



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 16, 2017 – 06:53 PM EST

PDB ID : 5UME
Title : Crystal Structure of 5,10-Methylenetetrahydrofolate Reductase MetF from Haemophilus influenzae
Authors : Kim, Y.; Mulligan, R.; Maltseva, N.; Grimshaw, S.; Anderson, W.F.; Joachimiak, A.; Center for Structural Genomics of Infectious Diseases (CS-GID)
Deposited on : 2017-01-27
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-20028442
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-20028442

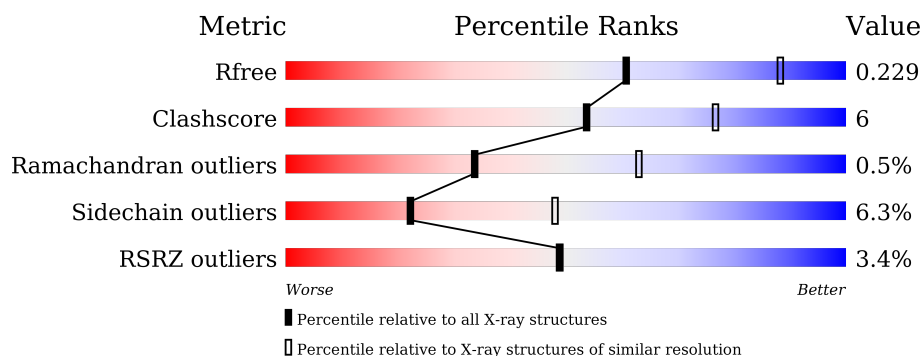
1 Overall quality at a glance

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	295	<div> <div>4%</div> <div> <div></div> <div>78%</div> <div>17%</div> <div>• •</div> </div> </div>
1	B	295	<div> <div>5%</div> <div> <div></div> <div>71%</div> <div>15%</div> <div>• 11%</div> </div> </div>
1	C	295	<div> <div>%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div>• •</div> </div> </div>
1	D	295	<div> <div>4%</div> <div> <div></div> <div>76%</div> <div>19%</div> <div>• •</div> </div> </div>
1	E	295	<div> <div>4%</div> <div> <div></div> <div>81%</div> <div>16%</div> <div>• •</div> </div> </div>
1	F	295	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>18%</div> <div>• •</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	ACY	D	302	-	-	-	X
5	EDO	F	303	-	-	-	X
6	FMT	B	303	-	-	-	X
6	FMT	C	301	-	-	-	X
6	FMT	C	302	-	-	-	X
6	FMT	D	304	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 13964 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 5,10-methylenetetrahydrofolate reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	0	0	0
			2276	1456	393	420	7			
1	B	262	Total	C	N	O	S	0	0	0
			2092	1341	361	383	7			
1	C	286	Total	C	N	O	S	0	1	0
			2286	1461	396	421	8			
1	D	289	Total	C	N	O	S	0	0	0
			2302	1472	399	424	7			
1	E	292	Total	C	N	O	S	0	0	0
			2323	1484	402	429	8			
1	F	282	Total	C	N	O	S	0	1	0
			2253	1442	391	413	7			

There are 18 discrepancies between the modelled and reference sequences:

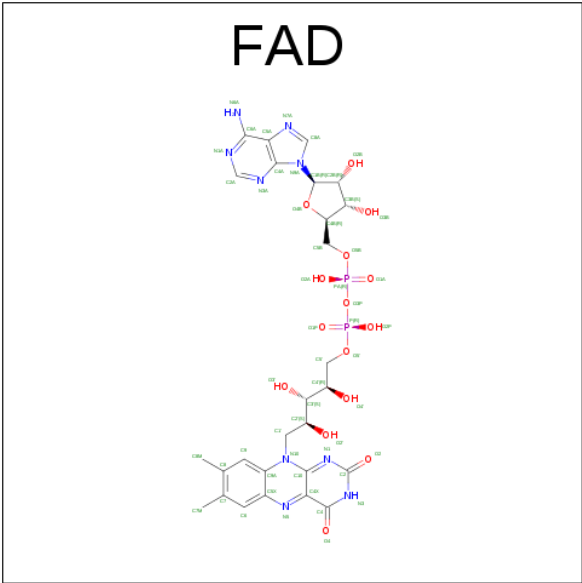
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP P45208
A	-1	ASN	-	expression tag	UNP P45208
A	0	ALA	-	expression tag	UNP P45208
B	-2	SER	-	expression tag	UNP P45208
B	-1	ASN	-	expression tag	UNP P45208
B	0	ALA	-	expression tag	UNP P45208
C	-2	SER	-	expression tag	UNP P45208
C	-1	ASN	-	expression tag	UNP P45208
C	0	ALA	-	expression tag	UNP P45208
D	-2	SER	-	expression tag	UNP P45208
D	-1	ASN	-	expression tag	UNP P45208
D	0	ALA	-	expression tag	UNP P45208
E	-2	SER	-	expression tag	UNP P45208
E	-1	ASN	-	expression tag	UNP P45208
E	0	ALA	-	expression tag	UNP P45208
F	-2	SER	-	expression tag	UNP P45208
F	-1	ASN	-	expression tag	UNP P45208

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Chain	Residue	Modelled	Actual	Comment	Reference
F	0	ALA	-	expression tag	UNP P45208

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



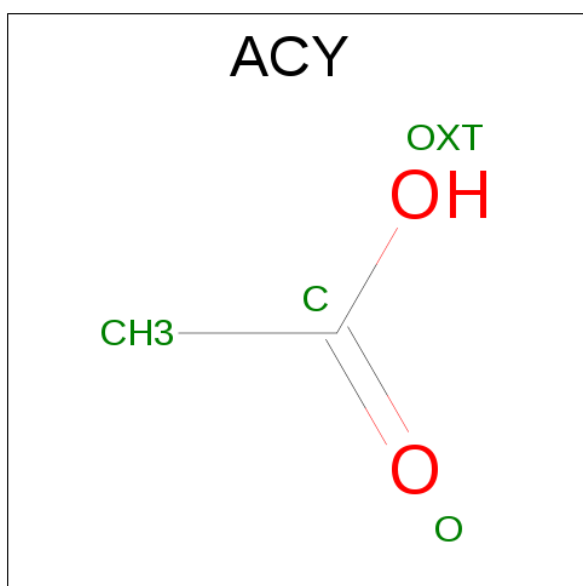
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	E	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	F	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is ACETIC ACID (three-letter code: ACY) (formula: C₂H₄O₂).



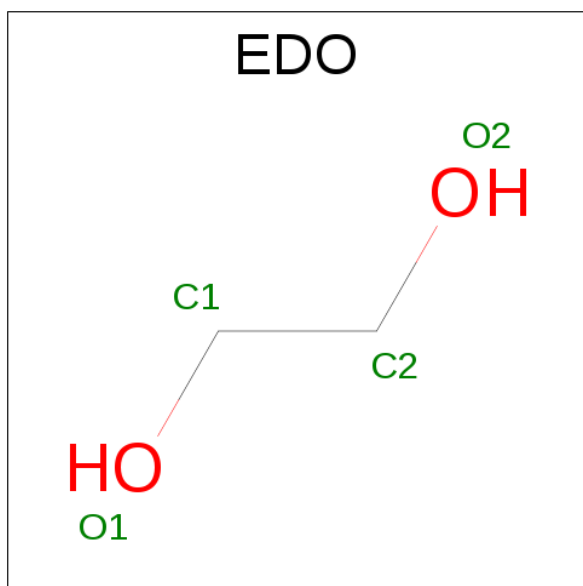
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		

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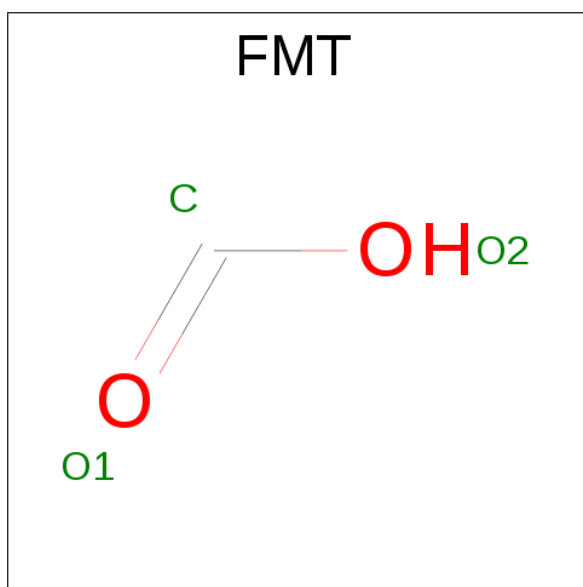
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	E	1	Total	C	O	0	0
			4	2	2		
4	F	1	Total	C	O	0	0
			4	2	2		
4	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is FORMIC ACID (three-letter code: FMT) (formula: CH₂O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			3	1	2		
6	C	1	Total	C	O	0	0
			3	1	2		
6	C	1	Total	C	O	0	0
			3	1	2		
6	D	1	Total	C	O	0	0
			3	1	2		

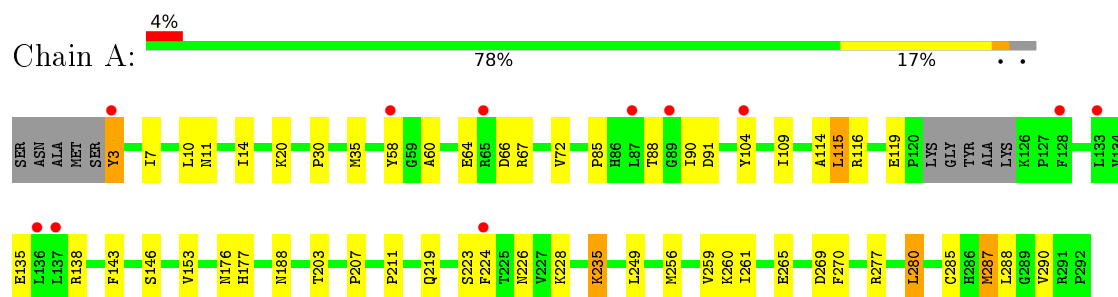
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	10	Total	O	0	0
			10	10		
7	B	4	Total	O	0	0
			4	4		
7	C	8	Total	O	0	0
			8	8		
7	D	8	Total	O	0	0
			8	8		
7	E	9	Total	O	0	0
			9	9		
7	F	9	Total	O	0	0
			9	9		

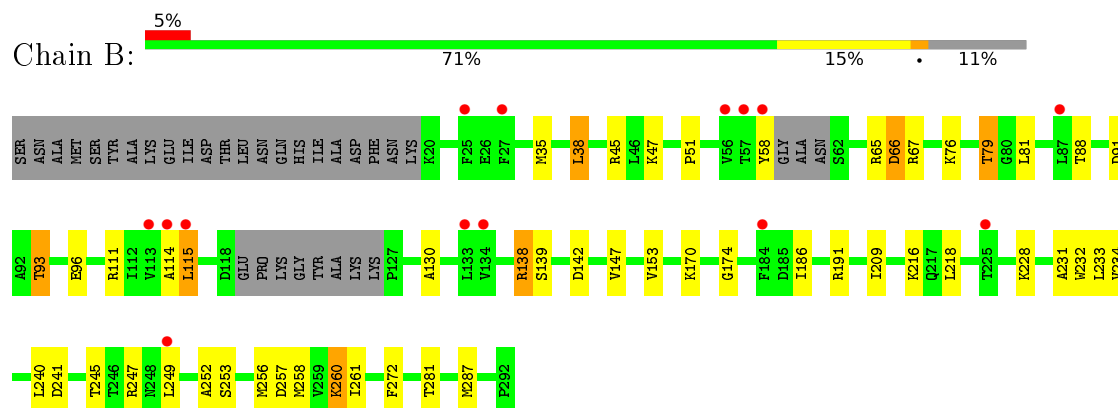
3 Residue-property plots [i](#)

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

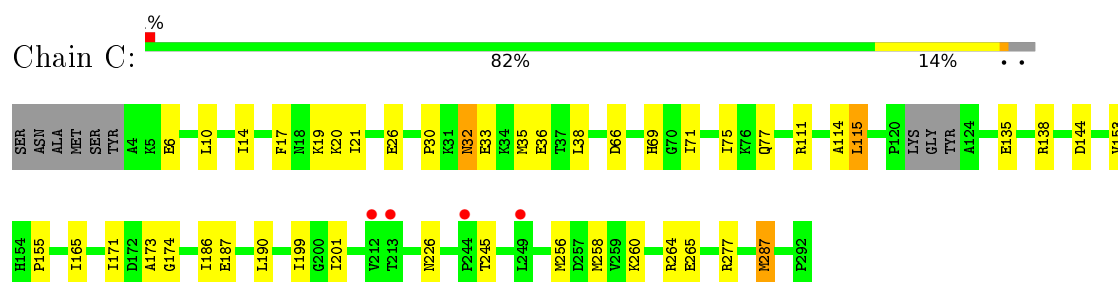
- Molecule 1: 5,10-methylenetetrahydrofolate reductase



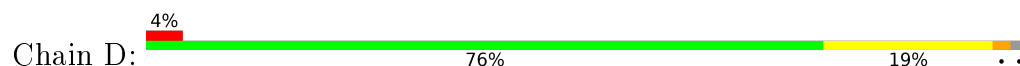
- Molecule 1: 5,10-methylenetetrahydrofolate reductase

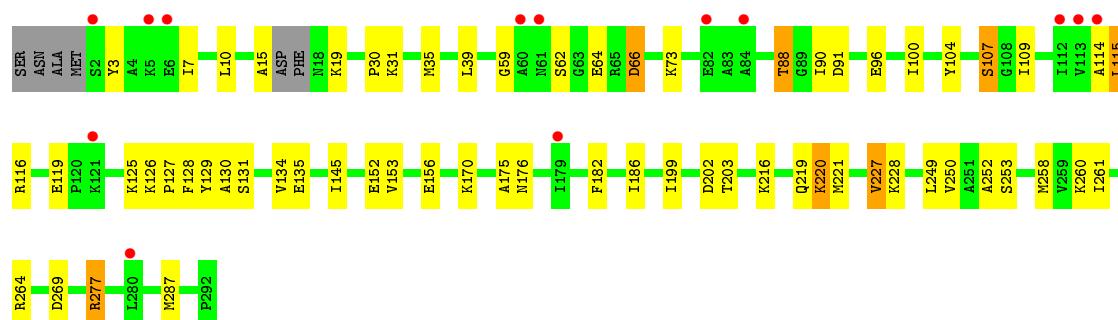


- Molecule 1: 5,10-methylenetetrahydrofolate reductase

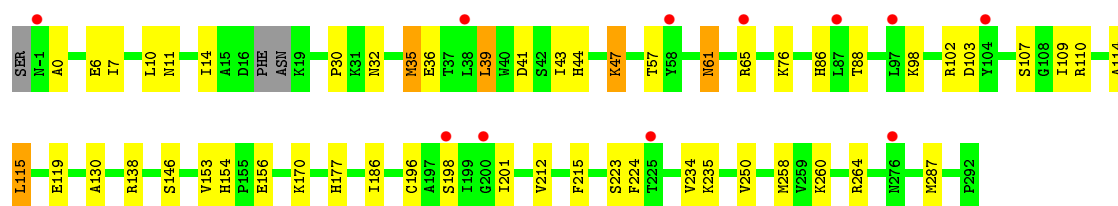
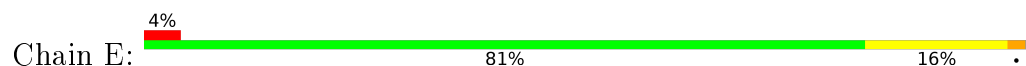


- Molecule 1: 5,10-methylenetetrahydrofolate reductase

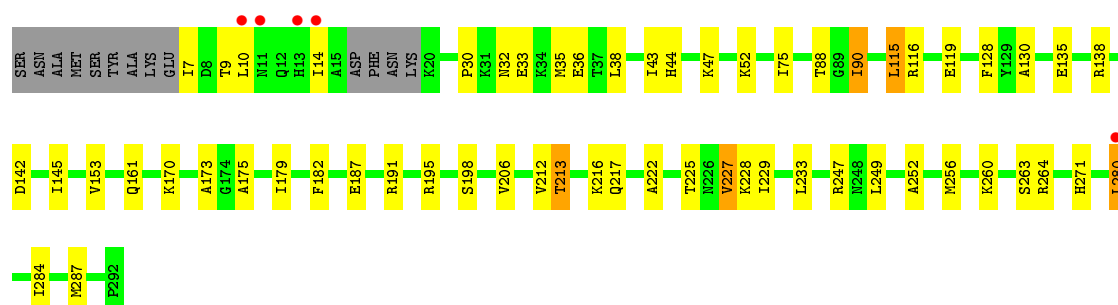
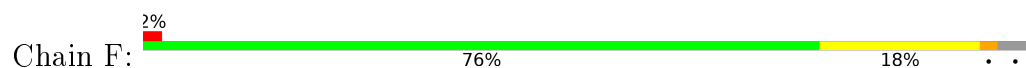




- Molecule 1: 5,10-methylenetetrahydrofolate reductase



- Molecule 1: 5,10-methylenetetrahydrofolate reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	90.06Å 134.96Å 182.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.73 – 2.70 47.30 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.5 (43.73-2.70) 98.6 (47.30-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 2.69Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, R_{free}	0.169 , 0.228 0.168 , 0.229	Depositor DCC
R_{free} test set	3022 reflections (4.95%)	DCC
Wilson B-factor (Å ²)	74.8	Xtriage
Anisotropy	0.354	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 48.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13964	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.91% 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: ACY, FMT, EDO, FAD, SO4

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	0.43	0/2326	0.59	0/3151
1	B	0.42	0/2137	0.56	0/2893
1	C	0.47	0/2335	0.60	0/3161
1	D	0.44	0/2352	0.59	0/3184
1	E	0.43	0/2373	0.57	0/3212
1	F	0.43	0/2302	0.59	0/3117
All	All	0.44	0/13825	0.58	0/18718

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	143	PHE	Peptide

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	2276	0	2277	34	0
1	B	2092	0	2102	26	0
1	C	2286	0	2294	26	0
1	D	2302	0	2312	44	0
1	E	2323	0	2333	31	0
1	F	2253	0	2267	33	0
2	A	53	0	31	2	0
2	B	53	0	31	1	0
2	C	53	0	31	2	0
2	D	53	0	31	2	0
2	E	53	0	31	4	0
2	F	53	0	31	2	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	4	0	3	0	0
4	B	4	0	3	0	0
4	C	4	0	3	1	0
4	D	12	0	9	0	0
4	E	4	0	3	0	0
4	F	8	0	6	0	0
5	A	4	0	6	0	0
5	F	4	0	6	1	0
6	B	3	0	1	0	0
6	C	6	0	2	0	0
6	D	3	0	1	1	0
7	A	10	0	0	1	0
7	B	4	0	0	0	0
7	C	8	0	0	2	0
7	D	8	0	0	0	0
7	E	9	0	0	1	0
7	F	9	0	0	0	0
All	All	13964	0	13814	180	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 6.

All (180) 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:D:116:ARG:HH22	1:D:125:LYS:HD3	1.50	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:161:GLN:OE1	1:F:195:ARG:NH2	2.22	0.72
1:C:26:GLU:OE2	7:C:401:HOH:O	2.10	0.69
1:E:11:ASN:HD21	1:E:260:LYS:HE3	1.60	0.67
1:C:138:ARG:HD3	1:C:174:GLY:HA3	1.78	0.66
4:C:303:ACY:O	7:C:401:HOH:O	2.14	0.66
1:A:10:LEU:HD13	1:D:3:TYR:HE2	1.64	0.63
1:B:249:LEU:HD22	1:F:252:ALA:HB2	1.82	0.62
1:A:235:LYS:HE2	1:D:15:ALA:HB1	1.82	0.61
1:A:176:ASN:O	1:A:203:THR:HG21	2.01	0.61
1:D:30:PRO:HB3	1:D:35:MET:HG3	1.81	0.61
1:A:14:ILE:HG13	1:A:260:LYS:HG2	1.83	0.60
1:B:261:ILE:HD11	1:F:7:ILE:HG12	1.84	0.60
1:A:10:LEU:HD13	1:D:3:TYR:CE2	2.37	0.59
1:F:52:LYS:HD2	5:F:303:EDO:H11	1.83	0.59
1:A:226:ASN:ND2	1:A:226:ASN:O	2.36	0.59
1:B:58:TYR:CD2	1:B:88:THR:HG22	2.37	0.59
1:A:104:TYR:CD1	1:A:109:ILE:HD12	2.38	0.58
1:F:10:LEU:HD11	1:F:260:LYS:HB3	1.86	0.58
1:D:115:LEU:HD13	2:D:300:FAD:C4X	2.34	0.57
1:D:88:THR:HG22	1:D:91:ASP:OD2	2.04	0.57
1:D:216:LYS:HA	1:D:219:GLN:HE21	1.70	0.57
1:A:207:PRO:HD2	1:A:269:ASP:O	2.04	0.57
1:D:66:ASP:OD1	1:D:66:ASP:N	2.39	0.56
1:E:11:ASN:ND2	1:E:260:LYS:HE3	2.21	0.56
1:E:88:THR:HA	1:E:115:LEU:O	2.05	0.56
1:A:115:LEU:HD13	2:A:300:FAD:C4X	2.36	0.56
1:E:98:LYS:O	1:E:102:ARG:HG2	2.05	0.56
1:D:90:ILE:HG12	1:D:116:ARG:O	2.06	0.55
1:F:256:MET:SD	1:F:287:MET:HG3	2.47	0.55
1:B:252:ALA:HB2	1:F:249:LEU:HD22	1.87	0.55
1:C:190:LEU:HD22	1:C:265:GLU:HG2	1.89	0.55
1:B:35:MET:HA	1:B:38:LEU:HD23	1.89	0.54
1:B:138:ARG:NH2	1:B:142:ASP:OD1	2.38	0.54
1:D:107:SER:HB3	1:D:109:ILE:HD12	1.90	0.54
1:C:35:MET:HE3	1:C:38:LEU:HB3	1.89	0.54
1:F:138:ARG:HH12	1:F:142:ASP:HA	1.73	0.54
1:E:39:LEU:HD22	1:E:43:ILE:HD11	1.90	0.53
1:F:33:GLU:H	1:F:33:GLU:CD	2.11	0.53
1:C:14:ILE:HG13	1:C:260:LYS:HG2	1.89	0.53
1:F:90:ILE:HB	1:F:116:ARG:O	2.09	0.53
1:E:61:ASN:ND2	1:E:61:ASN:O	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:176:ASN:O	1:D:203:THR:HG21	2.09	0.53
1:A:138:ARG:NH1	7:A:402:HOH:O	2.41	0.52
1:D:90:ILE:HG21	1:D:119:GLU:HG3	1.91	0.52
1:D:186:ILE:HD11	1:D:258:MET:HA	1.90	0.52
1:D:10:LEU:HD21	1:D:260:LYS:HB3	1.92	0.52
1:D:96:GLU:O	1:D:100:ILE:HG12	2.09	0.52
1:A:249:LEU:HD23	1:D:249:LEU:HD23	1.91	0.52
1:B:115:LEU:HD13	2:B:300:FAD:C4X	2.40	0.52
1:C:115:LEU:HD22	2:C:300:FAD:C9A	2.40	0.52
1:F:213:THR:O	1:F:247:ARG:HD2	2.09	0.52
1:A:211:PRO:HB3	1:A:280:LEU:HD12	1.92	0.52
1:C:10:LEU:O	1:C:14:ILE:HG12	2.09	0.52
2:E:300:FAD:O1P	7:E:401:HOH:O	2.18	0.51
1:B:186:ILE:HD11	1:B:258:MET:HG2	1.93	0.51
1:D:130:ALA:HB3	1:D:170:LYS:HD3	1.92	0.51
1:C:135:GLU:HG3	1:C:173:ALA:HB1	1.92	0.50
1:F:130:ALA:HB3	1:F:170:LYS:HD3	1.93	0.50
1:A:249:LEU:HD22	1:D:252:ALA:HB2	1.92	0.50
1:A:11:ASN:ND2	1:A:260:LYS:HD3	2.27	0.50
1:B:58:TYR:HD2	1:B:88:THR:HG22	1.76	0.50
1:E:57:THR:HA	1:E:86:HIS:CG	2.47	0.50
1:C:165:ILE:HD13	1:C:199:ILE:HD11	1.93	0.50
1:D:126:LYS:O	1:D:128:PHE:N	2.40	0.50
1:E:154:HIS:CE1	1:E:156:GLU:HB2	2.47	0.49
1:B:240:LEU:O	1:B:247:ARG:NH1	2.42	0.49
1:E:115:LEU:HD22	2:E:300:FAD:C9A	2.42	0.49
1:E:196:CYS:HB3	1:E:201:ILE:HD12	1.94	0.49
1:C:186:ILE:HD11	1:C:258:MET:HG2	1.95	0.49
1:F:30:PRO:HG2	1:F:36:GLU:HG2	1.94	0.49
1:F:116:ARG:HD2	1:F:128:PHE:O	2.13	0.49
1:B:147:VAL:HG22	1:B:170:LYS:HE2	1.95	0.48
1:B:66:ASP:OD1	1:B:66:ASP:N	2.45	0.48
1:B:138:ARG:HD2	1:B:174:GLY:HA3	1.94	0.48
1:E:30:PRO:HG2	1:E:36:GLU:HG2	1.96	0.48
1:C:115:LEU:HD13	2:C:300:FAD:C4X	2.45	0.47
1:D:114:ALA:C	1:D:115:LEU:HG	2.34	0.47
1:E:215:PHE:HZ	1:E:234:VAL:HG22	1.79	0.47
1:E:76:LYS:HG2	1:E:76:LYS:H	1.55	0.47
1:D:125:LYS:NZ	1:D:156:GLU:OE2	2.38	0.47
1:A:85:PRO:HD3	1:A:109:ILE:HD13	1.97	0.47
1:D:59:GLY:HA3	1:D:62:SER:OG	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:285:CYS:HB3	1:A:290:VAL:HB	1.96	0.47
1:A:91:ASP:OD2	1:A:91:ASP:N	2.48	0.47
1:C:30:PRO:HB3	1:C:35:MET:HG3	1.97	0.46
1:A:259:VAL:HG13	1:A:288:LEU:HD21	1.97	0.46
1:B:191:ARG:HD2	1:C:265:GLU:OE1	2.15	0.46
1:B:232:TRP:CZ3	1:B:233:LEU:HD13	2.50	0.46
1:C:17:PHE:HD1	1:C:21:ILE:HD11	1.80	0.46
1:A:256:MET:SD	1:A:287:MET:SD	3.14	0.46
1:B:191:ARG:HH22	1:C:264:ARG:NH1	2.13	0.46
1:F:229:ILE:HG23	1:F:233:LEU:HD23	1.96	0.46
1:C:17:PHE:CD1	1:C:21:ILE:HD11	2.50	0.46
1:C:32:ASN:OD1	1:C:32:ASN:N	2.49	0.46
1:F:216:LYS:HE3	1:F:216:LYS:HB3	1.78	0.46
1:E:109:ILE:O	1:E:110:ARG:NH2	2.48	0.45
1:E:32:ASN:OD1	1:E:35:MET:HB2	2.16	0.45
1:F:145:ILE:HB	1:F:175:ALA:HA	1.98	0.45
1:E:115:LEU:HD13	2:E:300:FAD:C4X	2.46	0.45
1:C:111:ARG:HA	1:C:144:ASP:O	2.16	0.45
1:A:3:TYR:CE2	1:D:264:ARG:HG2	2.52	0.45
1:B:231:ALA:O	1:B:234:VAL:HG12	2.16	0.45
1:C:287[B]:MET:SD	1:E:250:VAL:HG22	2.57	0.45
1:E:10:LEU:O	1:E:14:ILE:HG12	2.16	0.45
1:A:287:MET:SD	1:D:250:VAL:HG22	2.56	0.45
1:F:10:LEU:HD22	1:F:264:ARG:NH1	2.31	0.45
1:B:272:PHE:CD2	1:B:281:THR:HG23	2.51	0.45
1:D:269:ASP:HB3	6:D:304:FMT:H	1.98	0.45
1:E:10:LEU:HD11	1:E:260:LYS:O	2.17	0.44
1:B:93:THR:HG23	1:B:96:GLU:OE2	2.18	0.44
1:F:187:GLU:OE2	1:F:191:ARG:NH1	2.50	0.44
1:A:259:VAL:HG22	1:A:270:PHE:CE1	2.52	0.44
1:A:261:ILE:CD1	1:D:7:ILE:HG12	2.47	0.44
1:E:130:ALA:HB3	1:E:170:LYS:HD3	1.99	0.44
1:E:235:LYS:O	1:E:235:LYS:HG3	2.17	0.44
1:F:30:PRO:HB3	1:F:35:MET:HG3	2.00	0.44
1:A:88:THR:HA	1:A:115:LEU:O	2.18	0.44
1:B:130:ALA:HB3	1:B:170:LYS:HE3	2.00	0.44
1:D:152:GLU:OE2	1:D:228:LYS:HE3	2.18	0.44
1:D:277:ARG:HE	1:D:277:ARG:HB3	1.26	0.44
1:E:114:ALA:C	1:E:115:LEU:HG	2.38	0.44
1:E:186:ILE:HD11	1:E:258:MET:HG2	2.00	0.44
1:F:179:ILE:HG21	1:F:271:HIS:CD2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:265:GLU:OE1	1:E:0:ALA:HB3	2.18	0.44
1:A:90:ILE:HB	1:A:116:ARG:O	2.17	0.43
1:D:104:TYR:HD1	1:D:109:ILE:HD13	1.82	0.43
1:A:72:VAL:HG11	1:A:109:ILE:HD11	2.00	0.43
1:A:90:ILE:HD12	1:A:90:ILE:HA	1.88	0.43
1:F:32:ASN:HB2	1:F:33:GLU:OE1	2.19	0.43
1:C:71:ILE:O	1:C:75:ILE:HG13	2.19	0.43
1:D:73:LYS:HE2	1:D:73:LYS:HB2	1.70	0.43
1:A:228:LYS:HB2	1:A:228:LYS:HE3	1.72	0.43
1:C:171:ILE:HG13	1:C:201:ILE:HG23	2.00	0.43
1:D:115:LEU:HD13	2:D:300:FAD:N5	2.34	0.43
1:D:131:SER:O	1:D:135:GLU:HG3	2.19	0.43
1:C:256:MET:SD	1:C:287[B]:MET:SD	3.17	0.42
1:F:43:ILE:HD13	1:F:75:ILE:HG12	2.00	0.42
1:E:103:ASP:O	1:E:107:SER:HB2	2.18	0.42
1:E:146:SER:HA	1:E:177:HIS:O	2.18	0.42
1:F:135:GLU:HG2	1:F:173:ALA:HB1	2.00	0.42
1:F:88:THR:HA	1:F:115:LEU:O	2.19	0.42
1:D:126:LYS:C	1:D:128:PHE:H	2.22	0.42
1:B:51:PRO:HD2	1:B:81:LEU:HD21	2.00	0.42
1:D:220:LYS:HG2	1:D:221:MET:N	2.35	0.42
1:F:280:LEU:O	1:F:284:ILE:HG13	2.20	0.42
1:A:114:ALA:C	1:A:115:LEU:HG	2.40	0.42
1:A:146:SER:HA	1:A:177:HIS:HB3	2.02	0.42
1:C:33:GLU:HA	1:C:36:GLU:HB2	2.01	0.42
1:D:116:ARG:HD2	1:D:129:TYR:HA	2.01	0.42
1:F:222:ALA:HA	1:F:225:THR:OG1	2.20	0.42
1:B:114:ALA:C	1:B:115:LEU:HG	2.40	0.41
1:D:10:LEU:HD13	1:D:264:ARG:HB2	2.01	0.41
1:F:44:HIS:HA	1:F:47:LYS:HD3	2.02	0.41
1:A:7:ILE:HG13	1:D:261:ILE:HD11	2.01	0.41
1:B:216:LYS:NZ	1:B:241:ASP:OD1	2.47	0.41
1:D:130:ALA:O	1:D:134:VAL:HG23	2.19	0.41
1:F:115:LEU:HD22	2:F:300:FAD:C9A	2.50	0.41
1:F:14:ILE:HD11	1:F:260:LYS:HG2	2.02	0.41
1:B:88:THR:HG23	1:B:91:ASP:OD2	2.20	0.41
1:F:182:PHE:CE1	1:F:227:VAL:HG11	2.54	0.41
1:E:30:PRO:HB2	1:E:35:MET:HB3	2.03	0.41
1:A:64:GLU:O	1:A:67:ARG:HB3	2.20	0.41
1:F:115:LEU:HD13	2:F:300:FAD:C4X	2.50	0.41
1:E:264:ARG:HD2	1:E:264:ARG:HA	1.77	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:104:TYR:CD1	1:D:109:ILE:HD13	2.56	0.41
1:D:30:PRO:HD3	1:D:39:LEU:HD22	2.03	0.41
1:E:44:HIS:HA	1:E:47:LYS:HE2	2.03	0.41
1:F:90:ILE:HD13	1:F:90:ILE:HA	1.77	0.41
1:A:30:PRO:HG2	1:A:67:ARG:NH2	2.35	0.40
1:D:182:PHE:CE1	1:D:227:VAL:HG11	2.57	0.40
1:A:115:LEU:HD13	2:A:300:FAD:N5	2.35	0.40
1:B:47:LYS:HD3	1:B:79:THR:HB	2.02	0.40
1:C:155:PRO:HG3	1:C:226:ASN:HB2	2.04	0.40
1:D:145:ILE:HB	1:D:175:ALA:HA	2.03	0.40
1:D:62:SER:C	1:D:64:GLU:H	2.24	0.40
1:B:256:MET:O	1:B:260:LYS:HB2	2.21	0.40
1:E:10:LEU:HD21	1:E:260:LYS:HB3	2.03	0.40
1:E:115:LEU:HD13	2:E:300:FAD:N5	2.37	0.40
1:C:114:ALA:C	1:C:115:LEU:HG	2.41	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	281/295 (95%)	274 (98%)	5 (2%)	2 (1%)	26	55
1	B	256/295 (87%)	246 (96%)	9 (4%)	1 (0%)	39	69
1	C	283/295 (96%)	272 (96%)	10 (4%)	1 (0%)	39	69
1	D	285/295 (97%)	275 (96%)	8 (3%)	2 (1%)	26	55
1	E	288/295 (98%)	277 (96%)	10 (4%)	1 (0%)	46	75
1	F	279/295 (95%)	269 (96%)	9 (3%)	1 (0%)	39	69
All	All	1672/1770 (94%)	1613 (96%)	51 (3%)	8 (0%)	34	63

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	153	VAL
1	B	153	VAL
1	D	153	VAL
1	F	153	VAL
1	C	153	VAL
1	E	153	VAL
1	A	60	ALA
1	D	127	PRO

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	246/253 (97%)	229 (93%)	17 (7%)	19	43
1	B	227/253 (90%)	207 (91%)	20 (9%)	12	28
1	C	247/253 (98%)	234 (95%)	13 (5%)	28	57
1	D	248/253 (98%)	235 (95%)	13 (5%)	29	58
1	E	250/253 (99%)	234 (94%)	16 (6%)	22	47
1	F	243/253 (96%)	229 (94%)	14 (6%)	25	52
All	All	1461/1518 (96%)	1368 (94%)	93 (6%)	22	47

All (93) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	TYR
1	A	20	LYS
1	A	35	MET
1	A	58	TYR
1	A	66	ASP
1	A	115	LEU
1	A	119	GLU
1	A	135	GLU
1	A	188	ASN
1	A	219	GLN
1	A	223	SER

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Mol	Chain	Res	Type
1	A	224	PHE
1	A	235	LYS
1	A	265	GLU
1	A	277	ARG
1	A	280	LEU
1	A	287	MET
1	B	38	LEU
1	B	45	ARG
1	B	65	ARG
1	B	66	ASP
1	B	67	ARG
1	B	76	LYS
1	B	79	THR
1	B	93	THR
1	B	111	ARG
1	B	115	LEU
1	B	138	ARG
1	B	139	SER
1	B	209	ILE
1	B	218	LEU
1	B	228	LYS
1	B	245	THR
1	B	253	SER
1	B	257	ASP
1	B	260	LYS
1	B	287	MET
1	C	6	GLU
1	C	19	LYS
1	C	20	LYS
1	C	32	ASN
1	C	66	ASP
1	C	69	HIS
1	C	77	GLN
1	C	115	LEU
1	C	187	GLU
1	C	245	THR
1	C	277	ARG
1	C	287[A]	MET
1	C	287[B]	MET
1	D	19	LYS
1	D	31	LYS
1	D	66	ASP

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Mol	Chain	Res	Type
1	D	88	THR
1	D	107	SER
1	D	115	LEU
1	D	199	ILE
1	D	202	ASP
1	D	220	LYS
1	D	227	VAL
1	D	253	SER
1	D	277	ARG
1	D	287	MET
1	E	6	GLU
1	E	7	ILE
1	E	35	MET
1	E	39	LEU
1	E	41	ASP
1	E	47	LYS
1	E	61	ASN
1	E	65	ARG
1	E	115	LEU
1	E	119	GLU
1	E	138	ARG
1	E	198	SER
1	E	212	VAL
1	E	223	SER
1	E	224	PHE
1	E	287	MET
1	F	9	THR
1	F	38	LEU
1	F	90	ILE
1	F	115	LEU
1	F	119	GLU
1	F	198	SER
1	F	206	VAL
1	F	212	VAL
1	F	213	THR
1	F	217	GLN
1	F	227	VAL
1	F	228	LYS
1	F	263	SER
1	F	280	LEU

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	11	ASN
1	A	18	ASN
1	A	77	GLN
1	B	217	GLN
1	C	248	ASN
1	D	188	ASN
1	D	219	GLN
1	E	11	ASN
1	E	61	ASN
1	F	44	HIS

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](#)

23 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	FAD	A	300	-	52,58,58	0.99	4 (7%)	52,89,89	2.48	5 (9%)
3	SO4	A	301	-	4,4,4	0.37	0	6,6,6	1.14	1 (16%)
4	ACY	A	302	-	0,3,3	0.00	-	0,3,3	0.00	-
5	EDO	A	303	-	3,3,3	0.70	0	2,2,2	0.78	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FAD	B	300	-	52,58,58	0.90	4 (7%)	52,89,89	2.44	5 (9%)
3	SO4	B	301	-	4,4,4	0.80	0	6,6,6	0.93	0
4	ACY	B	302	-	0,3,3	0.00	-	0,3,3	0.00	-
6	FMT	B	303	-	0,2,2	0.00	-	0,1,1	0.00	-
2	FAD	C	300	-	52,58,58	0.90	2 (3%)	52,89,89	2.43	5 (9%)
6	FMT	C	301	-	0,2,2	0.00	-	0,1,1	0.00	-
6	FMT	C	302	-	0,2,2	0.00	-	0,1,1	0.00	-
4	ACY	C	303	-	0,3,3	0.00	-	0,3,3	0.00	-
2	FAD	D	300	-	52,58,58	1.09	4 (7%)	52,89,89	2.44	5 (9%)
4	ACY	D	301	-	0,3,3	0.00	-	0,3,3	0.00	-
4	ACY	D	302	-	0,3,3	0.00	-	0,3,3	0.00	-
4	ACY	D	303	-	0,3,3	0.00	-	0,3,3	0.00	-
6	FMT	D	304	-	0,2,2	0.00	-	0,1,1	0.00	-
2	FAD	E	300	-	52,58,58	1.03	3 (5%)	52,89,89	2.43	5 (9%)
4	ACY	E	301	-	0,3,3	0.00	-	0,3,3	0.00	-
2	FAD	F	300	-	52,58,58	0.96	2 (3%)	52,89,89	2.47	6 (11%)
4	ACY	F	301	-	0,3,3	0.00	-	0,3,3	0.00	-
4	ACY	F	302	-	0,3,3	0.00	-	0,3,3	0.00	-
5	EDO	F	303	-	3,3,3	0.56	0	2,2,2	0.36	0

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	FAD	A	300	-	-	0/30/50/50	0/6/6/6
3	SO4	A	301	-	-	0/0/0/0	0/0/0/0
4	ACY	A	302	-	-	0/0/0/0	0/0/0/0
5	EDO	A	303	-	-	0/1/1/1	0/0/0/0
2	FAD	B	300	-	-	0/30/50/50	0/6/6/6
3	SO4	B	301	-	-	0/0/0/0	0/0/0/0
4	ACY	B	302	-	-	0/0/0/0	0/0/0/0
6	FMT	B	303	-	-	0/0/0/0	0/0/0/0
2	FAD	C	300	-	-	0/30/50/50	0/6/6/6
6	FMT	C	301	-	-	0/0/0/0	0/0/0/0
6	FMT	C	302	-	-	0/0/0/0	0/0/0/0
4	ACY	C	303	-	-	0/0/0/0	0/0/0/0
2	FAD	D	300	-	-	0/30/50/50	0/6/6/6
4	ACY	D	301	-	-	0/0/0/0	0/0/0/0
4	ACY	D	302	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ACY	D	303	-	-	0/0/0/0	0/0/0/0
6	FMT	D	304	-	-	0/0/0/0	0/0/0/0
2	FAD	E	300	-	-	0/30/50/50	0/6/6/6
4	ACY	E	301	-	-	0/0/0/0	0/0/0/0
2	FAD	F	300	-	-	0/30/50/50	0/6/6/6
4	ACY	F	301	-	-	0/0/0/0	0/0/0/0
4	ACY	F	302	-	-	0/0/0/0	0/0/0/0
5	EDO	F	303	-	-	0/1/1/1	0/0/0/0

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	300	FAD	C10-N10	-2.20	1.36	1.39
2	F	300	FAD	C4X-N5	-2.03	1.30	1.33
2	A	300	FAD	C4X-N5	-2.01	1.30	1.33
2	D	300	FAD	C4X-C10	2.09	1.44	1.40
2	B	300	FAD	C9A-N10	2.18	1.41	1.38
2	A	300	FAD	C5X-N5	2.21	1.38	1.35
2	C	300	FAD	C4-C4X	2.23	1.45	1.41
2	D	300	FAD	C9A-N10	2.24	1.42	1.38
2	B	300	FAD	C5X-N5	2.24	1.38	1.35
2	E	300	FAD	C9A-N10	2.25	1.42	1.38
2	B	300	FAD	C4-C4X	2.35	1.46	1.41
2	B	300	FAD	C4-N3	2.69	1.37	1.33
2	E	300	FAD	C5X-N5	2.81	1.39	1.35
2	A	300	FAD	C4-N3	3.13	1.38	1.33
2	C	300	FAD	C4-N3	3.34	1.39	1.33
2	D	300	FAD	C5X-N5	3.48	1.40	1.35
2	E	300	FAD	C4-N3	3.58	1.39	1.33
2	F	300	FAD	C4-N3	3.61	1.39	1.33
2	D	300	FAD	C4-N3	4.07	1.40	1.33

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	300	FAD	C4X-C4-N3	-7.69	113.46	123.52
2	B	300	FAD	C4X-C4-N3	-7.42	113.83	123.52
2	F	300	FAD	C4X-C4-N3	-7.41	113.84	123.52
2	A	300	FAD	C4X-C4-N3	-7.40	113.85	123.52
2	C	300	FAD	C4X-C4-N3	-7.19	114.12	123.52
2	E	300	FAD	C4X-C4-N3	-7.19	114.12	123.52
2	F	300	FAD	C4X-C10-N10	-5.46	116.55	120.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	300	FAD	C4X-C10-N10	-5.31	116.66	120.52
2	E	300	FAD	N3-C2-N1	-5.20	118.94	127.69
2	B	300	FAD	C4X-C10-N10	-5.16	116.77	120.52
2	E	300	FAD	C4X-C10-N10	-5.15	116.78	120.52
2	A	300	FAD	C4X-C10-N10	-5.12	116.80	120.52
2	A	300	FAD	N3-C2-N1	-5.08	119.13	127.69
2	D	300	FAD	C4X-C10-N10	-4.97	116.91	120.52
2	D	300	FAD	N3-C2-N1	-4.97	119.32	127.69
2	C	300	FAD	N3-C2-N1	-4.94	119.36	127.69
2	F	300	FAD	N3-C2-N1	-4.72	119.74	127.69
2	B	300	FAD	N3-C2-N1	-4.55	120.02	127.69
2	A	300	FAD	C4-C4X-C10	-4.18	117.27	119.94
2	E	300	FAD	C4-C4X-C10	-4.06	117.35	119.94
2	C	300	FAD	C4-C4X-C10	-3.98	117.39	119.94
2	F	300	FAD	C4-C4X-C10	-3.93	117.42	119.94
2	B	300	FAD	C4-C4X-C10	-3.80	117.51	119.94
2	D	300	FAD	C4-C4X-C10	-3.73	117.56	119.94
3	A	301	SO4	O2-S-O1	-2.25	102.07	109.59
2	F	300	FAD	O4'-C4'-C5'	-2.07	105.58	110.09
2	E	300	FAD	C4-N3-C2	12.94	125.96	115.16
2	C	300	FAD	C4-N3-C2	13.00	126.01	115.16
2	D	300	FAD	C4-N3-C2	13.05	126.05	115.16
2	B	300	FAD	C4-N3-C2	13.20	126.17	115.16
2	F	300	FAD	C4-N3-C2	13.22	126.19	115.16
2	A	300	FAD	C4-N3-C2	13.38	126.32	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	300	FAD	2	0
2	B	300	FAD	1	0
2	C	300	FAD	2	0
4	C	303	ACY	1	0
2	D	300	FAD	2	0
6	D	304	FMT	1	0
2	E	300	FAD	4	0
2	F	300	FAD	2	0
5	F	303	EDO	1	0

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, 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	285/295 (96%)	0.11	11 (3%)	43 43	56, 80, 141, 182	0
1	B	262/295 (88%)	0.20	14 (5%)	30 28	56, 91, 133, 159	0
1	C	286/295 (96%)	0.07	4 (1%)	78 77	55, 76, 133, 164	0
1	D	289/295 (97%)	0.25	13 (4%)	37 36	56, 82, 138, 193	0
1	E	292/295 (98%)	0.29	11 (3%)	44 44	58, 91, 155, 192	0
1	F	282/295 (95%)	0.15	5 (1%)	71 72	53, 79, 121, 221	0
All	All	1696/1770 (95%)	0.18	58 (3%)	49 49	53, 82, 141, 221	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	58	TYR	6.9
1	D	2	SER	6.6
1	F	13	HIS	6.6
1	F	14	ILE	5.7
1	F	10	LEU	5.4
1	D	61	ASN	4.3
1	E	-1	ASN	4.1
1	D	121	LYS	4.1
1	B	114	ALA	3.9
1	D	5	LYS	3.8
1	A	224	PHE	3.8
1	D	60	ALA	3.8
1	A	89	GLY	3.8
1	E	104	TYR	3.7
1	D	113	VAL	3.5
1	E	65	ARG	3.5
1	E	58	TYR	3.3
1	D	112	ILE	3.3
1	F	280	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
1	E	276	ASN	3.0
1	A	128	PHE	3.0
1	E	87	LEU	2.9
1	D	6	GLU	2.9
1	A	3	TYR	2.8
1	B	87	LEU	2.8
1	A	65	ARG	2.8
1	E	97	LEU	2.7
1	B	115	LEU	2.7
1	A	137	LEU	2.6
1	D	179	ILE	2.5
1	C	249	LEU	2.5
1	B	133	LEU	2.5
1	B	56	VAL	2.5
1	E	200	GLY	2.4
1	A	87	LEU	2.4
1	B	225	THR	2.4
1	C	244	PRO	2.4
1	C	213	THR	2.3
1	E	225	THR	2.3
1	B	25	PHE	2.3
1	B	113	VAL	2.3
1	A	136	LEU	2.3
1	B	184	PHE	2.3
1	D	114	ALA	2.3
1	E	38	LEU	2.3
1	B	134	VAL	2.3
1	E	198	SER	2.3
1	C	212	VAL	2.2
1	D	84	ALA	2.2
1	F	11	ASN	2.2
1	D	82	GLU	2.2
1	A	58	TYR	2.2
1	A	104	TYR	2.2
1	B	57	THR	2.2
1	B	27	PHE	2.1
1	B	249	LEU	2.1
1	A	133	LEU	2.1
1	D	280	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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(Å ²)	Q<0.9
6	FMT	C	302	3/3	0.72	0.69	21.30	115,115,120,121	0
5	EDO	F	303	4/4	0.90	0.76	11.93	82,84,91,95	0
6	FMT	D	304	3/3	0.92	0.76	9.89	90,90,95,101	0
6	FMT	C	301	3/3	0.96	0.33	6.01	85,85,90,100	0
4	ACY	D	302	4/4	0.84	0.35	4.22	110,112,114,114	0
6	FMT	B	303	3/3	0.98	0.38	3.36	95,95,98,98	0
5	EDO	A	303	4/4	0.74	0.23	0.84	100,102,106,111	0
2	FAD	B	300	53/53	0.96	0.17	-0.56	72,87,105,107	0
2	FAD	F	300	53/53	0.97	0.16	-0.60	57,65,83,90	0
2	FAD	D	300	53/53	0.96	0.15	-0.64	48,63,90,97	0
2	FAD	C	300	53/53	0.98	0.13	-0.93	50,72,86,90	0
2	FAD	E	300	53/53	0.97	0.14	-0.93	44,64,89,94	0
2	FAD	A	300	53/53	0.98	0.12	-1.22	60,77,106,109	0
3	SO4	A	301	5/5	0.94	0.06	-1.29	144,144,146,146	0
4	ACY	A	302	4/4	0.87	0.28	-	87,97,99,101	0
4	ACY	C	303	4/4	0.93	0.19	-	96,99,103,104	0
3	SO4	B	301	5/5	0.84	0.16	-	157,161,164,167	0
4	ACY	F	301	4/4	0.94	0.29	-	84,85,86,87	0
4	ACY	B	302	4/4	0.90	0.22	-	93,96,98,100	0
4	ACY	F	302	4/4	0.84	0.25	-	114,116,118,120	0
4	ACY	E	301	4/4	0.92	0.16	-	89,89,90,91	0
4	ACY	D	303	4/4	0.89	0.22	-	58,75,76,81	0
4	ACY	D	301	4/4	0.91	0.11	-	90,102,102,105	0

6.5 Other polymers [i](#)

There are no such residues in this entry.