



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 16, 2017 – 02:43 PM EST

PDB ID : 5TTK  
Title : Crystal Structure of Selenomethionine-incorporated Nicotine Oxidoreductase from *Pseudomonas putida*  
Authors : Tararina, M.A.; Janda, K.D.; Allen, K.N.  
Deposited on : 2016-11-03  
Resolution : 2.51 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7.1 (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

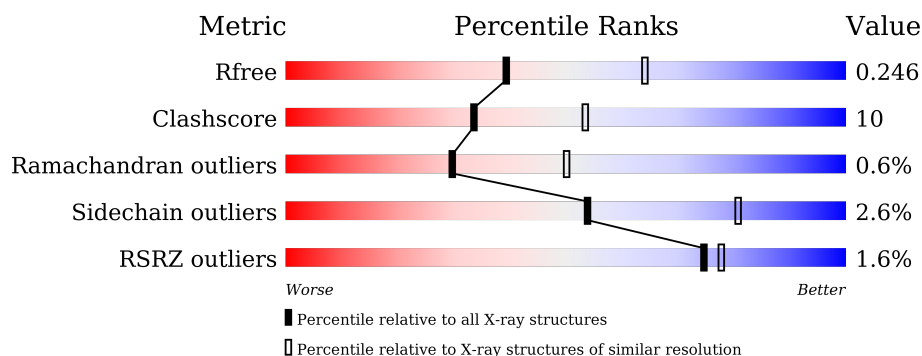
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.51 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	490	<div> <div>%</div> <div> <div></div> <div>70%</div> <div>17%</div> <div>••</div> <div>12%</div> </div> </div>
1	B	490	<div> <div>2%</div> <div> <div></div> <div>71%</div> <div>16%</div> <div>•</div> <div>12%</div> </div> </div>
1	C	490	<div> <div>%</div> <div> <div></div> <div>71%</div> <div>16%</div> <div>••</div> <div>11%</div> </div> </div>
1	D	490	<div> <div>%</div> <div> <div></div> <div>64%</div> <div>23%</div> <div>•</div> <div>11%</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13917 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 Amine oxidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	432	Total	C	N	O	S	Se	0	0	0
			3354	2126	583	633	3	9			
1	B	432	Total	C	N	O	S	Se	0	0	0
			3354	2126	583	633	3	9			
1	C	435	Total	C	N	O	S	Se	0	0	0
			3381	2143	588	638	3	9			
1	D	435	Total	C	N	O	S	Se	0	0	0
			3381	2143	588	638	3	9			

There are 32 discrepancies between the modelled and reference sequences:

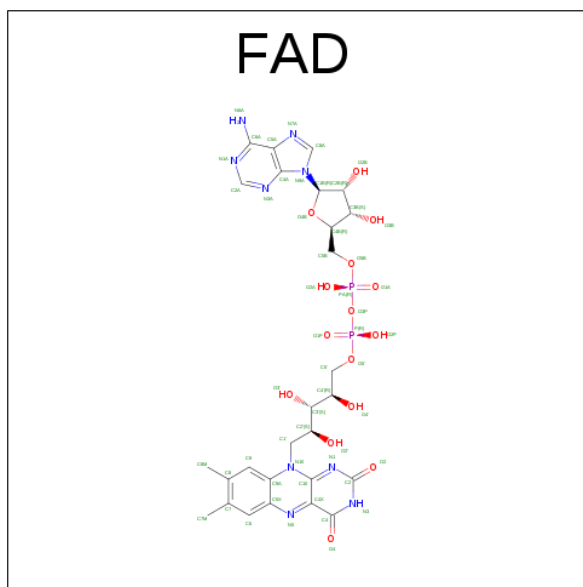
Chain	Residue	Modelled	Actual	Comment	Reference
A	483	LEU	-	expression tag	UNP F8G0P2
A	484	GLU	-	expression tag	UNP F8G0P2
A	485	HIS	-	expression tag	UNP F8G0P2
A	486	HIS	-	expression tag	UNP F8G0P2
A	487	HIS	-	expression tag	UNP F8G0P2
A	488	HIS	-	expression tag	UNP F8G0P2
A	489	HIS	-	expression tag	UNP F8G0P2
A	490	HIS	-	expression tag	UNP F8G0P2
B	483	LEU	-	expression tag	UNP F8G0P2
B	484	GLU	-	expression tag	UNP F8G0P2
B	485	HIS	-	expression tag	UNP F8G0P2
B	486	HIS	-	expression tag	UNP F8G0P2
B	487	HIS	-	expression tag	UNP F8G0P2
B	488	HIS	-	expression tag	UNP F8G0P2
B	489	HIS	-	expression tag	UNP F8G0P2
B	490	HIS	-	expression tag	UNP F8G0P2
C	483	LEU	-	expression tag	UNP F8G0P2
C	484	GLU	-	expression tag	UNP F8G0P2
C	485	HIS	-	expression tag	UNP F8G0P2
C	486	HIS	-	expression tag	UNP F8G0P2
C	487	HIS	-	expression tag	UNP F8G0P2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	488	HIS	-	expression tag	UNP F8G0P2
C	489	HIS	-	expression tag	UNP F8G0P2
C	490	HIS	-	expression tag	UNP F8G0P2
D	483	LEU	-	expression tag	UNP F8G0P2
D	484	GLU	-	expression tag	UNP F8G0P2
D	485	HIS	-	expression tag	UNP F8G0P2
D	486	HIS	-	expression tag	UNP F8G0P2
D	487	HIS	-	expression tag	UNP F8G0P2
D	488	HIS	-	expression tag	UNP F8G0P2
D	489	HIS	-	expression tag	UNP F8G0P2
D	490	HIS	-	expression tag	UNP F8G0P2

- 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
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		

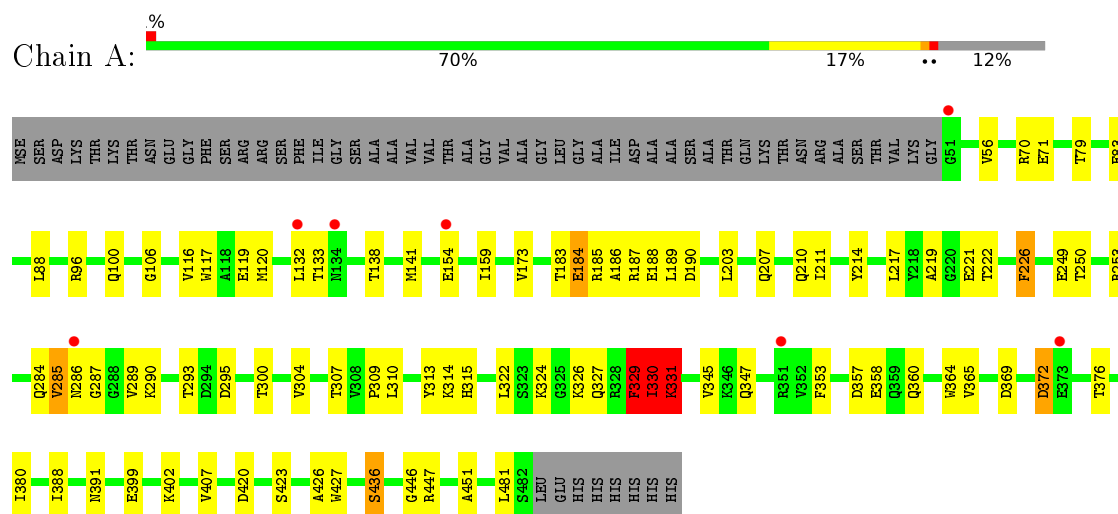
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	65	Total 65	O 65	0	0
3	B	64	Total 64	O 64	0	0
3	C	58	Total 58	O 58	0	0
3	D	48	Total 48	O 48	0	0

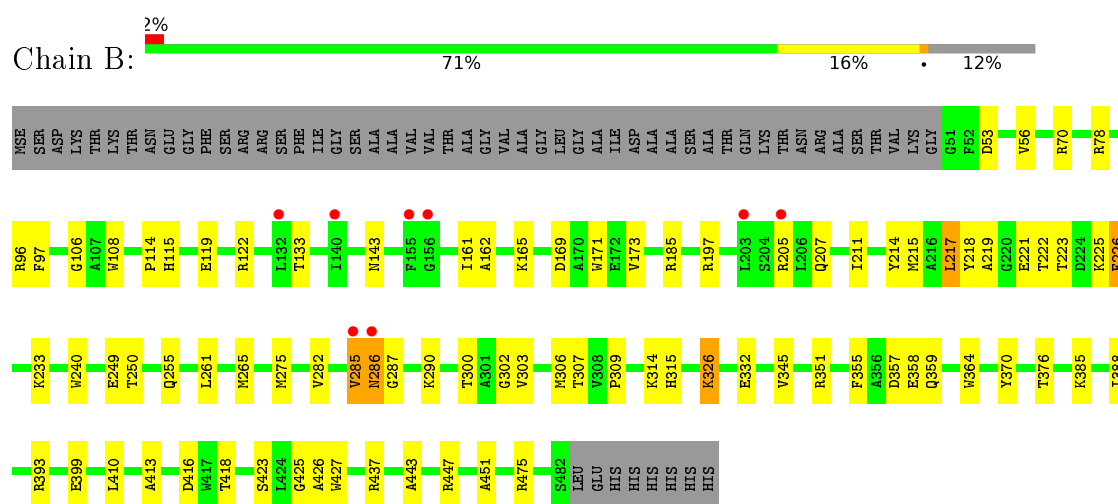
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

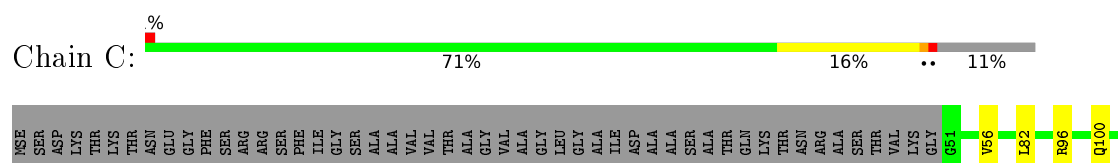
#### • Molecule 1: Amine oxidase

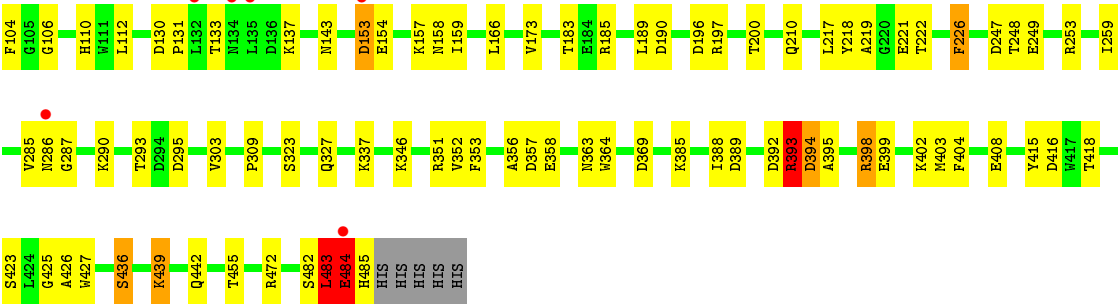


#### • Molecule 1: Amine oxidase

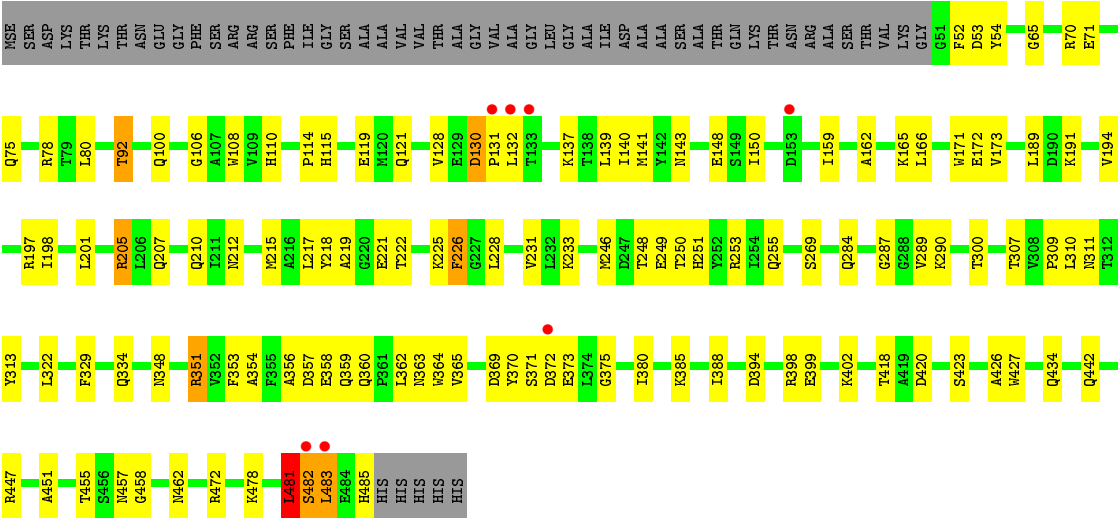


#### • Molecule 1: Amine oxidase





• Molecule 1: Amine oxidase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	97.10Å 126.13Å 168.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.55 – 2.51 48.55 – 2.51	Depositor EDS
% Data completeness (in resolution range)	95.0 (48.55-2.51) 95.0 (48.55-2.51)	Depositor EDS
$R_{merge}$	0.19	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.55 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.10.1 _2155: ???)	Depositor
R, $R_{free}$	0.203 , 0.246 0.202 , 0.246	Depositor DCC
$R_{free}$ test set	1997 reflections (2.94%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.7	Xtriage
Anisotropy	0.118	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 41.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	13917	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.29% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.47	1/3416 (0.0%)	0.68	4/4612 (0.1%)
1	B	0.30	0/3416	0.55	4/4612 (0.1%)
1	C	0.43	3/3444 (0.1%)	0.71	9/4650 (0.2%)
1	D	0.35	0/3444	0.57	3/4650 (0.1%)
All	All	0.39	4/13720 (0.0%)	0.63	20/18524 (0.1%)

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
1	C	0	1
All	All	0	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	393	ARG	CZ-NH1	-7.40	1.23	1.33
1	C	393	ARG	NE-CZ	-7.28	1.23	1.33
1	C	484	GLU	CA-C	-6.59	1.35	1.52
1	A	330	ILE	CA-C	5.16	1.66	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	331	LYS	N-CA-CB	-18.37	77.53	110.60
1	C	484	GLU	N-CA-C	-13.23	75.28	111.00
1	C	484	GLU	C-N-CA	-11.39	93.23	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	329	PHE	O-C-N	-10.43	106.01	122.70
1	B	217	LEU	CA-CB-CG	8.59	135.06	115.30
1	C	398	ARG	NE-CZ-NH1	8.02	124.31	120.30
1	C	398	ARG	CB-CG-CD	7.37	130.75	111.60
1	C	393	ARG	NE-CZ-NH1	-6.79	116.91	120.30
1	A	329	PHE	C-N-CA	-6.38	105.76	121.70
1	C	439	LYS	CA-CB-CG	6.21	127.06	113.40
1	C	483	LEU	C-N-CA	6.18	137.16	121.70
1	D	130	ASP	CB-CG-OD1	6.06	123.76	118.30
1	C	439	LYS	CD-CE-NZ	5.91	125.29	111.70
1	C	483	LEU	CA-C-N	-5.67	104.74	117.20
1	D	191	LYS	CB-CG-CD	-5.51	97.26	111.60
1	A	154	GLU	CB-CA-C	-5.49	99.42	110.40
1	B	217	LEU	CB-CG-CD2	-5.33	101.94	111.00
1	B	217	LEU	CB-CG-CD1	5.24	119.91	111.00
1	B	359	GLN	CA-CB-CG	-5.20	101.96	113.40
1	D	481	LEU	N-CA-C	-5.10	97.22	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	329	PHE	Mainchain
1	C	483	LEU	Mainchain

## 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	3354	0	3294	70	0
1	B	3354	0	3296	60	0
1	C	3381	0	3320	66	0
1	D	3381	0	3320	90	0
2	A	53	0	30	4	0
2	B	53	0	30	4	0
2	C	53	0	30	4	0
2	D	53	0	30	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	65	0	0	2	0
3	B	64	0	0	2	0
3	C	58	0	0	1	0
3	D	48	0	0	0	0
All	All	13917	0	13350	276	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 10.

All (276) 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:B:217:LEU:HD21	1:B:364:TRP:HB2	1.29	1.14
1:A:329:PHE:O	1:A:330:ILE:C	1.94	0.99
1:A:286:ASN:OD1	1:A:287:GLY:N	1.97	0.97
1:A:185:ARG:NH1	1:A:188:GLU:OE1	2.01	0.93
1:A:96:ARG:NH2	1:B:255:GLN:NE2	2.18	0.92
1:D:478:LYS:O	1:D:482:SER:HB3	1.71	0.90
1:A:284:GLN:NE2	1:A:447:ARG:H	1.75	0.83
1:A:96:ARG:NH2	1:B:255:GLN:HE21	1.77	0.83
1:C:393:ARG:O	1:C:395:ALA:N	2.12	0.81
1:C:483:LEU:C	1:C:485:HIS:N	2.27	0.80
1:D:70:ARG:NH1	1:D:71:GLU:OE1	2.16	0.79
1:D:137:LYS:HE3	1:D:139:LEU:HD21	1.66	0.78
1:D:70:ARG:NH2	1:D:119:GLU:OE2	2.17	0.77
1:C:484:GLU:O	1:C:485:HIS:O	2.02	0.76
1:D:482:SER:OG	1:D:483:LEU:N	2.12	0.76
1:C:110:HIS:HD2	1:C:112:LEU:H	1.33	0.75
1:D:482:SER:O	1:D:483:LEU:HG	1.86	0.75
1:D:222:THR:OG1	1:D:225:LYS:HG3	1.86	0.75
1:D:215:MSE:HE1	1:D:231:VAL:HG21	1.70	0.74
1:B:217:LEU:CD2	1:B:364:TRP:HB2	2.13	0.73
1:D:482:SER:C	1:D:483:LEU:HG	2.11	0.72
1:A:70:ARG:NH1	1:A:71:GLU:OE1	2.23	0.71
1:A:117:TRP:HA	1:A:120:MSE:HE3	1.71	0.71
1:B:357:ASP:OD2	1:B:358:GLU:N	2.22	0.71
1:B:332:GLU:OE1	1:B:437:ARG:NH2	2.24	0.70
1:D:394:ASP:OD2	1:D:398:ARG:NH2	2.25	0.70
1:B:169:ASP:OD2	1:B:197:ARG:NH2	2.26	0.69
1:D:485:HIS:ND1	1:D:485:HIS:O	2.24	0.69
1:D:233:LYS:NZ	1:D:458:GLY:O	2.24	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:481:LEU:O	1:D:482:SER:HB3	1.93	0.69
1:A:284:GLN:HE22	1:A:447:ARG:H	1.41	0.68
1:B:169:ASP:CG	1:B:197:ARG:HH22	1.97	0.68
1:C:439:LYS:HE2	1:C:442:GLN:OE1	1.94	0.68
1:A:190:ASP:OD2	1:A:436:SER:OG	2.11	0.67
1:B:217:LEU:HD21	1:B:364:TRP:CB	2.15	0.67
1:C:100:GLN:NE2	1:C:369:ASP:OD2	2.27	0.67
1:D:75:GLN:HG3	1:D:478:LYS:HD2	1.76	0.67
1:C:484:GLU:O	1:C:485:HIS:C	2.18	0.67
1:C:398:ARG:O	1:C:402:LYS:HG3	1.95	0.66
1:C:285:VAL:HG11	1:C:290:LYS:HE3	1.77	0.65
1:A:293:THR:OG1	1:A:295:ASP:OD1	2.15	0.64
1:D:354:ALA:HB3	1:D:365:VAL:HG22	1.80	0.64
1:A:285:VAL:HG11	1:A:290:LYS:HG3	1.78	0.64
1:C:159:ILE:HD13	1:C:210:GLN:HB3	1.79	0.64
1:D:309:PRO:HD3	2:D:501:FAD:H51A	1.79	0.64
1:C:385:LYS:HD3	1:C:418:THR:OG1	1.99	0.63
1:C:218:TYR:OH	1:C:249:GLU:OE2	2.13	0.62
1:A:100:GLN:NE2	1:A:369:ASP:OD2	2.33	0.61
1:A:96:ARG:HH22	1:B:255:GLN:HE21	1.48	0.61
1:A:183:THR:HG22	1:A:186:ALA:H	1.66	0.60
1:C:130:ASP:OD1	1:C:253:ARG:NH1	2.34	0.60
1:A:327:GLN:O	1:A:331:LYS:HB2	2.01	0.60
1:B:388:ILE:HD12	1:B:399:GLU:HG3	1.84	0.60
1:B:309:PRO:HD3	2:B:501:FAD:H51A	1.83	0.60
1:A:253:ARG:NH1	3:A:602:HOH:O	2.35	0.59
1:C:309:PRO:HD3	2:C:501:FAD:H51A	1.83	0.59
1:C:484:GLU:OE2	1:C:484:GLU:HA	2.01	0.59
1:D:307:THR:HG22	1:D:451:ALA:HB3	1.85	0.59
1:D:217:LEU:HD11	1:D:364:TRP:HB2	1.83	0.58
1:C:143:ASN:ND2	1:C:357:ASP:HB3	2.18	0.58
1:A:284:GLN:HE22	1:A:446:GLY:HA3	1.68	0.58
1:A:329:PHE:O	1:A:331:LYS:N	2.37	0.58
1:D:53:ASP:O	1:D:78:ARG:NH2	2.37	0.57
1:C:106:GLY:HA2	2:C:501:FAD:C4X	2.34	0.57
1:D:357:ASP:OD1	1:D:358:GLU:N	2.30	0.57
1:B:282:VAL:HG21	1:B:306:MSE:HE1	1.87	0.57
1:B:357:ASP:OD2	1:B:358:GLU:HG3	2.05	0.57
1:A:217:LEU:HD11	1:A:364:TRP:HB2	1.86	0.56
1:D:100:GLN:NE2	1:D:369:ASP:OD2	2.36	0.56
1:D:108:TRP:CG	1:D:250:THR:HA	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:309:PRO:HD3	2:A:501:FAD:H51A	1.88	0.56
1:B:223:THR:N	1:B:358:GLU:OE2	2.29	0.55
1:C:133:THR:HB	1:C:351:ARG:HD2	1.88	0.55
1:A:141:MSE:HE2	1:A:360:GLN:HB3	1.89	0.55
1:B:106:GLY:HA2	2:B:501:FAD:C4X	2.36	0.55
1:B:309:PRO:HG2	1:B:423:SER:HA	1.89	0.55
1:C:110:HIS:CD2	1:C:112:LEU:H	2.20	0.55
1:A:391:ASN:HB3	1:D:483:LEU:HD22	1.89	0.55
1:B:285:VAL:HG11	1:B:290:LYS:HE2	1.88	0.54
1:C:154:GLU:HG2	1:C:158:ASN:ND2	2.22	0.54
1:D:194:VAL:HG11	1:D:215:MSE:HE2	1.89	0.54
1:A:116:VAL:HG22	1:A:120:MSE:HE2	1.91	0.53
1:D:198:ILE:HA	1:D:201:LEU:HD23	1.91	0.53
1:B:108:TRP:CG	1:B:250:THR:HA	2.43	0.53
1:A:106:GLY:HA2	2:A:501:FAD:C4X	2.39	0.53
1:A:159:ILE:HD13	1:A:210:GLN:HB3	1.90	0.53
1:C:389:ASP:OD2	1:C:392:ASP:HB2	2.08	0.53
1:A:420:ASP:HB2	1:D:485:HIS:HB2	1.91	0.52
1:D:218:TYR:OH	1:D:249:GLU:OE2	2.17	0.52
1:A:219:ALA:HB2	1:A:226:PHE:CD2	2.44	0.52
1:C:346:LYS:HD2	1:C:408:GLU:OE2	2.10	0.52
1:B:173:VAL:HG22	1:B:185:ARG:HG2	1.92	0.52
1:B:233:LYS:NZ	3:B:603:HOH:O	2.40	0.52
1:D:159:ILE:HD13	1:D:210:GLN:HB3	1.91	0.52
1:D:478:LYS:O	1:D:481:LEU:O	2.28	0.52
1:B:314:LYS:HG3	1:B:315:HIS:ND1	2.25	0.52
1:C:484:GLU:C	1:C:485:HIS:O	2.48	0.52
1:D:140:ILE:HG22	1:D:150:ILE:HD11	1.92	0.51
1:B:169:ASP:HB2	1:B:173:VAL:HG23	1.93	0.51
1:B:70:ARG:NH2	1:B:119:GLU:OE2	2.43	0.51
1:D:171:TRP:HD1	1:D:172:GLU:OE1	1.94	0.51
1:C:153:ASP:O	1:C:157:LYS:HG3	2.10	0.51
1:D:115:HIS:O	1:D:119:GLU:HG2	2.11	0.51
1:C:418:THR:HA	1:C:425:GLY:N	2.25	0.51
1:D:371:SER:OG	1:D:373:GLU:O	2.25	0.50
1:A:173:VAL:HG11	1:A:189:LEU:HD12	1.93	0.50
1:C:219:ALA:HB2	1:C:226:PHE:CD1	2.46	0.50
1:B:302:GLY:O	1:B:447:ARG:HD3	2.10	0.50
1:D:373:GLU:O	1:D:375:GLY:N	2.41	0.50
1:B:217:LEU:CD2	1:B:364:TRP:CB	2.84	0.50
1:D:219:ALA:HB2	1:D:226:PHE:CD2	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:357:ASP:OD1	1:C:358:GLU:N	2.40	0.50
1:C:110:HIS:HE1	1:C:247:ASP:OD2	1.95	0.50
1:D:356:ALA:HB3	1:D:363:ASN:HA	1.93	0.50
1:D:442:GLN:NE2	1:D:457:ASN:OD1	2.37	0.50
1:B:307:THR:HG22	1:B:451:ALA:HB3	1.93	0.49
1:A:345:VAL:HG12	1:A:376:THR:HB	1.93	0.49
1:A:345:VAL:HG21	1:A:407:VAL:CG2	2.42	0.49
1:B:115:HIS:O	1:B:119:GLU:HG2	2.13	0.49
1:B:133:THR:HG21	1:B:351:ARG:HD2	1.95	0.49
1:A:426:ALA:HB1	1:A:427:TRP:CE3	2.47	0.49
1:D:351:ARG:HD2	1:D:370:TYR:OH	2.13	0.49
1:A:70:ARG:NH2	1:A:119:GLU:OE2	2.46	0.49
1:A:357:ASP:OD2	1:A:358:GLU:N	2.37	0.49
1:C:104:PHE:CG	1:C:415:TYR:HB2	2.47	0.49
1:C:455:THR:O	1:C:472:ARG:NH2	2.46	0.49
1:A:96:ARG:HH21	1:B:255:GLN:NE2	2.08	0.48
1:D:162:ALA:HB2	1:D:207:GLN:HB3	1.93	0.48
1:A:307:THR:HG22	1:A:451:ALA:HB3	1.95	0.48
1:C:110:HIS:CG	1:C:248:THR:HG22	2.49	0.48
1:C:183:THR:HG22	1:C:185:ARG:N	2.28	0.48
1:D:329:PHE:CE1	1:D:434:GLN:HG2	2.48	0.48
1:D:455:THR:O	1:D:472:ARG:NH2	2.46	0.48
1:B:393:ARG:NH2	1:B:410:LEU:O	2.47	0.48
1:C:293:THR:OG1	1:C:295:ASP:OD2	2.32	0.48
1:B:282:VAL:HG21	1:B:306:MSE:CE	2.44	0.47
1:C:483:LEU:HA	1:C:485:HIS:O	2.14	0.47
1:B:162:ALA:HB2	1:B:207:GLN:HB3	1.96	0.47
1:D:165:LYS:HB3	1:D:201:LEU:HD12	1.96	0.47
1:B:326:LYS:NZ	1:B:443:ALA:O	2.48	0.47
1:C:190:ASP:OD2	1:C:436:SER:OG	2.31	0.47
1:D:348:ASN:ND2	1:D:372:ASP:OD1	2.46	0.47
1:C:485:HIS:C	1:C:485:HIS:CD2	2.88	0.47
1:D:108:TRP:O	2:D:501:FAD:N3	2.46	0.47
1:A:214:TYR:OH	1:A:249:GLU:OE2	2.26	0.47
1:A:217:LEU:CD1	1:A:364:TRP:HB2	2.45	0.47
1:D:148:GLU:OE2	1:D:205:ARG:NH2	2.47	0.47
1:A:106:GLY:HA2	2:A:501:FAD:N5	2.30	0.47
1:C:137:LYS:HB2	1:C:352:VAL:HG12	1.96	0.47
1:A:284:GLN:HE22	1:A:447:ARG:N	2.11	0.47
1:A:184:GLU:HA	1:A:187:ARG:NH1	2.30	0.47
1:A:290:LYS:HG2	1:A:300:THR:HG22	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:166:LEU:HD12	1:D:197:ARG:HG2	1.96	0.46
1:D:197:ARG:O	1:D:201:LEU:HD22	2.14	0.46
1:A:56:VAL:HG13	1:A:79:THR:HG23	1.96	0.46
1:A:285:VAL:HG11	1:A:290:LYS:HE3	1.98	0.46
1:B:221:GLU:HG3	1:B:222:THR:H	1.81	0.46
1:B:218:TYR:OH	1:B:249:GLU:OE2	2.25	0.46
1:D:140:ILE:HG22	1:D:150:ILE:CD1	2.46	0.46
1:D:351:ARG:HD2	1:D:370:TYR:CZ	2.51	0.46
1:C:388:ILE:HD12	1:C:399:GLU:HG3	1.98	0.46
1:B:221:GLU:HG3	1:B:222:THR:N	2.31	0.46
1:C:221:GLU:HG3	1:C:222:THR:N	2.31	0.46
1:C:394:ASP:O	1:C:398:ARG:HG2	2.15	0.46
1:D:130:ASP:O	1:D:132:LEU:N	2.48	0.46
1:C:106:GLY:HA2	2:C:501:FAD:N5	2.31	0.46
1:A:203:LEU:HB3	1:A:207:GLN:HB2	1.98	0.45
1:C:286:ASN:OD1	1:C:287:GLY:N	2.49	0.45
1:C:403:MSE:HG2	1:C:404:PHE:CE2	2.51	0.45
1:B:143:ASN:ND2	3:B:607:HOH:O	2.48	0.45
1:A:185:ARG:HH11	1:A:188:GLU:CD	2.12	0.45
1:B:219:ALA:HB2	1:B:226:PHE:CD2	2.52	0.45
1:D:197:ARG:HH21	1:D:228:LEU:HD21	1.80	0.45
1:A:365:VAL:HG22	1:A:380:ILE:HG12	1.98	0.45
1:B:214:TYR:CD1	1:B:215:MSE:HE2	2.51	0.45
1:C:133:THR:HB	1:C:351:ARG:HH11	1.82	0.45
1:D:426:ALA:HB1	1:D:427:TRP:CE3	2.52	0.45
1:A:289:VAL:HG21	1:A:304:VAL:HB	1.98	0.45
1:B:418:THR:HA	1:B:425:GLY:N	2.31	0.45
1:C:196:ASP:O	1:C:200:THR:HG23	2.17	0.44
1:B:426:ALA:HB1	1:B:427:TRP:CE3	2.52	0.44
1:B:214:TYR:HD1	1:B:215:MSE:HE2	1.82	0.44
1:C:131:PRO:O	1:C:133:THR:HG23	2.17	0.44
1:D:132:LEU:HD12	1:D:251:HIS:CE1	2.52	0.44
1:D:284:GLN:HG2	1:D:289:VAL:HG12	1.98	0.44
1:A:221:GLU:HG3	1:A:222:THR:H	1.81	0.44
1:B:70:ARG:NH2	1:B:265:MSE:HE3	2.32	0.44
1:C:323:SER:O	1:C:327:GLN:HG3	2.16	0.44
1:C:484:GLU:N	1:C:485:HIS:O	2.50	0.44
1:A:345:VAL:CG2	1:A:407:VAL:HB	2.47	0.44
1:B:261:LEU:O	1:B:265:MSE:HG3	2.17	0.44
1:C:154:GLU:HG2	1:C:158:ASN:HD21	1.82	0.44
1:D:143:ASN:HD22	1:D:359:GLN:H	1.66	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:290:LYS:HG2	1:D:300:THR:HG22	2.00	0.44
1:A:221:GLU:HG3	1:A:222:THR:N	2.32	0.44
1:B:475:ARG:NH1	1:D:121:GLN:O	2.51	0.44
1:C:166:LEU:HD12	1:C:197:ARG:HG2	2.00	0.44
1:D:106:GLY:HA2	2:D:501:FAD:N5	2.33	0.44
1:A:250:THR:N	3:A:601:HOH:O	2.34	0.44
1:D:388:ILE:HD12	1:D:399:GLU:HG3	2.00	0.44
1:D:54:TYR:CE2	1:D:80:LEU:HD22	2.53	0.44
1:A:138:THR:HA	1:A:353:PHE:O	2.18	0.44
1:A:309:PRO:HG2	1:A:423:SER:HA	2.00	0.44
1:B:287:GLY:O	1:B:447:ARG:NH1	2.51	0.44
1:C:337:LYS:HA	1:C:418:THR:HG21	2.00	0.44
1:A:116:VAL:O	1:A:120:MSE:HG3	2.19	0.43
1:A:310:LEU:HA	1:A:313:TYR:CD2	2.53	0.43
1:B:122:ARG:HD3	1:D:71:GLU:OE2	2.18	0.43
1:B:53:ASP:HB2	1:B:300:THR:O	2.18	0.43
1:D:70:ARG:HA	1:D:269:SER:CB	2.48	0.43
1:B:345:VAL:HG22	1:B:376:THR:HB	2.01	0.43
1:C:356:ALA:HB3	1:C:363:ASN:HA	2.00	0.43
1:A:372:ASP:OD1	1:A:372:ASP:N	2.52	0.43
1:C:221:GLU:HG3	1:C:222:THR:H	1.84	0.43
1:C:56:VAL:HG13	1:C:303:VAL:HG13	1.99	0.43
1:D:106:GLY:HA2	2:D:501:FAD:C4X	2.49	0.43
1:D:70:ARG:HH22	1:D:119:GLU:CD	2.21	0.43
1:D:143:ASN:ND2	1:D:359:GLN:H	2.17	0.43
1:D:385:LYS:HB2	1:D:385:LYS:HE3	1.64	0.43
1:D:420:ASP:HB3	1:D:423:SER:OG	2.18	0.43
1:D:108:TRP:CD1	1:D:250:THR:HA	2.53	0.43
1:D:128:VAL:N	1:D:253:ARG:O	2.47	0.43
1:D:362:LEU:HD22	1:D:380:ILE:HG21	2.01	0.43
1:A:314:LYS:HG3	1:A:315:HIS:ND1	2.34	0.42
1:A:322:LEU:HD12	1:A:330:ILE:HD11	2.01	0.42
1:B:385:LYS:NZ	1:B:416:ASP:OD1	2.46	0.42
1:C:472:ARG:HD2	3:C:631:HOH:O	2.19	0.42
1:D:250:THR:HG21	1:D:364:TRP:HH2	1.84	0.42
1:D:217:LEU:CD1	1:D:364:TRP:HB2	2.48	0.42
1:D:194:VAL:O	1:D:198:ILE:HG13	2.18	0.42
1:A:185:ARG:HD2	1:A:188:GLU:OE1	2.19	0.42
1:B:161:ILE:O	1:B:165:LYS:HG3	2.20	0.42
1:C:385:LYS:HE2	1:C:416:ASP:OD2	2.19	0.42
1:A:295:ASP:OD2	1:B:78:ARG:NH2	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:393:ARG:HG2	1:C:394:ASP:H	1.84	0.42
1:D:110:HIS:CG	1:D:248:THR:HG22	2.55	0.42
1:D:219:ALA:HB2	1:D:226:PHE:CG	2.53	0.42
1:A:211:ILE:HA	1:A:211:ILE:HD12	1.92	0.42
1:B:106:GLY:HA2	2:B:501:FAD:N5	2.34	0.42
2:C:501:FAD:H9	2:C:501:FAD:H1'1	1.77	0.42
1:A:347:GLN:OE1	1:A:407:VAL:HB	2.20	0.42
1:D:65:GLY:HA2	1:D:307:THR:HG21	2.02	0.42
1:A:426:ALA:HB1	2:A:501:FAD:HM83	2.02	0.42
1:C:219:ALA:HB2	1:C:226:PHE:CG	2.55	0.42
1:D:365:VAL:HG12	1:D:380:ILE:HG23	2.02	0.42
1:C:96:ARG:NH2	1:D:255:GLN:HG2	2.35	0.41
1:D:310:LEU:HA	1:D:313:TYR:CD2	2.55	0.41
1:D:353:PHE:HE2	1:D:364:TRP:HE1	1.66	0.41
1:B:426:ALA:HB1	2:B:501:FAD:HM83	2.02	0.41
1:C:426:ALA:HB1	1:C:427:TRP:CE3	2.55	0.41
1:D:385:LYS:HD2	1:D:418:THR:HB	2.01	0.41
1:D:398:ARG:O	1:D:402:LYS:HG3	2.20	0.41
1:A:345:VAL:HG21	1:A:407:VAL:HB	2.01	0.41
1:D:141:MSE:HE2	1:D:360:GLN:CD	2.41	0.41
1:B:56:VAL:HG13	1:B:303:VAL:HG13	2.02	0.41
1:C:353:PHE:HE1	1:C:364:TRP:HE1	1.67	0.41
1:D:212:ASN:ND2	1:D:358:GLU:OE1	2.36	0.41
1:C:173:VAL:HG11	1:C:189:LEU:HD12	2.02	0.41
1:A:314:LYS:HD2	1:A:315:HIS:CE1	2.56	0.41
1:B:351:ARG:NE	1:B:370:TYR:OH	2.53	0.41
1:A:83:GLU:HG2	1:A:88:LEU:HD23	2.02	0.41
1:C:104:PHE:CD1	1:C:415:TYR:HB2	2.55	0.41
1:C:183:THR:HG22	1:C:185:ARG:H	1.85	0.41
1:A:327:GLN:HA	1:A:330:ILE:HB	2.02	0.40
1:B:171:TRP:HA	1:B:240:TRP:CD1	2.56	0.40
1:C:259:ILE:HD12	1:C:259:ILE:HA	1.97	0.40
1:B:114:PRO:HB3	1:D:114:PRO:HA	2.03	0.40
1:B:97:PHE:CD1	1:B:413:ALA:HB2	2.56	0.40
1:D:92:THR:OG1	1:D:106:GLY:O	2.36	0.40
1:D:173:VAL:HG11	1:D:189:LEU:HD12	2.03	0.40
1:A:388:ILE:HD12	1:A:399:GLU:HG3	2.04	0.40
1:C:309:PRO:HG2	1:C:423:SER:HA	2.04	0.40
1:D:311:ASN:OD1	1:D:334:GLN:N	2.39	0.40
1:A:420:ASP:HB3	1:A:423:SER:OG	2.21	0.40
1:D:287:GLY:O	1:D:447:ARG:NE	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:462:ASN:HA	2:D:501:FAD:N1	2.36	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	430/490 (88%)	418 (97%)	8 (2%)	4 (1%)	21	37
1	B	430/490 (88%)	419 (97%)	9 (2%)	2 (0%)	34	55
1	C	433/490 (88%)	421 (97%)	10 (2%)	2 (0%)	34	55
1	D	433/490 (88%)	420 (97%)	11 (2%)	2 (0%)	34	55
All	All	1726/1960 (88%)	1678 (97%)	38 (2%)	10 (1%)	30	50

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	133	THR
1	B	286	ASN
1	C	394	ASP
1	D	482	SER
1	C	393	ARG
1	A	132	LEU
1	A	330	ILE
1	D	131	PRO
1	B	285	VAL
1	A	285	VAL

### 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	347/380 (91%)	338 (97%)	9 (3%)	54	81
1	B	347/380 (91%)	338 (97%)	9 (3%)	54	81
1	C	350/380 (92%)	342 (98%)	8 (2%)	58	83
1	D	350/380 (92%)	340 (97%)	10 (3%)	50	77
All	All	1394/1520 (92%)	1358 (97%)	36 (3%)	54	81

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	184	GLU
1	A	226	PHE
1	A	324	LYS
1	A	326	LYS
1	A	331	LYS
1	A	372	ASP
1	A	402	LYS
1	A	436	SER
1	A	481	LEU
1	B	96	ARG
1	B	205	ARG
1	B	211	ILE
1	B	225	LYS
1	B	226	PHE
1	B	275	MSE
1	B	286	ASN
1	B	326	LYS
1	B	355	PHE
1	C	82	LEU
1	C	153	ASP
1	C	217	LEU
1	C	226	PHE
1	C	436	SER
1	C	482	SER

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Mol	Chain	Res	Type
1	C	483	LEU
1	C	484	GLU
1	D	52	PHE
1	D	92	THR
1	D	205	ARG
1	D	221	GLU
1	D	226	PHE
1	D	246	MSE
1	D	322	LEU
1	D	351	ARG
1	D	481	LEU
1	D	483	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	284	GLN
1	B	255	GLN
1	B	360	GLN
1	C	110	HIS
1	C	143	ASN
1	C	485	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](#)

4 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	501	-	52,58,58	4.39	22 (42%)	52,89,89	2.43	10 (19%)
2	FAD	B	501	-	52,58,58	4.40	22 (42%)	52,89,89	2.48	10 (19%)
2	FAD	C	501	-	52,58,58	4.40	22 (42%)	52,89,89	2.42	10 (19%)
2	FAD	D	501	-	52,58,58	4.42	22 (42%)	52,89,89	2.47	10 (19%)

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	501	-	-	0/30/50/50	0/6/6/6
2	FAD	B	501	-	-	0/30/50/50	0/6/6/6
2	FAD	C	501	-	-	0/30/50/50	0/6/6/6
2	FAD	D	501	-	-	0/30/50/50	0/6/6/6

All (88) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	FAD	C2B-C1B	-14.93	1.29	1.53
2	C	501	FAD	C2B-C1B	-14.89	1.30	1.53
2	A	501	FAD	C2B-C1B	-14.83	1.30	1.53
2	D	501	FAD	C2B-C1B	-14.82	1.30	1.53
2	B	501	FAD	O4B-C4B	-6.03	1.31	1.45
2	D	501	FAD	O4B-C4B	-5.99	1.31	1.45
2	C	501	FAD	O4B-C4B	-5.98	1.31	1.45
2	A	501	FAD	O4B-C4B	-5.90	1.31	1.45
2	A	501	FAD	O2'-C2'	-2.82	1.37	1.43
2	D	501	FAD	O3B-C3B	-2.82	1.36	1.43
2	B	501	FAD	O3B-C3B	-2.78	1.36	1.43
2	D	501	FAD	C5A-C4A	-2.76	1.34	1.40
2	B	501	FAD	C5A-C4A	-2.76	1.34	1.40
2	C	501	FAD	O3B-C3B	-2.75	1.36	1.43
2	D	501	FAD	O2'-C2'	-2.71	1.37	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	FAD	C5A-C4A	-2.71	1.34	1.40
2	A	501	FAD	C5A-C4A	-2.68	1.34	1.40
2	B	501	FAD	O2'-C2'	-2.63	1.37	1.43
2	A	501	FAD	O3B-C3B	-2.62	1.36	1.43
2	C	501	FAD	O2'-C2'	-2.56	1.37	1.43
2	C	501	FAD	O4-C4	-2.38	1.18	1.24
2	B	501	FAD	O4-C4	-2.37	1.18	1.24
2	D	501	FAD	O4-C4	-2.31	1.18	1.24
2	A	501	FAD	O4-C4	-2.28	1.18	1.24
2	C	501	FAD	O4'-C4'	-2.24	1.38	1.43
2	D	501	FAD	O4'-C4'	-2.21	1.38	1.43
2	A	501	FAD	O4'-C4'	-2.17	1.38	1.43
2	B	501	FAD	O4'-C4'	-2.13	1.38	1.43
2	B	501	FAD	C2A-N3A	2.27	1.36	1.32
2	D	501	FAD	C2A-N3A	2.33	1.36	1.32
2	C	501	FAD	C2A-N3A	2.39	1.36	1.32
2	D	501	FAD	C4X-C10	2.41	1.45	1.40
2	A	501	FAD	C4X-C10	2.44	1.45	1.40
2	B	501	FAD	C4X-C10	2.53	1.45	1.40
2	A	501	FAD	C2A-N3A	2.54	1.36	1.32
2	C	501	FAD	C4X-C10	2.57	1.45	1.40
2	C	501	FAD	O2B-C2B	2.85	1.49	1.43
2	A	501	FAD	C8M-C8	2.87	1.56	1.51
2	A	501	FAD	O2B-C2B	2.93	1.49	1.43
2	D	501	FAD	O2B-C2B	2.96	1.49	1.43
2	B	501	FAD	C8M-C8	2.98	1.56	1.51
2	C	501	FAD	C8M-C8	2.98	1.56	1.51
2	B	501	FAD	O2B-C2B	3.01	1.50	1.43
2	D	501	FAD	C8M-C8	3.17	1.57	1.51
2	B	501	FAD	C6A-N6A	3.27	1.47	1.34
2	C	501	FAD	C6A-N6A	3.30	1.47	1.34
2	A	501	FAD	C6A-N6A	3.32	1.47	1.34
2	D	501	FAD	C6A-N6A	3.33	1.47	1.34
2	D	501	FAD	C2-N1	5.01	1.48	1.38
2	B	501	FAD	C2-N1	5.06	1.48	1.38
2	C	501	FAD	C2-N1	5.11	1.48	1.38
2	A	501	FAD	C2-N1	5.15	1.48	1.38
2	C	501	FAD	C4-C4X	5.16	1.51	1.41
2	A	501	FAD	C4-C4X	5.23	1.51	1.41
2	B	501	FAD	C4-C4X	5.34	1.52	1.41
2	D	501	FAD	C4-C4X	5.35	1.52	1.41
2	A	501	FAD	C2-N3	5.35	1.49	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	FAD	C2-N3	5.35	1.49	1.38
2	D	501	FAD	C2-N3	5.52	1.49	1.38
2	B	501	FAD	C2-N3	5.54	1.49	1.38
2	A	501	FAD	C10-N10	5.55	1.45	1.39
2	B	501	FAD	C10-N10	5.74	1.45	1.39
2	D	501	FAD	C10-N10	5.88	1.46	1.39
2	C	501	FAD	C10-N10	6.23	1.46	1.39
2	A	501	FAD	C10-N1	6.45	1.46	1.35
2	D	501	FAD	C10-N1	6.51	1.46	1.35
2	C	501	FAD	C10-N1	6.57	1.46	1.35
2	B	501	FAD	C10-N1	6.58	1.46	1.35
2	C	501	FAD	C4-N3	7.13	1.45	1.33
2	A	501	FAD	C4-N3	7.20	1.46	1.33
2	B	501	FAD	C4-N3	7.35	1.46	1.33
2	D	501	FAD	C4-N3	7.46	1.46	1.33
2	B	501	FAD	C9A-N10	8.28	1.50	1.38
2	C	501	FAD	C9A-N10	8.38	1.50	1.38
2	D	501	FAD	C9A-N10	8.46	1.50	1.38
2	A	501	FAD	C9A-N10	8.50	1.50	1.38
2	D	501	FAD	C4X-N5	8.50	1.46	1.33
2	A	501	FAD	C4X-N5	8.51	1.46	1.33
2	C	501	FAD	C4X-N5	8.54	1.46	1.33
2	B	501	FAD	C4X-N5	8.58	1.46	1.33
2	A	501	FAD	C5X-N5	9.27	1.49	1.35
2	C	501	FAD	C5X-N5	9.34	1.50	1.35
2	B	501	FAD	C5X-N5	9.41	1.50	1.35
2	D	501	FAD	C5X-N5	9.52	1.50	1.35
2	B	501	FAD	O4B-C1B	14.13	1.61	1.41
2	C	501	FAD	O4B-C1B	14.21	1.61	1.41
2	D	501	FAD	O4B-C1B	14.32	1.61	1.41
2	A	501	FAD	O4B-C1B	14.42	1.61	1.41

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	FAD	N3A-C2A-N1A	-11.45	119.88	128.87
2	D	501	FAD	N3A-C2A-N1A	-11.39	119.92	128.87
2	A	501	FAD	N3A-C2A-N1A	-11.05	120.19	128.87
2	C	501	FAD	N3A-C2A-N1A	-11.04	120.19	128.87
2	B	501	FAD	N6A-C6A-N1A	-6.48	107.65	118.52
2	D	501	FAD	N6A-C6A-N1A	-6.15	108.20	118.52
2	A	501	FAD	N6A-C6A-N1A	-6.02	108.41	118.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	FAD	N6A-C6A-N1A	-5.71	108.93	118.52
2	D	501	FAD	N3-C2-N1	-4.73	119.72	127.69
2	A	501	FAD	N3-C2-N1	-4.71	119.75	127.69
2	B	501	FAD	N3-C2-N1	-4.69	119.78	127.69
2	C	501	FAD	N3-C2-N1	-4.54	120.05	127.69
2	D	501	FAD	C4B-O4B-C1B	-4.28	105.10	109.64
2	C	501	FAD	C7M-C7-C6	-3.85	109.46	120.33
2	C	501	FAD	C4B-O4B-C1B	-3.79	105.62	109.64
2	D	501	FAD	C7M-C7-C6	-3.74	109.77	120.33
2	B	501	FAD	C7M-C7-C6	-3.73	109.79	120.33
2	A	501	FAD	C7M-C7-C6	-3.52	110.38	120.33
2	B	501	FAD	C4B-O4B-C1B	-3.38	106.06	109.64
2	A	501	FAD	C4X-C4-N3	-3.27	119.25	123.52
2	A	501	FAD	C4B-O4B-C1B	-3.22	106.23	109.64
2	D	501	FAD	C4X-C4-N3	-3.09	119.48	123.52
2	C	501	FAD	C4X-C4-N3	-3.03	119.57	123.52
2	B	501	FAD	C4X-C4-N3	-3.03	119.57	123.52
2	B	501	FAD	C4X-N5-C5X	2.59	119.77	116.72
2	C	501	FAD	C4X-N5-C5X	2.65	119.84	116.72
2	A	501	FAD	C4X-N5-C5X	2.68	119.88	116.72
2	D	501	FAD	C4X-N5-C5X	2.80	120.02	116.72
2	D	501	FAD	C5X-C9A-N10	2.94	119.78	117.58
2	A	501	FAD	C5X-C9A-N10	3.43	120.15	117.58
2	B	501	FAD	C5X-C9A-N10	3.47	120.18	117.58
2	C	501	FAD	C5X-C9A-N10	3.59	120.27	117.58
2	A	501	FAD	C7M-C7-C8	4.21	129.79	120.73
2	D	501	FAD	C7M-C7-C8	4.30	129.97	120.73
2	B	501	FAD	C7M-C7-C8	4.44	130.27	120.73
2	C	501	FAD	C7M-C7-C8	4.52	130.45	120.73
2	C	501	FAD	C4-N3-C2	5.36	119.63	115.16
2	B	501	FAD	C4-N3-C2	5.55	119.78	115.16
2	D	501	FAD	C4-N3-C2	5.56	119.80	115.16
2	A	501	FAD	C4-N3-C2	5.96	120.13	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	FAD	4	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	FAD	4	0
2	C	501	FAD	4	0
2	D	501	FAD	5	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

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	423/490 (86%)	-0.11	7 (1%) 73 76	12, 26, 51, 71	0
1	B	423/490 (86%)	-0.10	8 (1%) 70 73	12, 27, 51, 75	0
1	C	426/490 (86%)	-0.04	6 (1%) 78 80	13, 26, 51, 77	0
1	D	426/490 (86%)	-0.02	7 (1%) 74 78	14, 34, 57, 75	0
All	All	1698/1960 (86%)	-0.07	28 (1%) 74 78	12, 28, 53, 77	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	132	LEU	4.6
1	C	132	LEU	3.8
1	A	286	ASN	3.8
1	B	286	ASN	3.6
1	A	132	LEU	3.5
1	C	484	GLU	3.1
1	D	483	LEU	3.0
1	B	156	GLY	2.9
1	D	131	PRO	2.8
1	C	134	ASN	2.7
1	B	203	LEU	2.6
1	A	51	GLY	2.5
1	A	373	GLU	2.5
1	C	135	LEU	2.4
1	B	285	VAL	2.3
1	C	153	ASP	2.3
1	D	372	ASP	2.2
1	B	205	ARG	2.2
1	C	286	ASN	2.2
1	D	132	LEU	2.2
1	A	134	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	154	GLU	2.2
1	D	133	THR	2.1
1	A	351	ARG	2.1
1	B	155	PHE	2.1
1	B	140	ILE	2.1
1	D	153	ASP	2.0
1	D	482	SER	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, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	FAD	A	501	53/53	0.96	0.17	1.09	10,14,23,25	0
2	FAD	B	501	53/53	0.96	0.14	0.73	9,15,22,25	0
2	FAD	D	501	53/53	0.96	0.14	0.21	14,22,28,30	0
2	FAD	C	501	53/53	0.96	0.14	-0.07	9,15,23,27	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.