



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 1, 2021 – 02:57 AM EDT

PDB ID : 2REH  
Title : Mechanistic and Structural Analyses of the Roles of Arg409 and Asp402 in the Reaction of the Flavoprotein Nitroalkane Oxidase  
Authors : Fitzpatrick, P.F.; Bozinovski, D.M.; Heroux, A.; Shaw, P.G.; Valley, M.P.; Orville, A.M.  
Deposited on : 2007-09-26  
Resolution : 2.40 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.23.2  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.23.2

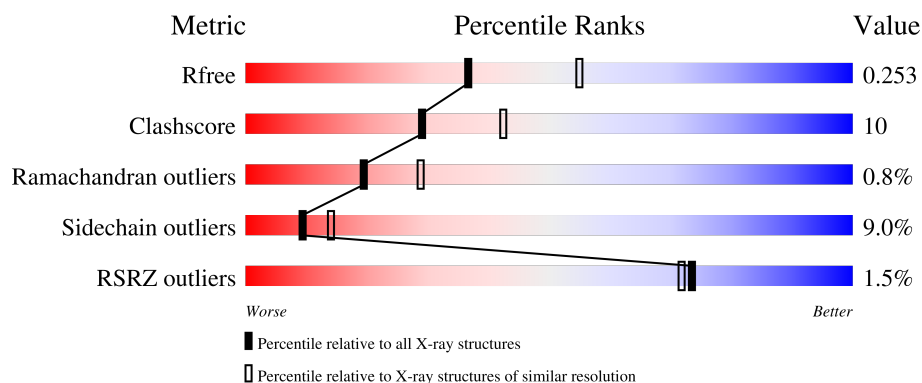
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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))
$R_{free}$	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. 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	439	<div> <div>%</div> <div> <div></div> <div>76%</div> <div>19%</div> <div>..</div> </div> </div>
1	B	439	<div> <div>4%</div> <div> <div></div> <div>65%</div> <div>27%</div> <div>5%</div> <div>..</div> </div> </div>
1	C	439	<div> <div></div> <div> <div></div> <div>77%</div> <div>17%</div> <div>..</div> </div> </div>
1	D	439	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>17%</div> <div>..</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13472 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 Nitroalkane oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	430	Total	C	N	O	S	0	0	0
			3302	2093	564	625	20			
1	B	430	Total	C	N	O	S	0	0	0
			3298	2091	564	623	20			
1	C	430	Total	C	N	O	S	0	0	0
			3302	2093	564	625	20			
1	D	430	Total	C	N	O	S	0	0	0
			3302	2093	564	625	20			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	402	GLU	ASP	engineered mutation	UNP Q8X1D8
B	402	GLU	ASP	engineered mutation	UNP Q8X1D8
C	402	GLU	ASP	engineered mutation	UNP Q8X1D8
D	402	GLU	ASP	engineered mutation	UNP Q8X1D8

- 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		

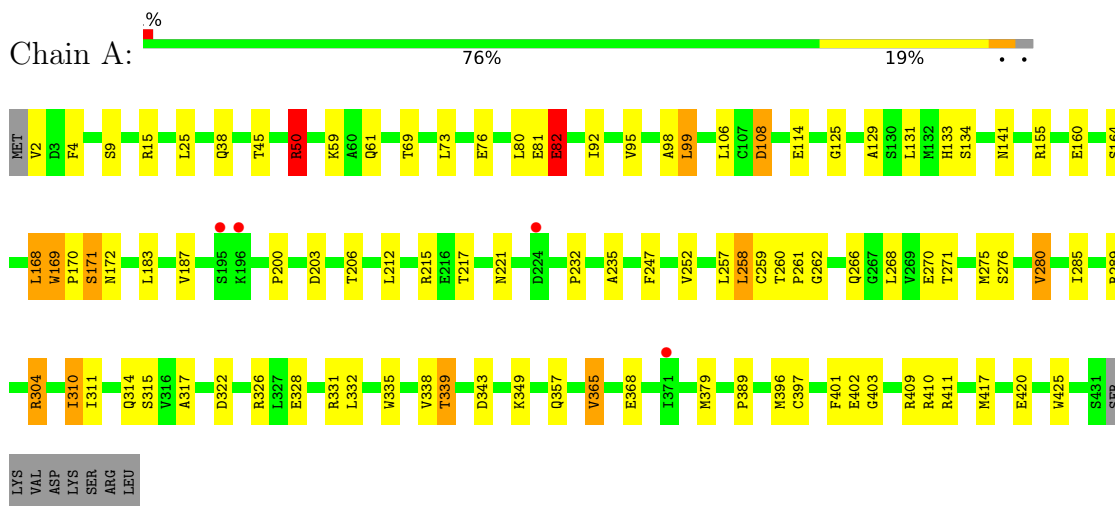
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	O	0	0
			11	11		
3	B	8	Total	O	0	0
			8	8		
3	C	16	Total	O	0	0
			16	16		
3	D	21	Total	O	0	0
			21	21		

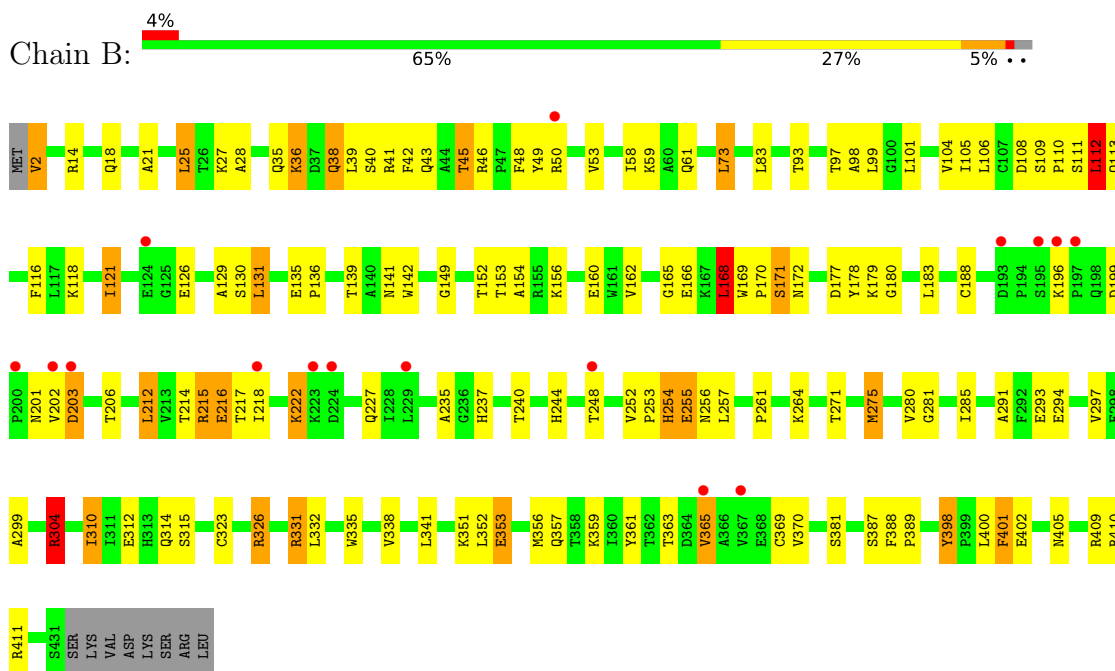
### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Nitroalkane oxidase

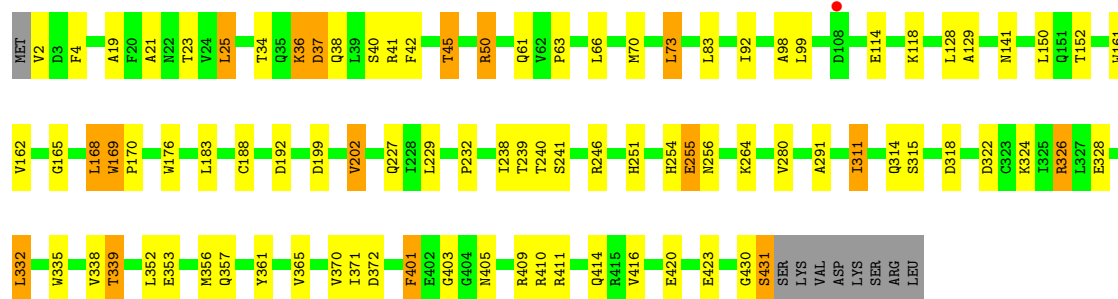


#### • Molecule 1: Nitroalkane oxidase




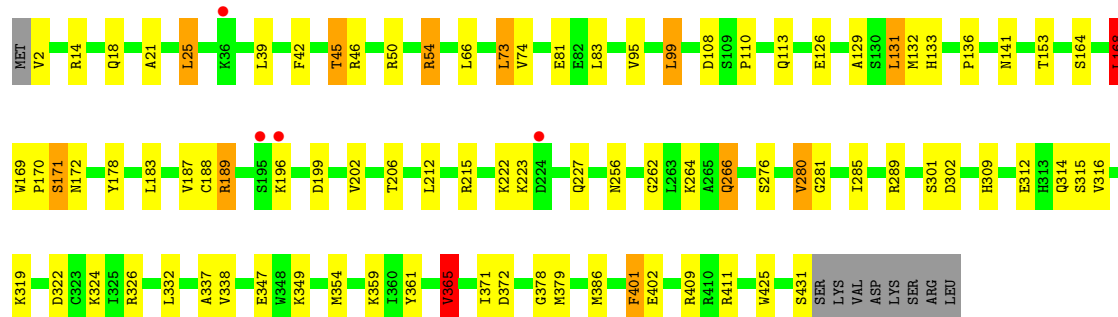
#### • Molecule 1: Nitroalkane oxidase

Chain C:  77% 17% . .



• Molecule 1: Nitroalkane oxidase

Chain D:  78% 17% . .



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	108.93Å 108.93Å 337.67Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.40 49.03 – 2.40	Depositor EDS
% Data completeness (in resolution range)	88.1 (50.00-2.40) 88.1 (49.03-2.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.43 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.204 , 0.263 0.199 , 0.253	Depositor DCC
$R_{free}$ test set	4051 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	50.2	Xtriage
Anisotropy	0.021	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 27.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.019 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	13472	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.46% 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 [i](#)

### 5.1 Standard geometry [i](#)

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	1.01	4/3371 (0.1%)	1.01	10/4575 (0.2%)
1	B	1.06	4/3367 (0.1%)	1.00	11/4570 (0.2%)
1	C	0.99	2/3371 (0.1%)	0.95	5/4575 (0.1%)
1	D	0.98	0/3371	0.94	6/4575 (0.1%)
All	All	1.01	10/13480 (0.1%)	0.98	32/18295 (0.2%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	216	GLU	CG-CD	7.34	1.62	1.51
1	A	114	GLU	CG-CD	6.72	1.62	1.51
1	B	216	GLU	CB-CG	6.17	1.63	1.52
1	B	188	CYS	CB-SG	-5.87	1.72	1.81
1	A	368	GLU	CG-CD	5.67	1.60	1.51
1	C	188	CYS	CB-SG	-5.67	1.72	1.81
1	A	247	PHE	CE2-CZ	5.38	1.47	1.37
1	C	420	GLU	CG-CD	5.30	1.59	1.51
1	B	398	TYR	CD1-CE1	5.23	1.47	1.39
1	A	420	GLU	CG-CD	5.12	1.59	1.51

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	410	ARG	NE-CZ-NH1	10.96	125.78	120.30
1	A	410	ARG	NE-CZ-NH2	-10.71	114.94	120.30
1	A	331	ARG	NE-CZ-NH2	-10.21	115.20	120.30
1	B	304	ARG	NE-CZ-NH1	8.74	124.67	120.30
1	B	326	ARG	NE-CZ-NH2	-7.83	116.39	120.30
1	B	331	ARG	NE-CZ-NH2	-7.63	116.49	120.30
1	B	310	ILE	CA-CB-CG2	7.26	125.43	110.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	112	LEU	CA-CB-CG	7.00	131.40	115.30
1	A	417	MET	CG-SD-CE	6.75	111.01	100.20
1	A	215	ARG	NE-CZ-NH2	-6.52	117.04	120.30
1	C	332	LEU	CA-CB-CG	6.31	129.81	115.30
1	B	310	ILE	CB-CA-C	6.30	124.20	111.60
1	B	331	ARG	NE-CZ-NH1	6.25	123.43	120.30
1	D	46	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	C	37	ASP	CB-CG-OD1	5.95	123.65	118.30
1	B	326	ARG	NE-CZ-NH1	5.88	123.24	120.30
1	B	168	LEU	CA-CB-CG	5.84	128.73	115.30
1	A	50	ARG	NE-CZ-NH1	5.82	123.21	120.30
1	D	46	ARG	NE-CZ-NH1	5.81	123.20	120.30
1	D	365	VAL	CG1-CB-CG2	5.45	119.62	110.90
1	C	150	LEU	CB-CG-CD1	-5.38	101.86	111.00
1	A	396	MET	CG-SD-CE	5.32	108.72	100.20
1	D	189	ARG	NE-CZ-NH1	5.25	122.93	120.30
1	C	326	ARG	NE-CZ-NH2	-5.25	117.68	120.30
1	A	343	ASP	CB-CG-OD1	5.21	122.98	118.30
1	A	215	ARG	NE-CZ-NH1	5.13	122.86	120.30
1	C	229	LEU	CA-CB-CG	5.11	127.06	115.30
1	B	304	ARG	NE-CZ-NH2	-5.10	117.75	120.30
1	A	106	LEU	CB-CG-CD1	-5.08	102.36	111.00
1	D	168	LEU	CA-CB-CG	5.06	126.94	115.30
1	B	131	LEU	CA-CB-CG	5.05	126.92	115.30
1	D	386	MET	CG-SD-CE	5.02	108.23	100.20

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3302	0	3310	74	0
1	B	3298	0	3306	102	0
1	C	3302	0	3310	59	0
1	D	3302	0	3310	51	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	53	0	31	5	0
2	B	53	0	31	7	0
2	C	53	0	31	6	0
2	D	53	0	31	1	0
3	A	11	0	0	0	0
3	B	8	0	0	0	0
3	C	16	0	0	0	0
3	D	21	0	0	0	0
All	All	13472	0	13360	258	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 (258) 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:215:ARG:HG2	1:B:215:ARG:HH11	1.00	1.15
1:B:156:LYS:HD2	1:B:160:GLU:O	1.51	1.10
1:D:54:ARG:HG2	1:D:54:ARG:HH11	1.18	1.04
1:B:215:ARG:HG2	1:B:215:ARG:NH1	1.69	0.99
1:A:379:MET:HE3	1:B:401:PHE:HA	1.42	0.98
1:A:280:VAL:HG21	1:A:402:GLU:HG2	1.50	0.93
1:A:81:GLU:O	1:A:82:GLU:HB2	1.67	0.92
1:A:168:LEU:C	1:A:170:PRO:HD3	1.91	0.91
1:B:215:ARG:HH11	1:B:215:ARG:CG	1.84	0.89
1:C:141:ASN:HD21	2:C:500:FAD:H61A	1.21	0.88
2:C:500:FAD:HM81	1:D:379:MET:HE2	1.56	0.88
1:A:379:MET:HE1	1:B:400:LEU:C	1.96	0.85
1:B:141:ASN:HD21	2:B:500:FAD:H61A	1.21	0.85
2:C:500:FAD:HM81	1:D:379:MET:CE	2.09	0.82
1:D:141:ASN:HD21	2:D:500:FAD:H61A	1.27	0.82
1:A:326:ARG:HB3	1:A:365:VAL:HG13	1.59	0.82
1:C:38:GLN:HG3	1:C:232:PRO:O	1.85	0.77
1:B:168:LEU:C	1:B:170:PRO:HD3	2.07	0.75
1:D:54:ARG:HG2	1:D:54:ARG:NH1	1.96	0.74
1:A:389:PRO:HB3	1:B:235:ALA:HB2	1.70	0.74
1:A:108:ASP:OD1	1:A:108:ASP:N	2.19	0.74
1:A:379:MET:HE3	1:B:401:PHE:CA	2.18	0.74
1:C:326:ARG:HB3	1:C:365:VAL:HG13	1.70	0.73
1:B:35:GLN:HE21	1:B:40:SER:HB3	1.53	0.73
1:D:14:ARG:O	1:D:18:GLN:HG3	1.90	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:GLN:HG3	1:A:232:PRO:O	1.90	0.72
1:C:61:GLN:HE21	1:C:98:ALA:HB2	1.54	0.72
1:A:141:ASN:HD21	2:A:500:FAD:H61A	1.39	0.70
1:B:141:ASN:ND2	2:B:500:FAD:H61A	1.90	0.69
1:B:314:GLN:HE21	1:C:315:SER:H	1.38	0.69
1:C:168:LEU:O	1:C:170:PRO:HD3	1.93	0.69
1:C:92:ILE:HG21	1:C:240:THR:HG23	1.73	0.69
1:D:99:LEU:HD22	1:D:131:LEU:HD23	1.75	0.68
1:C:73:LEU:HB3	1:C:338:VAL:HB	1.76	0.68
1:A:314:GLN:HB2	1:D:314:GLN:HB2	1.76	0.67
1:B:53:VAL:HG22	1:B:58:ILE:HG13	1.77	0.67
1:B:326:ARG:HB3	1:B:365:VAL:HG13	1.76	0.67
1:A:271:THR:O	1:A:275:MET:HG3	1.96	0.66
1:D:326:ARG:HB3	1:D:365:VAL:HG13	1.78	0.66
1:A:4:PHE:HD1	1:C:339:THR:HG21	1.61	0.66
1:C:322:ASP:O	1:C:326:ARG:HG3	1.95	0.65
1:D:73:LEU:HB3	1:D:338:VAL:HB	1.77	0.65
1:A:81:GLU:O	1:A:82:GLU:CB	2.38	0.65
1:B:299:ALA:HA	1:B:310:ILE:HG22	1.78	0.65
1:C:199:ASP:HB3	1:C:202:VAL:HG13	1.79	0.65
2:A:500:FAD:O5B	1:B:304:ARG:HG2	1.97	0.65
1:A:168:LEU:O	1:A:170:PRO:HD3	1.95	0.64
1:B:359:LYS:HD2	1:B:409:ARG:HG3	1.79	0.63
1:B:253:PRO:HB2	1:B:255:GLU:HG2	1.79	0.63
1:C:168:LEU:C	1:C:170:PRO:HD3	2.19	0.62
1:B:314:GLN:HB2	1:C:314:GLN:HB2	1.82	0.62
1:A:314:GLN:HE21	1:D:315:SER:H	1.48	0.62
1:B:291:ALA:HB2	1:B:370:VAL:HG22	1.82	0.62
1:B:109:SER:CB	1:B:112:LEU:HD22	2.30	0.61
1:B:168:LEU:O	1:B:170:PRO:HD3	2.00	0.61
1:D:81:GLU:OE1	1:D:289:ARG:NH1	2.32	0.61
1:A:379:MET:CE	1:B:400:LEU:C	2.68	0.61
1:A:61:GLN:NE2	1:A:76:GLU:OE2	2.32	0.61
1:D:199:ASP:HB3	1:D:202:VAL:HG12	1.83	0.60
1:B:353:GLU:O	1:B:357:GLN:HG3	2.01	0.60
1:D:54:ARG:HH11	1:D:54:ARG:CG	2.04	0.60
1:A:317:ALA:CB	1:C:414:GLN:HG3	2.32	0.60
1:D:322:ASP:O	1:D:326:ARG:HG3	2.00	0.60
1:B:73:LEU:HB3	1:B:338:VAL:HB	1.84	0.60
1:A:59:LYS:HE3	1:A:69:THR:HG23	1.83	0.59
1:D:280:VAL:HG21	1:D:402:GLU:HG2	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:281:GLY:O	1:B:285:ILE:HG13	2.03	0.59
1:C:326:ARG:NH2	1:C:372:ASP:OD2	2.36	0.59
1:B:21:ALA:HA	1:B:25:LEU:HB2	1.84	0.58
1:B:405:ASN:O	1:B:410:ARG:HG3	2.03	0.58
1:C:37:ASP:OD1	1:C:40:SER:OG	2.15	0.58
1:B:257:LEU:HD21	1:B:261:PRO:HD3	1.85	0.57
1:A:328:GLU:HG2	1:C:357:GLN:HE22	1.70	0.57
1:B:271:THR:O	1:B:275:MET:HG2	2.04	0.57
1:A:322:ASP:O	1:A:326:ARG:HG3	2.03	0.57
1:D:326:ARG:NH2	1:D:372:ASP:OD2	2.38	0.57
1:A:141:ASN:ND2	2:A:500:FAD:H61A	2.04	0.56
1:A:315:SER:H	1:D:314:GLN:HE21	1.53	0.56
1:B:109:SER:HB3	1:B:112:LEU:HD22	1.88	0.56
1:B:177:ASP:OD1	1:B:215:ARG:NH2	2.39	0.56
1:C:423:GLU:OE1	1:C:431:SER:HB3	2.06	0.55
1:B:180:GLY:HA3	1:B:218:ILE:CD1	2.35	0.55
1:B:237:HIS:O	1:B:240:THR:HG22	2.07	0.55
1:D:50:ARG:HD2	1:D:126:GLU:OE1	2.06	0.55
1:A:171:SER:O	1:A:172:ASN:HB2	2.05	0.55
1:B:42:PHE:O	1:B:45:THR:HB	2.07	0.55
1:C:141:ASN:ND2	2:C:500:FAD:H61A	1.98	0.55
1:B:14:ARG:O	1:B:18:GLN:HB2	2.07	0.55
1:B:199:ASP:O	1:B:202:VAL:HG12	2.06	0.54
1:C:50:ARG:HH11	1:C:50:ARG:HB2	1.71	0.54
1:B:361:TYR:O	1:B:365:VAL:HB	2.08	0.54
1:C:168:LEU:C	1:C:170:PRO:CD	2.76	0.54
1:B:341:LEU:HD23	1:B:351:LYS:HB3	1.89	0.54
1:A:379:MET:CE	1:B:400:LEU:O	2.57	0.53
1:C:371:ILE:HD11	1:D:371:ILE:HD11	1.91	0.53
1:B:61:GLN:HE21	1:B:98:ALA:HB2	1.74	0.53
1:A:317:ALA:HB2	1:C:414:GLN:HG3	1.91	0.53
1:B:129:ALA:HA	1:B:183:LEU:O	2.08	0.53
1:B:315:SER:H	1:C:314:GLN:HE21	1.55	0.53
1:A:50:ARG:NH2	1:A:125:GLY:O	2.23	0.52
1:A:258:LEU:HD21	1:A:268:LEU:HD13	1.92	0.52
1:B:203:ASP:HB3	1:B:206:THR:HG23	1.92	0.52
1:C:361:TYR:O	1:C:365:VAL:HB	2.10	0.52
1:A:425:TRP:CZ3	1:C:324:LYS:HE2	2.45	0.52
1:B:35:GLN:NE2	1:B:40:SER:HB3	2.23	0.52
1:B:109:SER:O	1:B:113:GLN:HG3	2.09	0.52
1:C:405:ASN:O	1:C:410:ARG:HG3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:GLN:HE21	1:B:41:ARG:HH21	1.58	0.51
1:D:199:ASP:HB3	1:D:202:VAL:CG1	2.40	0.51
1:C:255:GLU:HG2	1:C:255:GLU:O	2.08	0.51
1:D:309:HIS:O	1:D:312:GLU:HB2	2.11	0.51
1:B:110:PRO:HA	1:B:113:GLN:HE21	1.76	0.51
1:C:61:GLN:HE21	1:C:98:ALA:CB	2.20	0.51
1:B:401:PHE:C	1:B:401:PHE:CD2	2.84	0.51
1:B:49:TYR:O	1:B:53:VAL:HG23	2.11	0.50
1:C:352:LEU:O	1:C:356:MET:HG2	2.11	0.50
1:B:50:ARG:HD2	1:B:126:GLU:OE1	2.11	0.50
1:B:291:ALA:CB	1:B:370:VAL:HG22	2.40	0.50
1:C:401:PHE:C	1:C:401:PHE:CD2	2.84	0.50
1:A:168:LEU:HD13	1:A:169:TRP:NE1	2.27	0.50
1:A:379:MET:SD	2:B:500:FAD:HM81	2.52	0.50
1:D:110:PRO:HA	1:D:113:GLN:HE21	1.77	0.49
1:C:335:TRP:O	1:C:339:THR:HB	2.12	0.49
1:B:2:VAL:HG11	1:D:74:VAL:HG11	1.95	0.49
1:A:169:TRP:N	1:A:170:PRO:CD	2.76	0.49
1:B:28:ALA:HA	1:B:48:PHE:CZ	2.48	0.49
1:A:235:ALA:HB2	1:B:389:PRO:HB3	1.95	0.48
1:A:379:MET:HE1	1:B:400:LEU:O	2.13	0.48
1:A:379:MET:HG2	2:B:500:FAD:H3'	1.94	0.48
2:A:500:FAD:O2P	2:A:500:FAD:H52A	2.13	0.48
1:A:169:TRP:N	1:A:170:PRO:HD3	2.26	0.48
1:A:99:LEU:HD22	1:A:131:LEU:HD23	1.94	0.48
1:B:39:LEU:O	1:B:40:SER:C	2.50	0.48
1:B:46:ARG:NE	1:B:126:GLU:OE2	2.45	0.48
1:C:169:TRP:N	1:C:170:PRO:CD	2.77	0.47
1:B:152:THR:HA	1:B:165:GLY:HA3	1.95	0.47
1:C:92:ILE:CG2	1:C:240:THR:CG2	2.92	0.47
1:B:43:GLN:C	1:B:45:THR:H	2.17	0.47
1:B:136:PRO:HA	1:B:168:LEU:HD12	1.95	0.47
1:B:199:ASP:OD1	1:B:201:ASN:HB2	2.13	0.47
2:C:500:FAD:H52A	1:D:378:GLY:HA3	1.95	0.47
1:A:134:SER:HB2	1:A:168:LEU:O	2.14	0.47
1:A:335:TRP:O	1:A:339:THR:HB	2.15	0.47
1:B:315:SER:OG	1:C:318:ASP:OD1	2.27	0.47
1:C:21:ALA:HA	1:C:25:LEU:HB2	1.96	0.47
1:B:108:ASP:O	1:B:110:PRO:HD3	2.14	0.46
1:B:156:LYS:CD	1:B:160:GLU:O	2.43	0.46
1:D:401:PHE:C	1:D:401:PHE:CD2	2.88	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:92:ILE:HG21	1:C:240:THR:CG2	2.41	0.46
2:C:500:FAD:C8M	1:D:379:MET:CE	2.88	0.46
1:B:353:GLU:OE1	1:D:324:LYS:HE3	2.15	0.46
1:A:80:LEU:HD22	1:A:95:VAL:CG1	2.45	0.46
1:A:257:LEU:HG	1:A:259:CYS:O	2.15	0.46
1:A:257:LEU:HD21	1:A:261:PRO:HD3	1.97	0.46
1:A:304:ARG:HB2	1:A:310:ILE:CG1	2.46	0.46
1:A:129:ALA:HA	1:A:183:LEU:O	2.15	0.46
1:B:93:THR:HG23	1:B:172:ASN:OD1	2.16	0.46
1:B:179:LYS:CB	1:B:215:ARG:HH12	2.28	0.46
1:B:35:GLN:HB3	1:B:40:SER:HB2	1.98	0.46
1:B:154:ALA:HA	1:B:162:VAL:O	2.15	0.46
1:C:129:ALA:HA	1:C:183:LEU:O	2.16	0.46
1:D:349:LYS:HE2	1:D:349:LYS:HB2	1.81	0.46
1:A:160:GLU:HB2	1:A:252:VAL:O	2.16	0.45
1:A:217:THR:O	1:A:221:ASN:ND2	2.41	0.45
1:B:341:LEU:HD23	1:B:341:LEU:HA	1.78	0.45
1:A:304:ARG:HG2	2:B:500:FAD:O5B	2.16	0.45
1:B:331:ARG:NH2	1:D:425:TRP:CE3	2.80	0.45
1:B:357:GLN:HE21	1:B:357:GLN:HB3	1.51	0.45
1:D:54:ARG:NH1	1:D:54:ARG:CG	2.69	0.45
1:B:217:THR:HG21	1:B:252:VAL:HG22	1.99	0.45
1:D:337:ALA:HA	1:D:354:MET:HG3	1.98	0.45
1:A:4:PHE:CD1	1:C:339:THR:HG21	2.47	0.45
1:A:304:ARG:HA	1:A:304:ARG:NE	2.32	0.45
1:C:50:ARG:O	1:C:50:ARG:HG3	2.17	0.45
1:B:183:LEU:HD11	1:B:212:LEU:HG	1.99	0.44
1:C:353:GLU:HB2	1:C:416:VAL:HG11	1.99	0.44
1:B:352:LEU:O	1:B:356:MET:HG2	2.18	0.44
1:D:378:GLY:O	1:D:379:MET:C	2.55	0.44
1:B:61:GLN:HE21	1:B:98:ALA:CB	2.30	0.44
1:B:180:GLY:HA3	1:B:218:ILE:HD12	1.99	0.44
1:B:294:GLU:OE1	1:B:387:SER:HB2	2.17	0.44
1:C:42:PHE:CD2	1:C:241:SER:HB2	2.53	0.44
1:A:80:LEU:HD21	1:A:95:VAL:HG12	1.98	0.44
1:C:38:GLN:HE22	1:C:238:ILE:HA	1.83	0.44
1:A:183:LEU:HD11	1:A:212:LEU:HG	1.99	0.44
1:A:203:ASP:HB3	1:A:206:THR:HG23	2.00	0.44
1:B:36:LYS:HA	1:B:36:LYS:HD3	1.75	0.44
1:C:326:ARG:HB3	1:C:365:VAL:CG1	2.43	0.44
1:A:80:LEU:CD2	1:A:95:VAL:CG1	2.96	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:314:GLN:NE2	1:D:315:SER:H	2.13	0.44
1:B:381:SER:O	1:B:388:PHE:HB2	2.18	0.44
1:A:304:ARG:HB2	1:A:310:ILE:HG12	1.99	0.43
1:B:104:VAL:C	1:B:106:LEU:N	2.71	0.43
1:C:152:THR:HA	1:C:165:GLY:HA3	1.99	0.43
1:D:129:ALA:HA	1:D:183:LEU:O	2.18	0.43
1:A:310:ILE:HD13	1:A:310:ILE:HA	1.33	0.43
1:B:293:GLU:O	1:B:297:VAL:HG23	2.18	0.43
1:D:39:LEU:HD22	1:D:178:TYR:CZ	2.53	0.43
1:D:276:SER:OG	1:D:402:GLU:OE1	2.27	0.43
1:D:402:GLU:OE2	1:D:402:GLU:HA	2.19	0.43
1:A:206:THR:C	1:A:262:GLY:HA2	2.39	0.43
1:B:315:SER:H	1:C:314:GLN:NE2	2.15	0.43
1:C:36:LYS:HB2	1:C:36:LYS:HE2	1.21	0.43
1:D:21:ALA:HA	1:D:25:LEU:HB2	2.00	0.43
1:D:171:SER:O	1:D:172:ASN:HB2	2.19	0.43
1:A:379:MET:CE	2:B:500:FAD:H3'	2.48	0.43
1:A:339:THR:HG21	1:C:4:PHE:HD1	1.83	0.42
1:C:41:ARG:HB3	1:C:239:THR:HG22	2.01	0.42
1:C:42:PHE:O	1:C:45:THR:HB	2.19	0.42
1:A:397:CYS:SG	2:A:500:FAD:HM83	2.59	0.42
1:A:73:LEU:HB3	1:A:338:VAL:HB	2.00	0.42
1:B:166:GLU:OE2	1:B:244:HIS:NE2	2.50	0.42
1:D:316:VAL:O	1:D:319:LYS:HB2	2.20	0.42
1:B:109:SER:HB3	1:B:112:LEU:CD2	2.49	0.42
1:C:19:ALA:O	1:C:23:THR:HG23	2.19	0.42
1:D:153:THR:HA	1:D:189:ARG:O	2.19	0.42
1:A:379:MET:HB2	1:A:379:MET:HE2	1.73	0.42
1:B:73:LEU:CB	1:B:338:VAL:HB	2.50	0.42
1:D:42:PHE:O	1:D:45:THR:HB	2.19	0.42
1:B:171:SER:O	1:B:172:ASN:HB2	2.20	0.42
1:C:161:TRP:HD1	1:C:254:HIS:HA	1.84	0.42
1:A:276:SER:OG	1:A:402:GLU:OE1	2.19	0.42
1:B:222:LYS:HB2	1:B:222:LYS:HE3	1.78	0.42
1:C:291:ALA:HB2	1:C:370:VAL:HG22	2.02	0.42
1:D:132:MET:HA	1:D:170:PRO:HB3	2.01	0.42
1:A:357:GLN:HE22	1:C:328:GLU:HG2	1.85	0.42
1:B:335:TRP:HA	1:B:338:VAL:HG22	2.01	0.42
1:B:400:LEU:HA	1:B:400:LEU:HD23	1.66	0.41
1:D:133:HIS:HD2	1:D:187:VAL:HG11	1.85	0.41
1:C:291:ALA:CB	1:C:370:VAL:HG22	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:136:PRO:HA	1:D:168:LEU:HD12	2.02	0.41
1:B:256:ASN:N	1:B:256:ASN:HD22	2.18	0.41
1:A:168:LEU:C	1:A:170:PRO:CD	2.76	0.41
1:C:63:PRO:HA	1:C:70:MET:HB3	2.01	0.41
1:B:363:THR:HG22	1:B:398:TYR:HB3	2.02	0.41
1:B:39:LEU:HD22	1:B:178:TYR:CE2	2.55	0.41
1:D:281:GLY:O	1:D:285:ILE:HG13	2.21	0.41
1:A:15:ARG:HG2	1:A:15:ARG:HH11	1.85	0.41
1:D:361:TYR:O	1:D:365:VAL:HB	2.20	0.41
1:A:266:GLN:HE22	1:A:270:GLU:HB2	1.86	0.41
1:C:162:VAL:HG22	1:C:251:HIS:CE1	2.55	0.41
1:A:92:ILE:O	1:A:92:ILE:HG13	2.21	0.41
1:C:311:ILE:O	1:C:311:ILE:HG13	2.20	0.41
1:B:116:PHE:CE2	1:B:212:LEU:HD23	2.56	0.41
1:C:128:LEU:HD11	1:C:176:TRP:CE2	2.56	0.41
1:B:59:LYS:HG3	1:B:121:ILE:HA	2.04	0.40
1:B:179:LYS:HB3	1:B:215:ARG:HH12	1.86	0.40
1:D:266:GLN:NE2	1:D:266:GLN:O	2.54	0.40
1:D:359:LYS:HD2	1:D:409:ARG:HG3	2.02	0.40
1:A:61:GLN:HE21	1:A:98:ALA:HB2	1.86	0.40
1:B:104:VAL:O	1:B:106:LEU:N	2.55	0.40
1:B:156:LYS:HE3	1:B:254:HIS:HB2	2.04	0.40
1:B:323:CYS:HB3	1:B:369:CYS:SG	2.61	0.40
1:A:379:MET:HE2	2:B:500:FAD:H3'	2.02	0.40
1:B:97:THR:O	1:B:101:LEU:HG	2.21	0.40
1:D:206:THR:O	1:D:262:GLY:HA2	2.21	0.40
1:A:133:HIS:HD2	1:A:187:VAL:HG11	1.87	0.40
1:A:285:ILE:O	1:A:289:ARG:HB2	2.21	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	428/439 (98%)	411 (96%)	14 (3%)	3 (1%)	22	32
1	B	428/439 (98%)	396 (92%)	26 (6%)	6 (1%)	11	15
1	C	428/439 (98%)	412 (96%)	13 (3%)	3 (1%)	22	32
1	D	428/439 (98%)	416 (97%)	11 (3%)	1 (0%)	47	62
All	All	1712/1756 (98%)	1635 (96%)	64 (4%)	13 (1%)	19	29

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	169	TRP
1	B	169	TRP
1	C	169	TRP
1	D	169	TRP
1	A	82	GLU
1	B	105	ILE
1	B	142	TRP
1	B	254	HIS
1	B	402	GLU
1	C	430	GLY
1	A	403	GLY
1	B	149	GLY
1	C	403	GLY

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	354/364 (97%)	328 (93%)	26 (7%)	14	22
1	B	353/364 (97%)	313 (89%)	40 (11%)	6	8
1	C	354/364 (97%)	326 (92%)	28 (8%)	12	19
1	D	354/364 (97%)	321 (91%)	33 (9%)	9	13
All	All	1415/1456 (97%)	1288 (91%)	127 (9%)	9	14

All (127) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	VAL
1	A	9	SER
1	A	25	LEU
1	A	45	THR
1	A	50	ARG
1	A	82	GLU
1	A	99	LEU
1	A	108	ASP
1	A	155	ARG
1	A	164	SER
1	A	168	LEU
1	A	171	SER
1	A	200	PRO
1	A	258	LEU
1	A	260	THR
1	A	280	VAL
1	A	304	ARG
1	A	310	ILE
1	A	311	ILE
1	A	332	LEU
1	A	339	THR
1	A	349	LYS
1	A	365	VAL
1	A	401	PHE
1	A	409	ARG
1	A	411	ARG
1	B	2	VAL
1	B	25	LEU
1	B	27	LYS
1	B	36	LYS
1	B	38	GLN
1	B	45	THR
1	B	73	LEU
1	B	83	LEU
1	B	99	LEU
1	B	111	SER
1	B	112	LEU
1	B	118	LYS
1	B	121	ILE
1	B	130	SER
1	B	131	LEU
1	B	135	GLU
1	B	139	THR

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Mol	Chain	Res	Type
1	B	153	THR
1	B	168	LEU
1	B	171	SER
1	B	196	LYS
1	B	203	ASP
1	B	212	LEU
1	B	214	THR
1	B	215	ARG
1	B	216	GLU
1	B	222	LYS
1	B	227	GLN
1	B	248	THR
1	B	255	GLU
1	B	264	LYS
1	B	275	MET
1	B	280	VAL
1	B	304	ARG
1	B	312	GLU
1	B	332	LEU
1	B	353	GLU
1	B	365	VAL
1	B	401	PHE
1	B	411	ARG
1	C	2	VAL
1	C	25	LEU
1	C	34	THR
1	C	36	LYS
1	C	45	THR
1	C	50	ARG
1	C	66	LEU
1	C	73	LEU
1	C	83	LEU
1	C	99	LEU
1	C	114	GLU
1	C	118	LYS
1	C	168	LEU
1	C	192	ASP
1	C	202	VAL
1	C	227	GLN
1	C	246	ARG
1	C	255	GLU
1	C	256	ASN

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Mol	Chain	Res	Type
1	C	264	LYS
1	C	280	VAL
1	C	311	ILE
1	C	332	LEU
1	C	339	THR
1	C	401	PHE
1	C	409	ARG
1	C	411	ARG
1	C	431	SER
1	D	2	VAL
1	D	25	LEU
1	D	45	THR
1	D	54	ARG
1	D	66	LEU
1	D	73	LEU
1	D	83	LEU
1	D	95	VAL
1	D	99	LEU
1	D	108	ASP
1	D	131	LEU
1	D	164	SER
1	D	168	LEU
1	D	171	SER
1	D	188	CYS
1	D	196	LYS
1	D	212	LEU
1	D	215	ARG
1	D	222	LYS
1	D	223	LYS
1	D	227	GLN
1	D	256	ASN
1	D	264	LYS
1	D	266	GLN
1	D	280	VAL
1	D	301	SER
1	D	302	ASP
1	D	332	LEU
1	D	347	GLU
1	D	365	VAL
1	D	401	PHE
1	D	411	ARG
1	D	431	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	ASN
1	A	35	GLN
1	A	38	GLN
1	A	43	GLN
1	A	113	GLN
1	A	137	ASN
1	A	141	ASN
1	A	159	ASN
1	A	256	ASN
1	A	266	GLN
1	A	314	GLN
1	A	357	GLN
1	B	35	GLN
1	B	38	GLN
1	B	43	GLN
1	B	61	GLN
1	B	113	GLN
1	B	137	ASN
1	B	141	ASN
1	B	144	GLN
1	B	256	ASN
1	B	266	GLN
1	B	314	GLN
1	C	16	HIS
1	C	22	ASN
1	C	38	GLN
1	C	43	GLN
1	C	61	GLN
1	C	113	GLN
1	C	137	ASN
1	C	141	ASN
1	C	198	GLN
1	C	266	GLN
1	C	314	GLN
1	C	357	GLN
1	D	35	GLN
1	D	43	GLN
1	D	113	GLN
1	D	137	ASN
1	D	141	ASN
1	D	198	GLN

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Mol	Chain	Res	Type
1	D	227	GLN
1	D	251	HIS
1	D	256	ASN
1	D	266	GLN
1	D	313	HIS
1	D	314	GLN
1	D	357	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

### 5.6 Ligand geometry ⓘ

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	C	500	-	51,58,58	1.71	9 (17%)	60,89,89	2.14	14 (23%)
2	FAD	D	500	-	51,58,58	2.03	11 (21%)	60,89,89	2.33	16 (26%)
2	FAD	A	500	-	51,58,58	1.64	11 (21%)	60,89,89	2.04	16 (26%)
2	FAD	B	500	-	51,58,58	1.92	10 (19%)	60,89,89	2.51	16 (26%)

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	C	500	-	-	3/30/50/50	0/6/6/6
2	FAD	D	500	-	-	4/30/50/50	0/6/6/6
2	FAD	A	500	-	-	5/30/50/50	0/6/6/6
2	FAD	B	500	-	-	6/30/50/50	0/6/6/6

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	500	FAD	C1'-N10	8.88	1.57	1.48
2	B	500	FAD	C1'-N10	6.46	1.54	1.48
2	C	500	FAD	C4X-N5	5.35	1.41	1.33
2	B	500	FAD	C4X-N5	5.29	1.40	1.33
2	D	500	FAD	C4X-N5	5.02	1.40	1.33
2	C	500	FAD	C2A-N3A	4.97	1.40	1.32
2	D	500	FAD	C2A-N3A	4.42	1.39	1.32
2	C	500	FAD	C10-N1	4.36	1.38	1.33
2	A	500	FAD	C4X-N5	4.19	1.39	1.33
2	B	500	FAD	C10-N1	4.17	1.38	1.33
2	A	500	FAD	C10-N1	4.00	1.38	1.33
2	A	500	FAD	C4-N3	3.89	1.39	1.33
2	A	500	FAD	C2A-N3A	3.82	1.38	1.32
2	C	500	FAD	C1'-N10	3.77	1.52	1.48
2	D	500	FAD	C10-N1	3.61	1.37	1.33
2	D	500	FAD	C4-N3	3.54	1.39	1.33
2	A	500	FAD	O4B-C1B	3.38	1.45	1.41
2	B	500	FAD	C5'-C4'	3.32	1.56	1.51
2	B	500	FAD	C4-N3	3.31	1.38	1.33
2	B	500	FAD	C2A-N3A	3.26	1.37	1.32
2	A	500	FAD	C1'-N10	3.17	1.51	1.48
2	B	500	FAD	C4X-C10	3.10	1.41	1.38
2	B	500	FAD	C5X-N5	3.06	1.40	1.35
2	A	500	FAD	C2A-N1A	2.81	1.39	1.33
2	B	500	FAD	C2A-N1A	2.75	1.39	1.33
2	D	500	FAD	C4X-C10	2.63	1.41	1.38
2	A	500	FAD	C5'-C4'	2.58	1.55	1.51
2	A	500	FAD	C6-C5X	-2.58	1.37	1.41
2	C	500	FAD	C2A-N1A	2.52	1.38	1.33
2	C	500	FAD	C4X-C10	2.51	1.41	1.38
2	A	500	FAD	C2B-C1B	-2.45	1.50	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	500	FAD	C2B-C1B	-2.40	1.50	1.53
2	D	500	FAD	C2A-N1A	2.31	1.38	1.33
2	D	500	FAD	C9A-N10	2.20	1.41	1.38
2	D	500	FAD	C5X-N5	2.15	1.38	1.35
2	B	500	FAD	C6-C5X	-2.13	1.38	1.41
2	C	500	FAD	C8A-N7A	2.09	1.38	1.34
2	C	500	FAD	C9A-N10	2.09	1.41	1.38
2	D	500	FAD	C2B-C3B	-2.06	1.47	1.53
2	C	500	FAD	C4-N3	2.04	1.36	1.33
2	A	500	FAD	C5X-N5	2.03	1.38	1.35

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	500	FAD	N3A-C2A-N1A	-9.49	113.84	128.68
2	D	500	FAD	C1'-N10-C9A	9.02	125.39	118.29
2	B	500	FAD	C10-C4X-N5	-7.69	115.94	121.26
2	B	500	FAD	C4-N3-C2	7.58	121.54	115.14
2	A	500	FAD	N3A-C2A-N1A	-7.44	117.05	128.68
2	B	500	FAD	N3A-C2A-N1A	-6.94	117.83	128.68
2	D	500	FAD	C1'-N10-C10	-6.41	112.67	118.41
2	B	500	FAD	C1'-N10-C9A	5.91	122.94	118.29
2	A	500	FAD	C4-N3-C2	5.67	119.93	115.14
2	D	500	FAD	C4-N3-C2	5.21	119.54	115.14
2	D	500	FAD	N3A-C2A-N1A	-5.20	120.55	128.68
2	C	500	FAD	C4X-N5-C5X	5.06	121.83	116.77
2	B	500	FAD	C4X-N5-C5X	4.74	121.51	116.77
2	B	500	FAD	C4X-C4-N3	-4.61	117.13	123.43
2	C	500	FAD	C10-C4X-N5	-4.56	118.11	121.26
2	B	500	FAD	C4-C4X-N5	4.50	123.74	118.60
2	A	500	FAD	C10-C4X-N5	-4.33	118.26	121.26
2	D	500	FAD	C10-C4X-N5	-4.31	118.28	121.26
2	C	500	FAD	C4-N3-C2	4.07	118.58	115.14
2	C	500	FAD	C1'-N10-C9A	4.01	121.45	118.29
2	D	500	FAD	O2'-C2'-C1'	3.97	119.15	109.59
2	C	500	FAD	C5X-C9A-N10	3.69	120.39	117.72
2	D	500	FAD	C4X-N5-C5X	3.63	120.39	116.77
2	A	500	FAD	O4'-C4'-C5'	-3.43	102.20	109.92
2	D	500	FAD	C9A-N10-C10	-3.24	117.66	121.91
2	A	500	FAD	C4X-N5-C5X	3.18	119.95	116.77
2	A	500	FAD	C4-C4X-N5	3.14	122.19	118.60
2	C	500	FAD	O2A-PA-O5B	3.11	122.19	107.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	FAD	O2B-C2B-C3B	-3.10	101.79	111.82
2	C	500	FAD	O2'-C2'-C1'	3.07	116.98	109.59
2	A	500	FAD	C1'-N10-C10	-2.99	115.73	118.41
2	A	500	FAD	C1'-N10-C9A	2.87	120.55	118.29
2	B	500	FAD	C7M-C7-C8	2.82	126.52	120.74
2	B	500	FAD	C1'-N10-C10	-2.80	115.91	118.41
2	D	500	FAD	C5X-C9A-N10	2.74	119.70	117.72
2	B	500	FAD	O4'-C4'-C3'	-2.73	102.47	109.10
2	B	500	FAD	O2'-C2'-C3'	-2.63	102.72	109.10
2	C	500	FAD	C4-C4X-N5	2.59	121.56	118.60
2	A	500	FAD	O5B-C5B-C4B	2.53	117.69	108.99
2	D	500	FAD	O2B-C2B-C3B	-2.51	103.69	111.82
2	A	500	FAD	O2'-C2'-C3'	-2.45	103.15	109.10
2	A	500	FAD	C4X-C4-N3	-2.43	120.11	123.43
2	A	500	FAD	C6-C5X-N5	-2.42	116.38	119.05
2	D	500	FAD	C4-C4X-N5	2.40	121.34	118.60
2	B	500	FAD	C4X-C10-N10	2.40	122.77	120.30
2	D	500	FAD	O4'-C4'-C5'	-2.37	104.60	109.92
2	D	500	FAD	C4X-C4-N3	-2.35	120.22	123.43
2	C	500	FAD	C9A-N10-C10	-2.34	118.85	121.91
2	B	500	FAD	O2'-C2'-C1'	2.31	115.15	109.59
2	C	500	FAD	C6-C7-C8	-2.29	116.05	119.91
2	A	500	FAD	O5'-P-O1P	-2.29	100.13	109.07
2	A	500	FAD	O4B-C4B-C3B	-2.24	100.68	105.11
2	D	500	FAD	C4X-C10-N10	2.21	122.57	120.30
2	A	500	FAD	C2A-N1A-C6A	2.20	122.51	118.75
2	D	500	FAD	C7M-C7-C6	-2.15	115.21	120.34
2	C	500	FAD	C1'-N10-C10	-2.12	116.51	118.41
2	D	500	FAD	C7M-C7-C8	2.10	125.04	120.74
2	C	500	FAD	O2'-C2'-C3'	-2.08	104.04	109.10
2	B	500	FAD	O2A-PA-O5B	2.06	117.33	107.75
2	A	500	FAD	C5A-C6A-N6A	2.06	123.48	120.35
2	C	500	FAD	C2A-N1A-C6A	2.02	122.22	118.75
2	B	500	FAD	C2A-N1A-C6A	2.02	122.21	118.75

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	500	FAD	C2'-C3'-C4'-O4'
2	A	500	FAD	O3'-C3'-C4'-O4'
2	B	500	FAD	O3'-C3'-C4'-O4'

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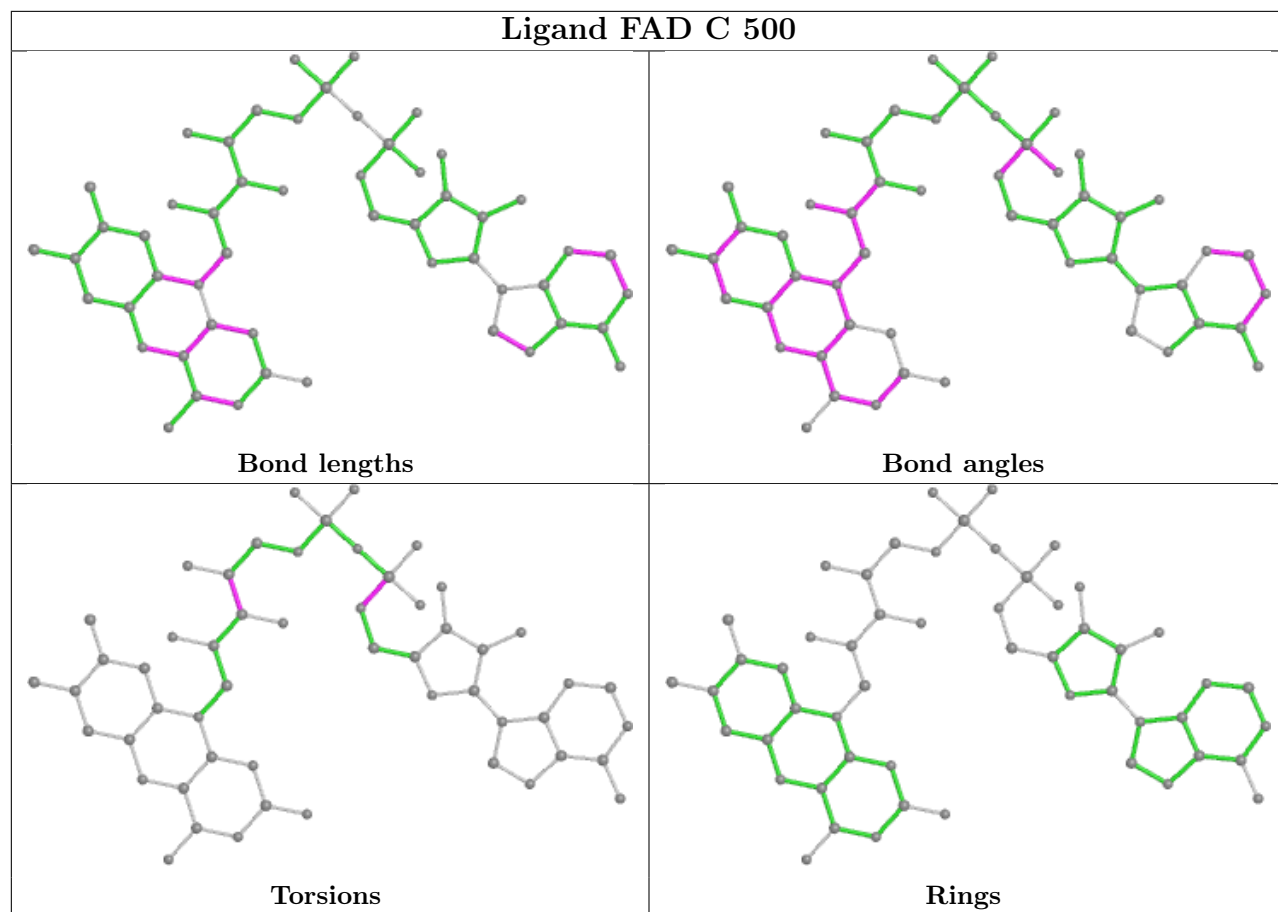
Mol	Chain	Res	Type	Atoms
2	D	500	FAD	O3'-C3'-C4'-O4'
2	A	500	FAD	O4B-C4B-C5B-O5B
2	D	500	FAD	C2'-C3'-C4'-O4'
2	A	500	FAD	C2'-C3'-C4'-O4'
2	B	500	FAD	O3'-C3'-C4'-C5'
2	B	500	FAD	C2'-C3'-C4'-C5'
2	B	500	FAD	P-O3P-PA-O1A
2	D	500	FAD	O3'-C3'-C4'-C5'
2	C	500	FAD	C5B-O5B-PA-O3P
2	C	500	FAD	C2'-C3'-C4'-O4'
2	C	500	FAD	O3'-C3'-C4'-O4'
2	A	500	FAD	O3'-C3'-C4'-C5'
2	A	500	FAD	C3B-C4B-C5B-O5B
2	D	500	FAD	C5B-O5B-PA-O3P
2	B	500	FAD	P-O3P-PA-O2A

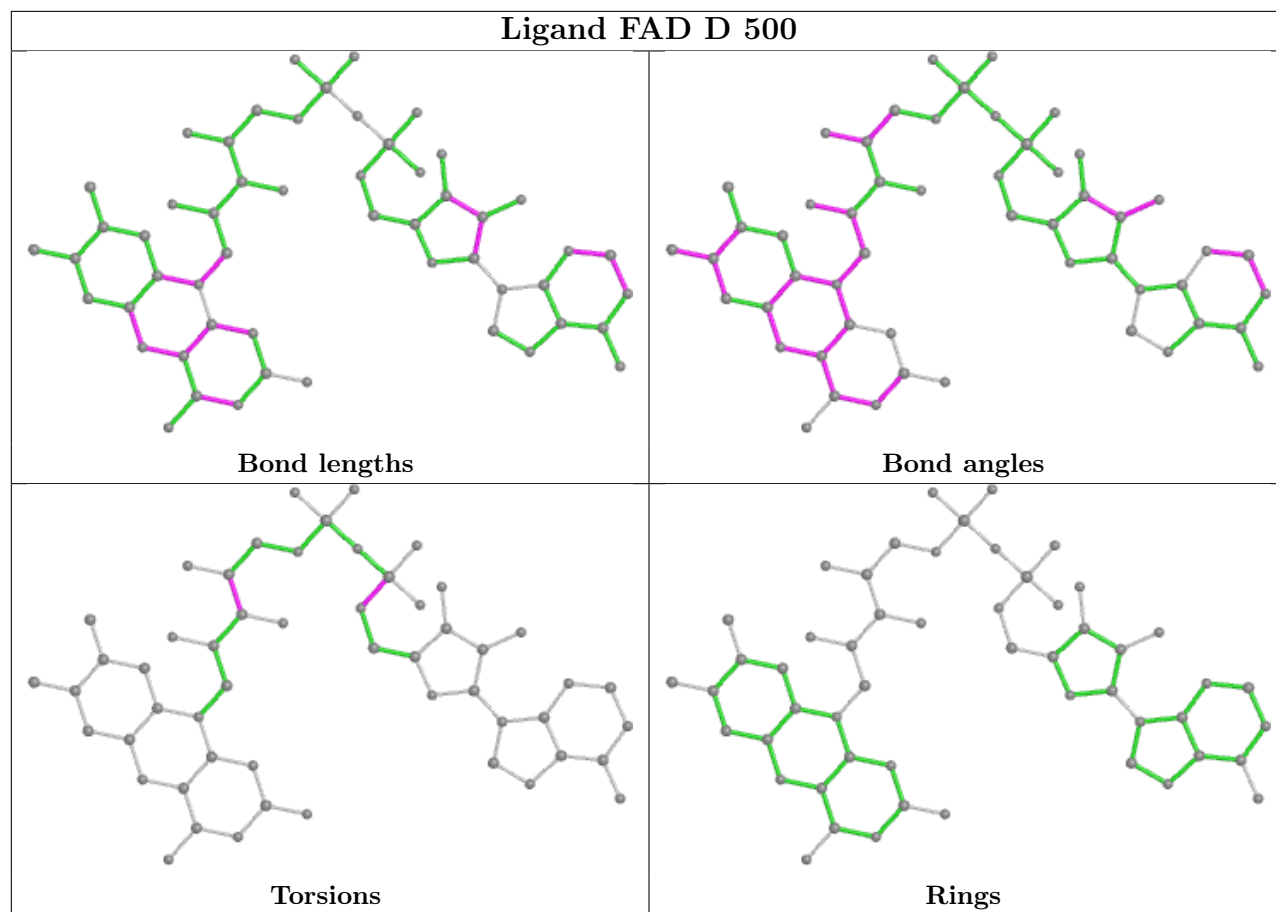
There are no ring outliers.

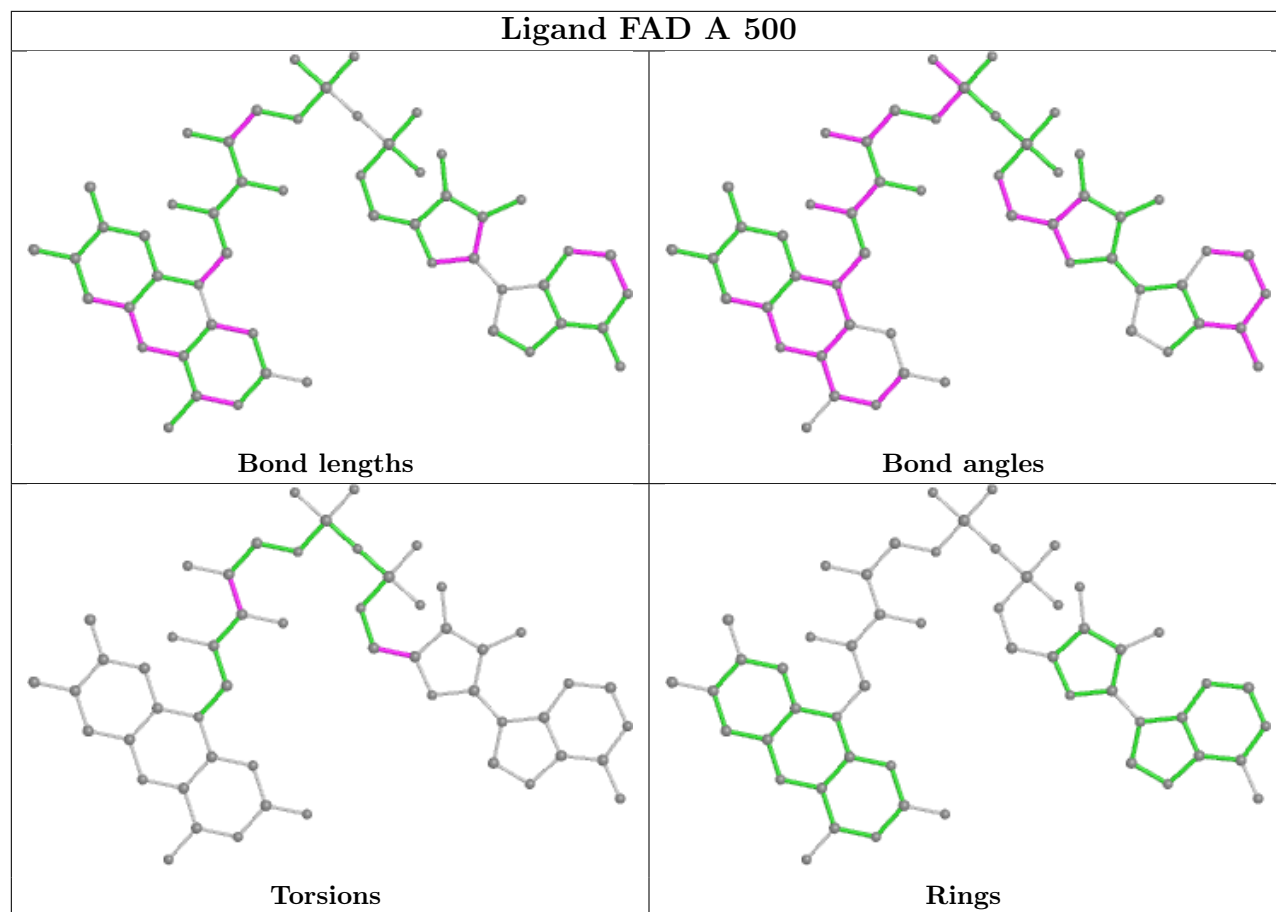
4 monomers are involved in 19 short contacts:

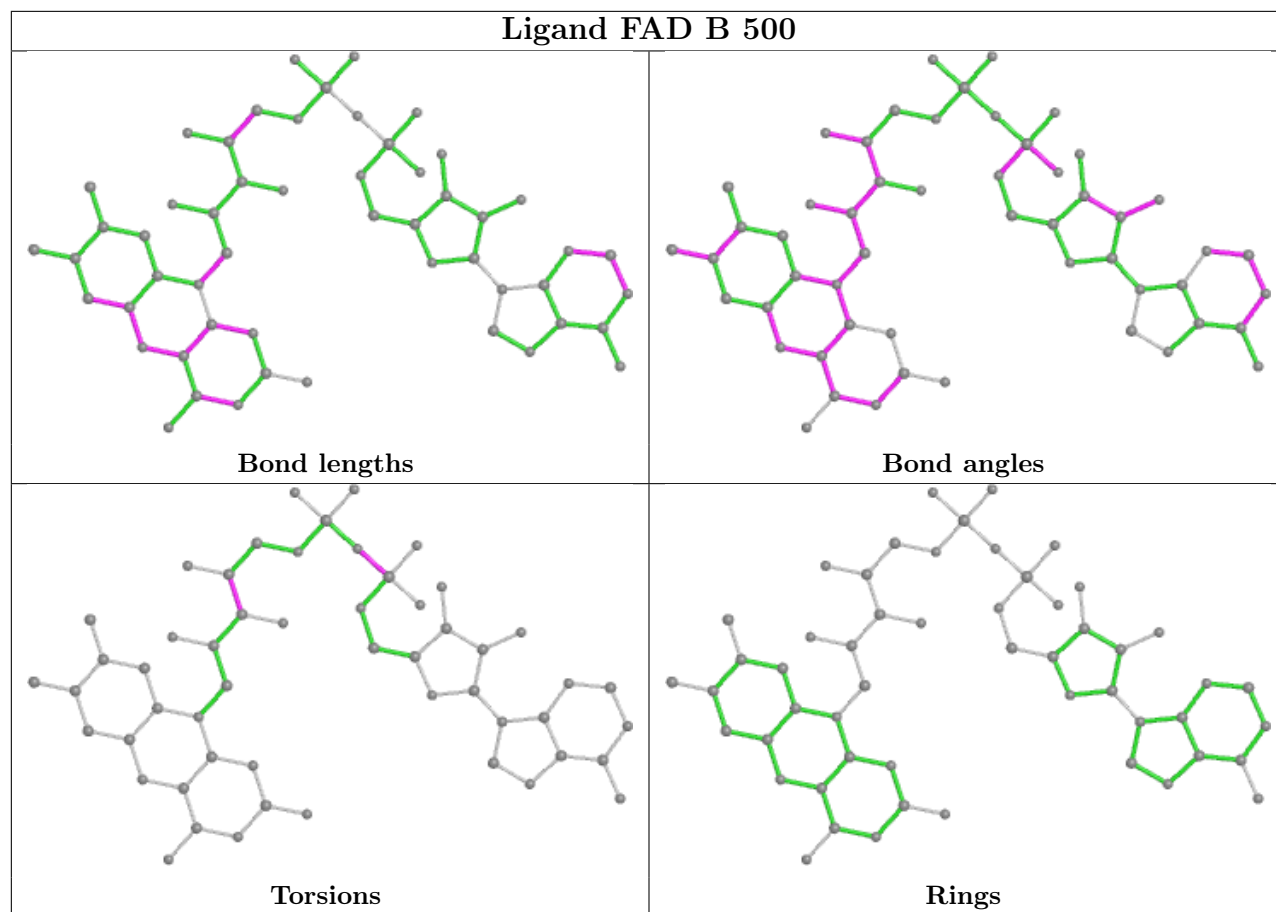
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	500	FAD	6	0
2	D	500	FAD	1	0
2	A	500	FAD	5	0
2	B	500	FAD	7	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









## 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	430/439 (97%)	-0.13	4 (0%) 84 82	31, 43, 58, 70	0
1	B	430/439 (97%)	0.06	16 (3%) 41 41	30, 44, 59, 72	0
1	C	430/439 (97%)	-0.27	1 (0%) 95 94	30, 42, 58, 71	0
1	D	430/439 (97%)	-0.22	4 (0%) 84 82	31, 42, 58, 70	1 (0%)
All	All	1720/1756 (97%)	-0.14	25 (1%) 73 72	30, 43, 58, 72	1 (0%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	202	VAL	3.7
1	A	224	ASP	3.5
1	C	108	ASP	3.4
1	B	195	SER	3.3
1	B	196	LYS	3.3
1	B	223	LYS	3.1
1	B	218	ILE	2.9
1	B	224	ASP	2.8
1	B	124	GLU	2.7
1	D	196	LYS	2.7
1	B	229	LEU	2.6
1	B	248	THR	2.5
1	B	50	ARG	2.5
1	D	36	LYS	2.4
1	A	195	SER	2.4
1	A	371	ILE	2.4
1	B	197	PRO	2.4
1	D	195	SER	2.3
1	B	365	VAL	2.2
1	B	367	VAL	2.1
1	B	203	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	193	ASP	2.0
1	B	200	PRO	2.0
1	D	224	ASP	2.0
1	A	196	LYS	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 monosaccharides 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, 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.

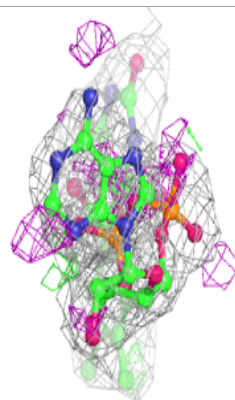
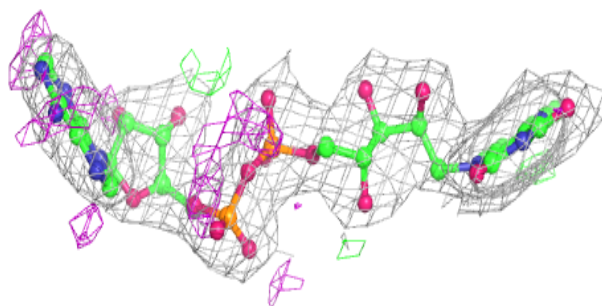
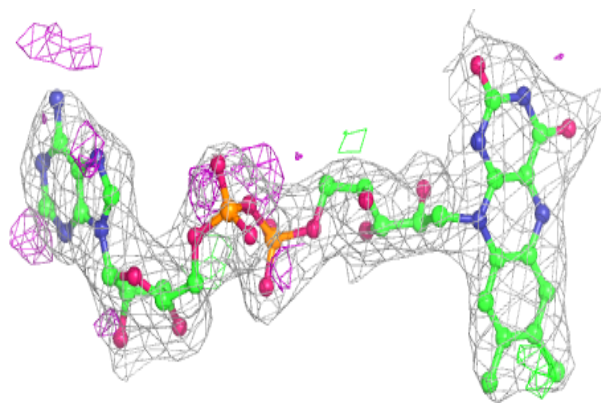
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	FAD	D	500	53/53	0.96	0.13	22,32,35,37	0
2	FAD	B	500	53/53	0.97	0.10	24,31,36,37	0
2	FAD	C	500	53/53	0.97	0.14	25,31,35,38	0
2	FAD	A	500	53/53	0.97	0.12	22,29,32,34	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

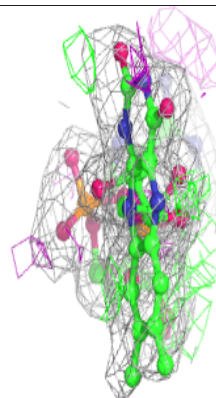
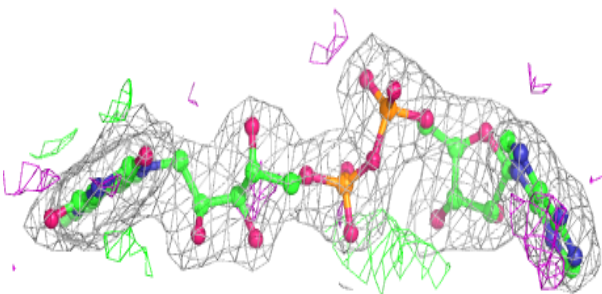
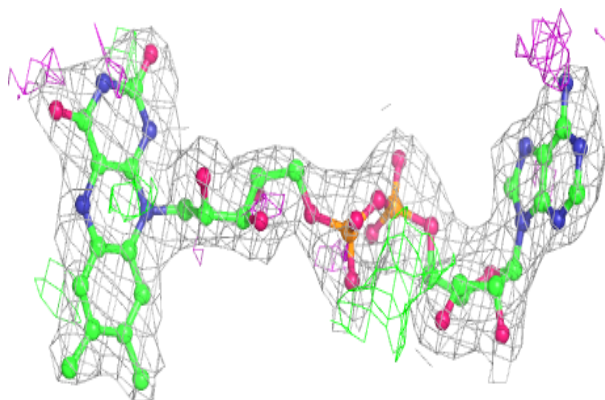


**Electron density around FAD D 500:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

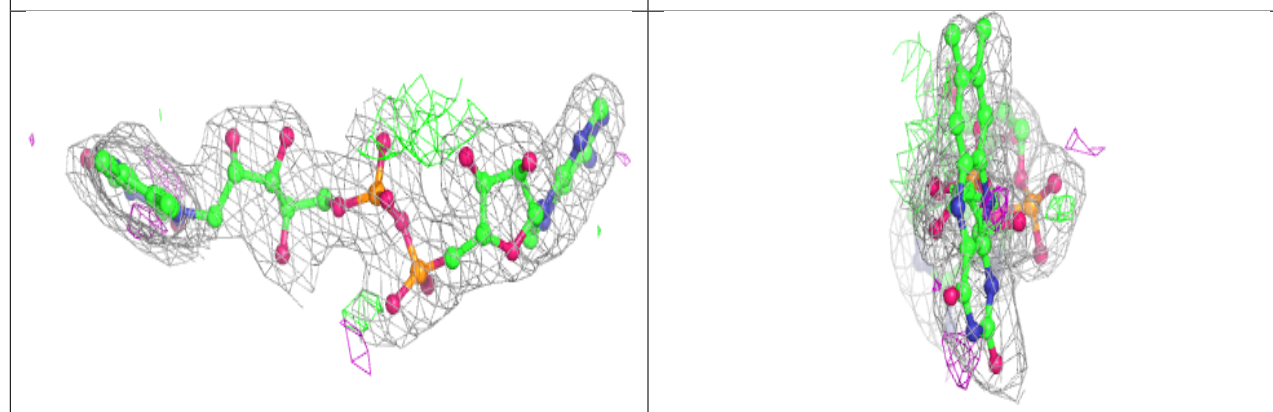
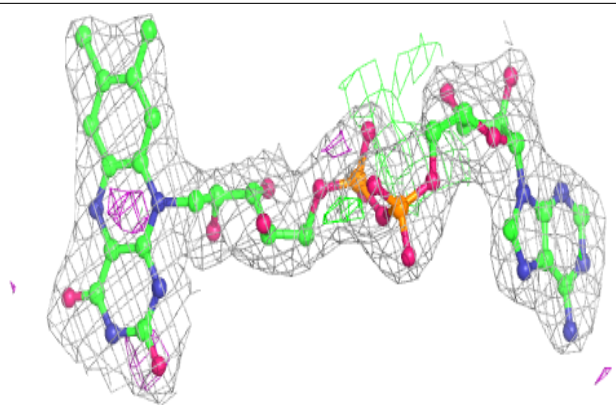
**Electron density around FAD B 500:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

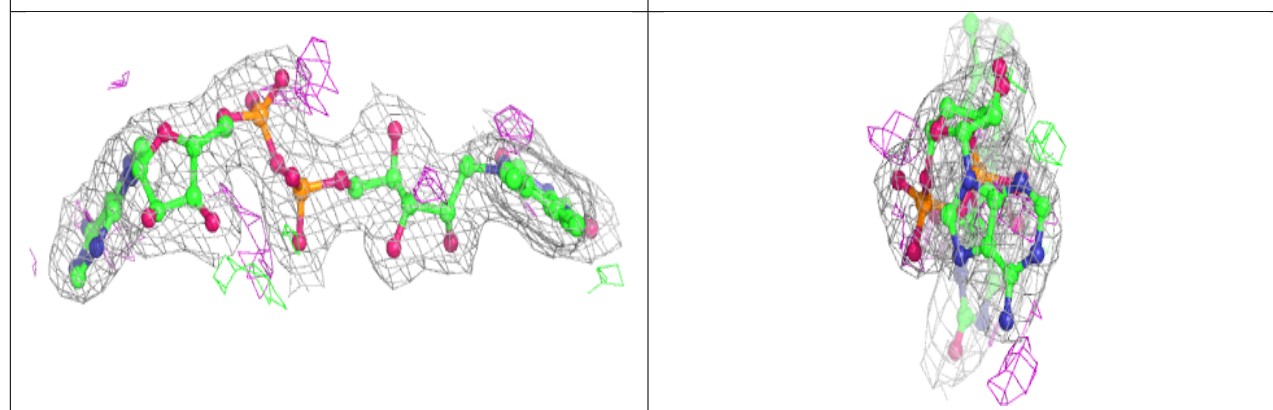
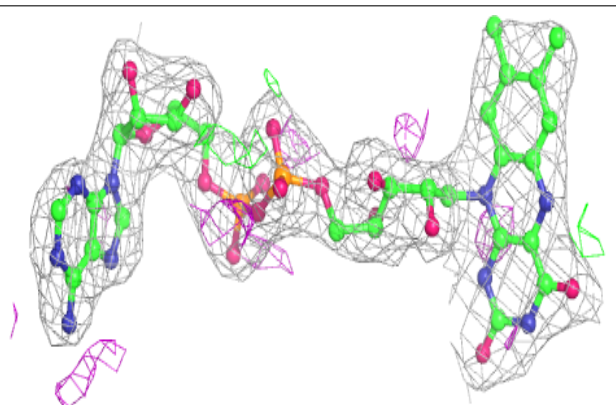


**Electron density around FAD C 500:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around FAD A 500:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.