



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 2, 2019 – 11:02 AM EST

PDB ID : 6BPY
Title : Aspergillus fumigatus Thioredoxin Reductase
Authors : Marshall, A.C.; Bruning, J.B.
Deposited on : 2017-11-27
Resolution : 3.20 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 : rb-20031633
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) : rb-20031633

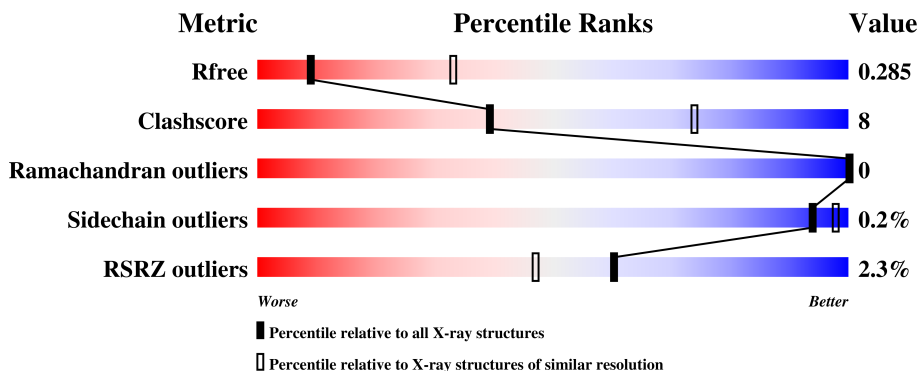
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1121 (3.22-3.18)
Clashscore	122126	1091 (3.20-3.20)
Ramachandran outliers	120053	1074 (3.20-3.20)
Sidechain outliers	120020	1073 (3.20-3.20)
RSRZ outliers	108989	1083 (3.22-3.18)

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	374	 67% 20% 13%
1	B	374	 72% 13% 15%

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
3	MLI	A	403	-	-	-	X
3	MLI	A	404	-	-	-	X
3	MLI	B	402	-	-	-	X
3	MLI	B	403	-	-	X	-
4	FMT	B	405	-	-	-	X
4	FMT	B	407	-	-	-	X

2 Entry composition i

There are 8 unique types of molecules in this entry. The entry contains 5003 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.

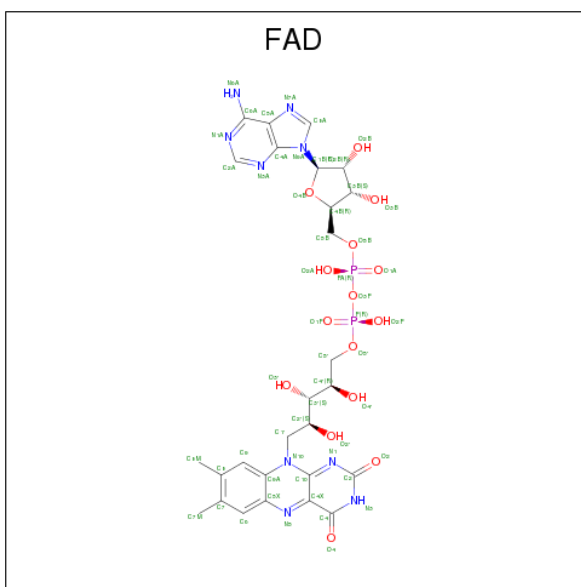
- Molecule 1 is a protein called Thioredoxin reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	324	Total	C	N	O	S	0	0	0
			2374	1487	413	464	10			
1	B	319	Total	C	N	O	S	0	0	0
			2347	1476	404	456	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP A0A229Y1X4
B	0	GLY	-	expression tag	UNP A0A229Y1X4

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



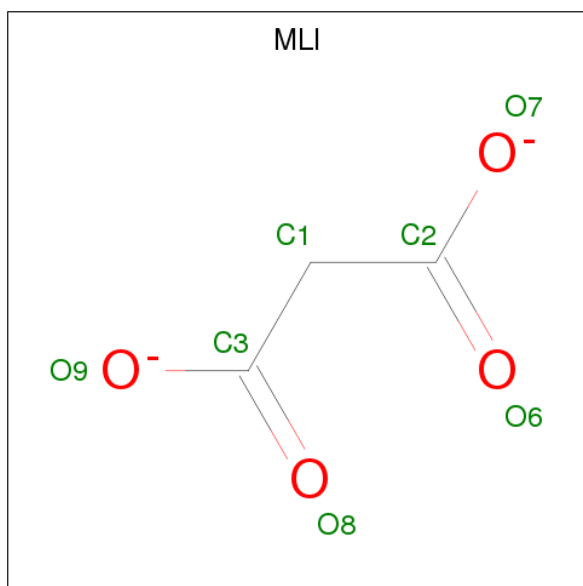
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

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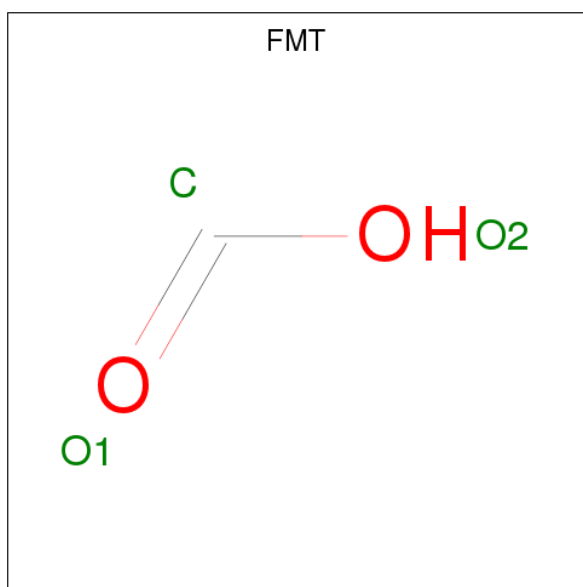
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	B	1	53	27	9	15	2	0	0

- Molecule 3 is MALONATE ION (three-letter code: MLI) (formula: $C_3H_2O_4$).



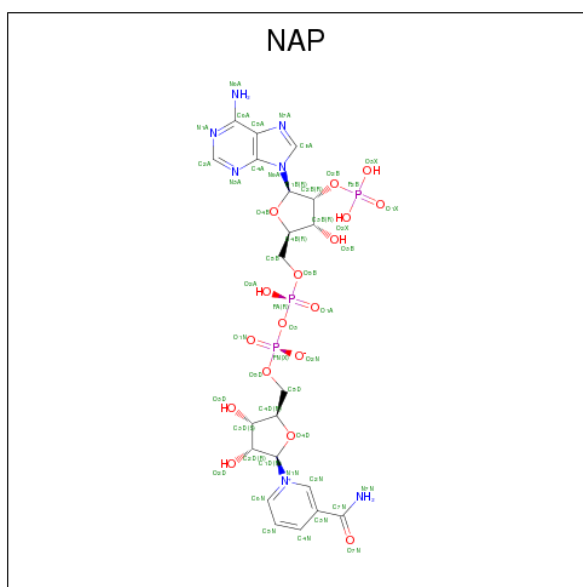
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	7	3	4	0	0
3	A	1	7	3	4	0	0
3	A	1	7	3	4	0	0
3	B	1	7	3	4	0	0
3	B	1	7	3	4	0	0

- Molecule 4 is FORMIC ACID (three-letter code: FMT) (formula: CH_2O_2).



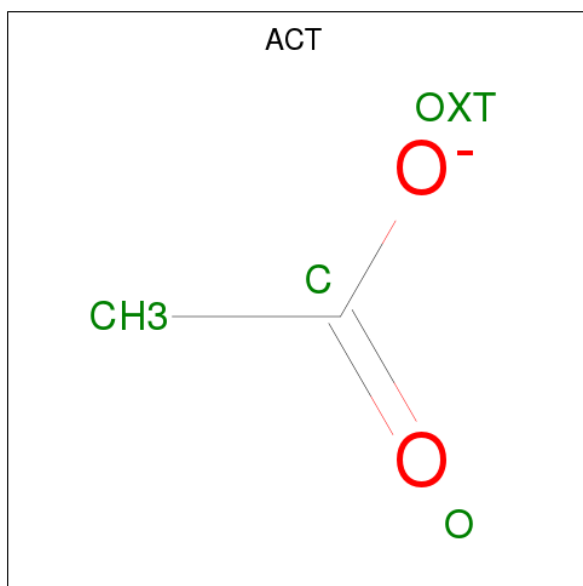
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 3 1 2	0	0
4	A	1	Total C O 3 1 2	0	0
4	B	1	Total C O 3 1 2	0	0
4	B	1	Total C O 3 1 2	0	0
4	B	1	Total C O 3 1 2	0	0
4	B	1	Total C O 3 1 2	0	0
4	B	1	Total C O 3 1 2	0	0

- Molecule 5 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



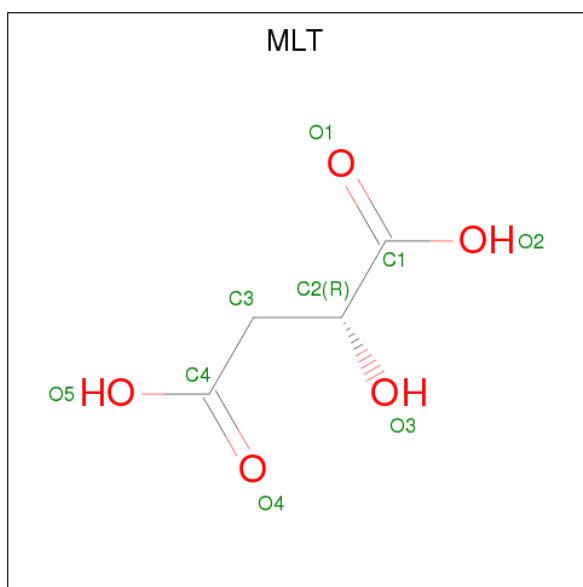
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	N	O			P
5	A	1	31	10	5	13	3	0	0
5	B	1	31	10	5	13	3	0	0

- Molecule 6 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 7 is D-MALATE (three-letter code: MLT) (formula: $C_4H_6O_5$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	B	1	Total C O 9 4 5	0	0

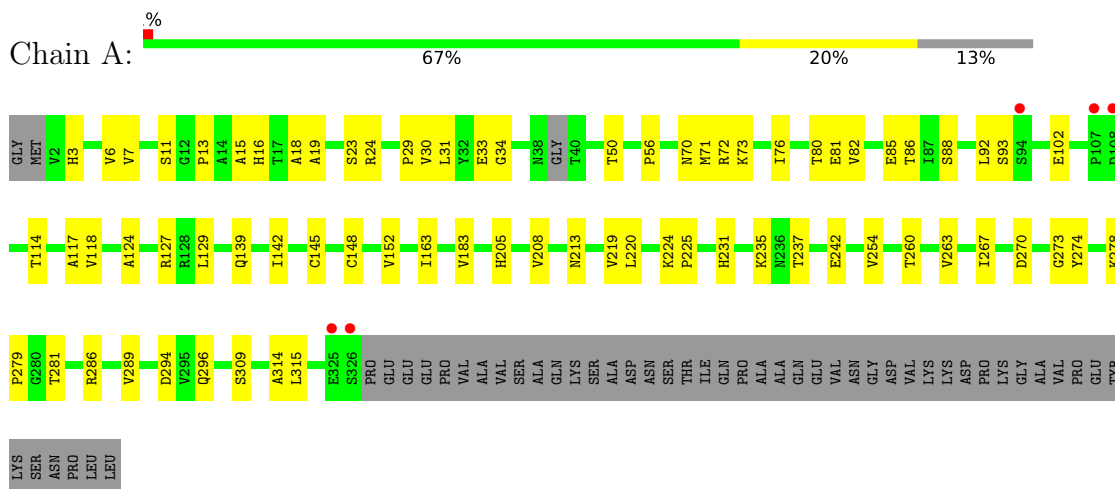
- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	22	Total O 22 22	0	0
8	B	23	Total O 23 23	0	0

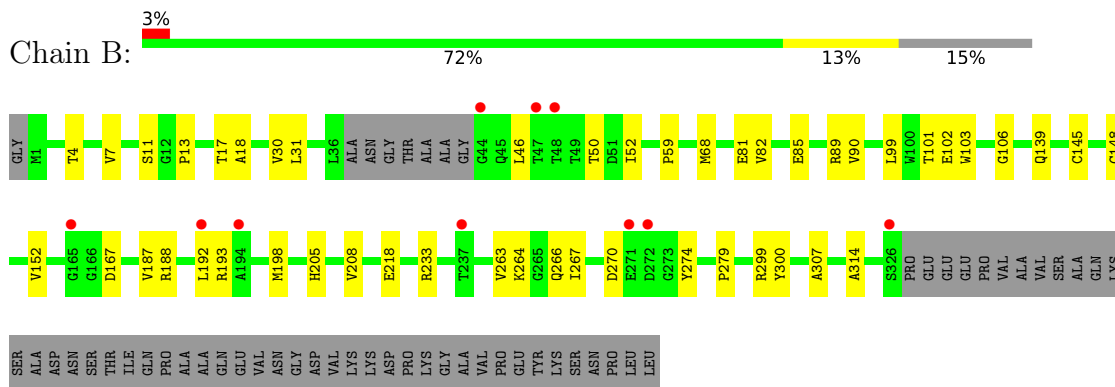
3 Residue-property plots [i](#)

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



- Molecule 1: Thioredoxin reductase



4 Data and refinement statistics i

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	159.22Å 159.22Å 121.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.97 – 3.20 29.97 – 3.20	Depositor EDS
% Data completeness (in resolution range)	95.3 (29.97-3.20) 79.8 (29.97-3.20)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.80 (at 3.18Å)	Xtrriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.231 , 0.261 0.266 , 0.285	Depositor DCC
R_{free} test set	1419 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	75.1	Xtrriage
Anisotropy	0.277	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 22.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.037 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	5003	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% 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: FMT, MLI, NAP, MLT, ACT, FAD

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.24	0/2414	0.45	0/3281
1	B	0.24	0/2389	0.43	0/3251
All	All	0.24	0/4803	0.44	0/6532

There are no bond length outliers.

There are no bond angle outliers.

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	2374	0	2301	46	0
1	B	2347	0	2276	36	0
2	A	53	0	30	0	0
2	B	53	0	30	3	0
3	A	21	0	6	1	0
3	B	14	0	4	2	0
4	A	6	0	2	0	0
4	B	15	0	5	0	0
5	A	31	0	11	0	0
5	B	31	0	11	2	0
6	A	4	0	3	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	9	0	4	0	0
8	A	22	0	0	0	0
8	B	23	0	0	0	0
All	All	5003	0	4683	77	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 (77) 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:279:PRO:HB2	1:B:279:PRO:HB2	1.74	0.70
1:B:89:ARG:NH2	1:B:106:GLY:O	2.27	0.67
1:A:92:LEU:O	1:A:286:ARG:NH1	2.28	0.65
1:B:30:VAL:HG12	1:B:81:GLU:HB3	1.79	0.64
1:A:19:ALA:HB1	1:A:80:THR:HG21	1.81	0.63
1:B:46:LEU:HD13	2:B:401:FAD:H2'	1.83	0.60
1:A:30:VAL:HG12	1:A:81:GLU:HB3	1.82	0.59
1:A:152:VAL:HG11	3:A:402:MLI:H11	1.84	0.59
1:A:270:ASP:OD1	1:A:274:TYR:N	2.39	0.56
1:B:90:VAL:HG12	1:B:99:LEU:HD22	1.88	0.55
1:B:18:ALA:HA	1:B:314:ALA:HB2	1.88	0.55
1:A:86:THR:HB	1:A:102:GLU:HG2	1.88	0.55
1:A:29:PRO:HG2	1:A:80:THR:HG22	1.86	0.55
1:A:73:LYS:HA	1:A:76:ILE:HG12	1.90	0.54
1:A:205:HIS:HB3	1:A:208:VAL:HG22	1.88	0.54
1:A:6:VAL:HG22	1:A:117:ALA:HB3	1.89	0.54
1:B:11:SER:H	2:B:401:FAD:H4B	1.72	0.54
1:B:13:PRO:HG3	1:B:68:MET:HE1	1.90	0.53
1:A:315:LEU:HD12	1:B:300:TYR:HD1	1.74	0.53
1:B:187:VAL:HG11	1:B:192:LEU:HD23	1.92	0.52
1:B:270:ASP:OD1	1:B:274:TYR:N	2.42	0.52
1:B:205:HIS:HB3	1:B:208:VAL:HG12	1.90	0.52
1:A:88:SER:HB2	1:A:102:GLU:HA	1.91	0.52
1:B:263:VAL:HG23	1:B:264:LYS:H	1.74	0.52
1:A:183:VAL:HB	1:A:208:VAL:HG12	1.92	0.51
1:A:18:ALA:HA	1:A:314:ALA:HB2	1.91	0.51
1:B:31:LEU:HB3	1:B:82:VAL:HG22	1.93	0.50
1:B:13:PRO:HD2	2:B:401:FAD:H5'2	1.93	0.49
1:B:167:ASP:HA	1:B:198:MET:HE1	1.93	0.49
1:B:193:ARG:NH1	5:B:409:NAP:O2X	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:LEU:HD11	1:A:163:ILE:HG21	1.95	0.48
1:A:224:LYS:HB2	1:A:225:PRO:HD2	1.96	0.48
1:B:188:ARG:NH2	5:B:409:NAP:O1X	2.36	0.47
1:A:296:GLN:NE2	1:A:309:SER:OG	2.47	0.47
1:B:46:LEU:HG	1:B:52:ILE:HD11	1.97	0.47
1:A:263:VAL:CG1	1:A:267:ILE:H	2.27	0.47
1:A:145:CYS:SG	1:A:148:CYS:N	2.84	0.47
1:A:142:ILE:HD11	1:A:219:VAL:HG11	1.96	0.47
1:A:278:LYS:HB3	1:A:281:THR:OG1	2.15	0.47
1:A:50:THR:HG23	1:A:139:GLN:HB3	1.97	0.47
1:A:23:SER:HB3	1:A:80:THR:HG23	1.96	0.46
1:A:15:ALA:HB1	1:A:31:LEU:HD11	1.96	0.46
1:A:213:ASN:HB3	1:A:237:THR:HG22	1.98	0.46
1:A:70:ASN:HB3	1:B:59:PRO:HD3	1.98	0.46
1:B:263:VAL:O	1:B:266:GLN:HG3	2.16	0.46
1:A:92:LEU:HD22	1:A:267:ILE:HD11	1.97	0.45
1:B:299:ARG:NH1	1:B:300:TYR:OH	2.48	0.45
1:A:33:GLU:OE2	1:A:72:ARG:NH1	2.49	0.45
1:B:17:THR:HG21	1:B:307:ALA:HB1	1.98	0.45
1:A:235:LYS:HB2	1:A:242:GLU:HG2	1.99	0.45
1:B:145:CYS:SG	1:B:148:CYS:N	2.85	0.45
1:A:13:PRO:HA	1:A:71:MET:SD	2.57	0.44
1:A:220:LEU:HB2	1:A:231:HIS:HB2	1.98	0.44
1:B:263:VAL:HG11	1:B:267:ILE:HD12	1.98	0.44
1:A:16:HIS:CE1	1:A:72:ARG:HA	2.52	0.44
1:A:124:ALA:HB3	1:A:294:ASP:HB3	2.00	0.44
1:B:4:THR:HG21	1:B:30:VAL:HG13	2.00	0.44
1:B:50:THR:HG23	1:B:139:GLN:HB3	1.98	0.44
1:A:7:VAL:HG23	1:A:30:VAL:HG23	2.00	0.44
1:A:3:HIS:ND1	1:A:114:THR:HG23	2.33	0.44
1:B:218:GLU:HG2	1:B:233:ARG:HD2	1.99	0.43
1:A:31:LEU:HB3	1:A:82:VAL:HG22	2.00	0.43
1:B:7:VAL:HG23	1:B:30:VAL:HG23	2.01	0.43
1:A:260:THR:HG21	1:A:273:GLY:O	2.18	0.43
1:B:46:LEU:HA	1:B:46:LEU:HD12	1.85	0.42
1:B:152:VAL:HG11	3:B:403:MLI:H11	2.02	0.42
1:B:85:GLU:HB3	1:B:101:THR:HG23	2.02	0.42
1:A:24:ARG:NH1	1:B:148:CYS:O	2.52	0.42
1:B:50:THR:OG1	3:B:403:MLI:O9	2.37	0.42
1:A:56:PRO:HB3	1:B:17:THR:HG22	2.02	0.42
1:A:127:ARG:HB2	1:A:254:VAL:HG23	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:GLU:O	1:B:103:TRP:HB2	2.20	0.41
1:A:118:VAL:HG13	1:A:289:VAL:HG23	2.03	0.40
1:A:183:VAL:O	1:A:208:VAL:HA	2.20	0.40
1:A:11:SER:HB3	1:A:33:GLU:HA	2.03	0.40
1:A:34:GLY:HA2	1:A:85:GLU:C	2.41	0.40
1:A:92:LEU:C	1:A:286:ARG:HH12	2.24	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	320/374 (86%)	305 (95%)	15 (5%)	0	100	100
1	B	315/374 (84%)	300 (95%)	15 (5%)	0	100	100
All	All	635/748 (85%)	605 (95%)	30 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	237/294 (81%)	236 (100%)	1 (0%)	92	96
1	B	236/294 (80%)	236 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	473/588 (80%)	472 (100%)	1 (0%)	94	98

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	93	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

18 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	401	-	51,58,58	3.86	19 (37%)	57,89,89	2.55	12 (21%)
3	MLI	A	402	-	0,6,6	0.00	-	0,7,7	0.00	-
3	MLI	A	403	-	0,6,6	0.00	-	0,7,7	0.00	-
3	MLI	A	404	-	0,6,6	0.00	-	0,7,7	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	FMT	A	405	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	A	406	-	0,2,2	0.00	-	0,1,1	0.00	-
5	NAP	A	407	-	29,33,52	3.58	9 (31%)	34,52,80	2.38	6 (17%)
6	ACT	A	408	-	1,3,3	6.47	1 (100%)	0,3,3	0.00	-
2	FAD	B	401	-	51,58,58	3.85	19 (37%)	57,89,89	2.60	12 (21%)
3	MLI	B	402	-	0,6,6	0.00	-	0,7,7	0.00	-
3	MLI	B	403	-	0,6,6	0.00	-	0,7,7	0.00	-
4	FMT	B	404	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	B	405	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	B	406	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	B	407	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	B	408	-	0,2,2	0.00	-	0,1,1	0.00	-
5	NAP	B	409	-	29,33,52	3.60	9 (31%)	34,52,80	2.41	6 (17%)
7	MLT	B	410	-	2,8,8	0.42	0	4,10,10	1.26	1 (25%)

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	401	-	-	0/28/50/50	0/6/6/6
3	MLI	A	402	-	-	0/0/4/4	0/0/0/0
3	MLI	A	403	-	-	0/0/4/4	0/0/0/0
3	MLI	A	404	-	-	0/0/4/4	0/0/0/0
4	FMT	A	405	-	-	0/0/0/0	0/0/0/0
4	FMT	A	406	-	-	0/0/0/0	0/0/0/0
5	NAP	A	407	-	-	0/17/37/67	0/3/3/5
6	ACT	A	408	-	-	0/0/0/0	0/0/0/0
2	FAD	B	401	-	-	0/28/50/50	0/6/6/6
3	MLI	B	402	-	-	0/0/4/4	0/0/0/0
3	MLI	B	403	-	-	0/0/4/4	0/0/0/0
4	FMT	B	404	-	-	0/0/0/0	0/0/0/0
4	FMT	B	405	-	-	0/0/0/0	0/0/0/0
4	FMT	B	406	-	-	0/0/0/0	0/0/0/0
4	FMT	B	407	-	-	0/0/0/0	0/0/0/0
4	FMT	B	408	-	-	0/0/0/0	0/0/0/0
5	NAP	B	409	-	-	0/17/37/67	0/3/3/5
7	MLT	B	410	-	-	0/2/8/8	0/0/0/0

All (57) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	FAD	O4B-C4B	-6.66	1.30	1.45
2	B	401	FAD	O4B-C4B	-6.64	1.30	1.45
5	B	409	NAP	O4B-C4B	-6.06	1.31	1.45
5	A	407	NAP	O4B-C4B	-5.89	1.31	1.45
2	B	401	FAD	O3B-C3B	-4.11	1.32	1.43
2	A	401	FAD	O3B-C3B	-4.07	1.33	1.43
5	B	409	NAP	C5A-C4A	-2.62	1.34	1.40
5	A	407	NAP	C5A-C4A	-2.61	1.34	1.40
5	A	407	NAP	O3B-C3B	-2.42	1.37	1.43
5	B	409	NAP	O3B-C3B	-2.38	1.37	1.43
2	A	401	FAD	C5A-C4A	-2.35	1.35	1.40
2	B	401	FAD	C5A-C4A	-2.35	1.35	1.40
2	A	401	FAD	O2'-C2'	-2.30	1.38	1.43
2	B	401	FAD	O2'-C2'	-2.25	1.38	1.43
2	A	401	FAD	C2A-N1A	2.14	1.37	1.33
5	A	407	NAP	C2A-N3A	2.14	1.35	1.32
5	B	409	NAP	C2A-N3A	2.16	1.35	1.32
2	B	401	FAD	C2A-N1A	2.19	1.38	1.33
2	A	401	FAD	C4A-N3A	2.19	1.38	1.35
2	B	401	FAD	C4A-N3A	2.23	1.38	1.35
5	A	407	NAP	PN-O5D	2.39	1.64	1.54
5	B	409	NAP	PN-O5D	2.39	1.64	1.54
2	B	401	FAD	C6A-N6A	2.44	1.44	1.34
2	A	401	FAD	C6A-N6A	2.46	1.44	1.34
5	B	409	NAP	PA-O5B	2.74	1.70	1.59
5	A	407	NAP	PA-O5B	2.78	1.70	1.59
2	A	401	FAD	C4X-C10	2.91	1.45	1.41
2	B	401	FAD	C4X-C10	3.00	1.46	1.41
2	A	401	FAD	O2B-C2B	3.14	1.50	1.43
2	B	401	FAD	O2B-C2B	3.16	1.50	1.43
2	B	401	FAD	C2A-N3A	3.21	1.37	1.32
2	A	401	FAD	C2A-N3A	3.25	1.37	1.32
5	A	407	NAP	C6A-N6A	3.49	1.48	1.34
5	B	409	NAP	C6A-N6A	3.51	1.48	1.34
2	A	401	FAD	C2-N1	4.68	1.47	1.38
2	B	401	FAD	C2-N1	4.71	1.47	1.38
2	A	401	FAD	C4-C4X	4.88	1.50	1.41
2	B	401	FAD	C4-C4X	4.93	1.50	1.41
2	A	401	FAD	C2-N3	5.44	1.48	1.38
2	B	401	FAD	C2-N3	5.46	1.49	1.38
2	A	401	FAD	C4-N3	6.22	1.44	1.33
2	B	401	FAD	C4-N3	6.28	1.44	1.33
5	A	407	NAP	PN-O3	6.45	1.69	1.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	409	NAP	PN-O3	6.46	1.69	1.60
6	A	408	ACT	CH3-C	6.47	1.57	1.48
2	A	401	FAD	C4X-N5	7.16	1.43	1.33
2	B	401	FAD	C4X-N5	7.33	1.44	1.33
2	A	401	FAD	C5X-N5	7.53	1.46	1.35
2	B	401	FAD	C5X-N5	7.66	1.47	1.35
2	A	401	FAD	C10-N1	8.93	1.45	1.33
2	B	401	FAD	C10-N1	8.93	1.45	1.33
2	B	401	FAD	C9A-N10	9.15	1.49	1.38
2	A	401	FAD	C9A-N10	9.32	1.50	1.38
2	B	401	FAD	O4B-C1B	15.09	1.62	1.41
2	A	401	FAD	O4B-C1B	15.32	1.62	1.41
5	A	407	NAP	O4B-C1B	15.40	1.62	1.41
5	B	409	NAP	O4B-C1B	15.44	1.62	1.41

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	FAD	N3A-C2A-N1A	-10.41	119.95	128.86
5	B	409	NAP	N3A-C2A-N1A	-10.37	119.98	128.86
5	A	407	NAP	N3A-C2A-N1A	-10.29	120.05	128.86
2	A	401	FAD	N3A-C2A-N1A	-10.07	120.25	128.86
2	A	401	FAD	N6A-C6A-N1A	-7.23	103.57	118.57
2	B	401	FAD	N6A-C6A-N1A	-7.16	103.70	118.57
2	B	401	FAD	C7M-C7-C6	-5.14	107.71	120.36
2	B	401	FAD	C4B-O4B-C1B	-4.91	104.71	109.83
2	A	401	FAD	C7M-C7-C6	-4.91	108.28	120.36
5	A	407	NAP	N6A-C6A-N1A	-4.90	108.39	118.57
5	B	409	NAP	N6A-C6A-N1A	-4.89	108.41	118.57
5	A	407	NAP	C1B-N9A-C4A	-3.66	120.31	126.64
5	B	409	NAP	C1B-N9A-C4A	-3.61	120.40	126.64
2	A	401	FAD	C4B-O4B-C1B	-3.35	106.34	109.83
5	B	409	NAP	PA-O3-PN	-2.87	122.98	132.63
2	A	401	FAD	C4X-C4-N3	-2.74	119.57	123.47
2	A	401	FAD	P-O3P-PA	-2.67	123.64	132.63
2	B	401	FAD	C4X-C4-N3	-2.64	119.72	123.47
5	A	407	NAP	PA-O3-PN	-2.48	124.30	132.63
7	B	410	MLT	C4-C3-C2	-2.40	110.89	113.96
5	B	409	NAP	C4B-O4B-C1B	-2.33	107.40	109.83
2	B	401	FAD	P-O3P-PA	-2.23	125.12	132.63
5	A	407	NAP	C4B-O4B-C1B	-2.05	107.69	109.83
2	B	401	FAD	C1'-N10-C9A	2.55	120.58	118.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	FAD	C5X-C9A-N10	2.70	119.77	117.71
2	A	401	FAD	C5X-C9A-N10	2.86	119.90	117.71
2	A	401	FAD	C4X-N5-C5X	3.09	120.00	116.76
2	A	401	FAD	C1'-N10-C9A	3.12	121.09	118.31
2	B	401	FAD	C4X-N5-C5X	3.15	120.06	116.76
2	B	401	FAD	C4-N3-C2	5.33	119.68	115.14
2	A	401	FAD	C7M-C7-C8	5.33	131.85	120.72
2	A	401	FAD	C4-N3-C2	5.37	119.72	115.14
2	B	401	FAD	C7M-C7-C8	5.40	132.01	120.72
5	A	407	NAP	C5A-C6A-N6A	5.41	131.49	120.47
5	B	409	NAP	C5A-C6A-N6A	5.42	131.52	120.47
2	B	401	FAD	C5A-C6A-N6A	7.72	136.21	120.47
2	A	401	FAD	C5A-C6A-N6A	7.82	136.41	120.47

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	402	MLI	1	0
2	B	401	FAD	3	0
3	B	403	MLI	2	0
5	B	409	NAP	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	324/374 (86%)	0.06	5 (1%) 73 61	3, 18, 33, 54	0
1	B	319/374 (85%)	0.11	10 (3%) 49 33	4, 19, 42, 73	0
All	All	643/748 (85%)	0.09	15 (2%) 60 47	3, 19, 37, 73	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	326	SER	5.0
1	B	194	ALA	3.8
1	B	326	SER	3.5
1	B	48	THR	2.9
1	B	192	LEU	2.7
1	B	44	GLY	2.6
1	B	165	GLY	2.6
1	A	108	ASP	2.4
1	A	107	PRO	2.3
1	A	325	GLU	2.3
1	B	47	THR	2.2
1	A	94	SER	2.2
1	B	271	GLU	2.1
1	B	272	ASP	2.1
1	B	237	THR	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
4	FMT	B	405	3/3	0.33	0.49	121,121,123,124	0
3	MLI	A	403	7/7	0.51	0.74	149,155,159,161	0
3	MLI	B	402	7/7	0.62	0.53	146,154,170,172	0
4	FMT	B	404	3/3	0.65	0.27	125,125,127,129	0
5	NAP	B	409	31/48	0.77	0.27	131,164,201,207	31
6	ACT	A	408	4/4	0.78	0.34	107,118,119,124	0
4	FMT	B	407	3/3	0.78	0.42	101,101,101,102	0
3	MLI	A	404	7/7	0.79	0.54	152,152,154,155	0
2	FAD	B	401	53/53	0.82	0.39	72,154,191,197	53
4	FMT	B	408	3/3	0.82	0.44	89,89,91,95	0
5	NAP	A	407	31/48	0.85	0.20	104,148,181,184	0
4	FMT	B	406	3/3	0.86	0.61	117,117,118,118	0
2	FAD	A	401	53/53	0.86	0.36	66,149,169,173	0
4	FMT	A	405	3/3	0.87	0.24	115,115,117,118	0
3	MLI	A	402	7/7	0.89	0.23	107,112,133,134	0
3	MLI	B	403	7/7	0.90	0.28	121,126,136,139	0
4	FMT	A	406	3/3	0.91	0.28	97,97,98,99	0
7	MLT	B	410	9/9	0.91	0.28	78,96,117,121	0

6.5 Other polymers [i](#)

There are no such residues in this entry.