A-N10	2.80	1.42	1.38
3	A	802	FMN	C10-N10	3.00	1.42	1.39
3	A	802	FMN	C9A-N10	3.14	1.43	1.38
4	A	803	ADP	C5-C4	3.27	1.47	1.40

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	803	ADP	N3-C2-N1	-6.66	123.64	128.87
4	B	803	ADP	N3-C2-N1	-6.29	123.93	128.87
3	A	802	FMN	N3-C2-N1	-5.93	117.70	127.69
3	B	802	FMN	C4-C4A-C10	-4.39	117.13	119.94
3	B	802	FMN	C4A-C4-N3	-4.29	117.91	123.52
3	A	802	FMN	C4-C4A-C10	-4.25	117.22	119.94
3	B	802	FMN	N3-C2-N1	-4.24	120.55	127.69
3	A	802	FMN	C4A-C4-N3	-3.42	119.05	123.52
4	B	803	ADP	C1'-N9-C4	-3.05	123.40	126.81
4	A	803	ADP	C1'-N9-C4	-2.85	123.63	126.81
3	A	802	FMN	C9A-C5A-N5	-2.69	117.81	122.18
3	B	802	FMN	O3P-P-O5'	-2.47	99.52	106.72
4	A	803	ADP	C4'-O4'-C1'	-2.42	107.08	109.64
3	A	802	FMN	C6-C7-C8	-2.19	115.68	119.97
4	A	803	ADP	O3B-PB-O2B	2.06	115.00	107.44
4	A	803	ADP	C2-N1-C6	2.13	122.56	118.77
3	A	802	FMN	O3P-P-O2P	2.13	115.27	107.44
3	A	802	FMN	C1'-N10-C9A	2.30	121.50	118.83
4	B	803	ADP	O2A-PA-O3A	2.35	115.33	105.27
3	B	802	FMN	O3P-P-O2P	2.38	116.20	107.44
3	B	802	FMN	O4'-C4'-C5'	2.45	115.43	110.09
3	B	802	FMN	O3P-P-O1P	2.47	118.69	110.63
3	B	802	FMN	C4-C4A-N5	2.92	122.25	118.70
3	B	802	FMN	C4A-N5-C5A	3.62	120.99	116.72
3	A	802	FMN	C4-C4A-N5	3.92	123.46	118.70
3	B	802	FMN	C1'-N10-C9A	4.25	123.75	118.83
3	A	802	FMN	C4A-N5-C5A	5.34	123.01	116.72
3	A	802	FMN	C4-N3-C2	5.45	119.70	115.16
3	B	802	FMN	C4-N3-C2	6.62	120.68	115.16

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	801	SF4	1	0
3	A	802	FMN	4	0
4	A	803	ADP	2	0
2	B	801	SF4	1	0
3	B	802	FMN	5	0
4	B	803	ADP	2	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	684/690 (99%)	0.01	1 (0%) 95 97	6, 21, 38, 45	0
1	B	684/690 (99%)	-0.09	2 (0%) 94 95	3, 19, 35, 47	0
All	All	1368/1380 (99%)	-0.04	3 (0%) 95 96	3, 20, 36, 47	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	9	ALA	2.5
1	A	284	GLU	2.2
1	B	136	ASP	2.1

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	LLDF	B-factors(\AA^2)	Q<0.9
3	FMN	B	802	31/31	0.98	0.19	0.50	14,18,22,25	0
3	FMN	A	802	31/31	0.97	0.19	0.14	12,20,23,23	0
4	ADP	B	803	27/27	0.99	0.15	-0.07	9,15,18,19	0
4	ADP	A	803	27/27	0.98	0.15	-0.36	12,17,27,28	0
2	SF4	A	801	8/8	0.97	0.07	-2.32	19,20,25,26	0
2	SF4	B	801	8/8	0.97	0.07	-3.19	15,18,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.