



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:12 PM UTC

PDB ID : 5VJ7 / pdb_00005vj7
Title : Ferredoxin NADP Oxidoreductase (Xfn)
Authors : Zadvornyy, O.A.; Nguyen, D.M.N.; Schut, G.J.; Lipscomb, G.L.; Tokmina-Lukaszewska, M.; Adams, M.W.W.; Peters, J.W.
Deposited on : 2017-04-18
Resolution : 2.55 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

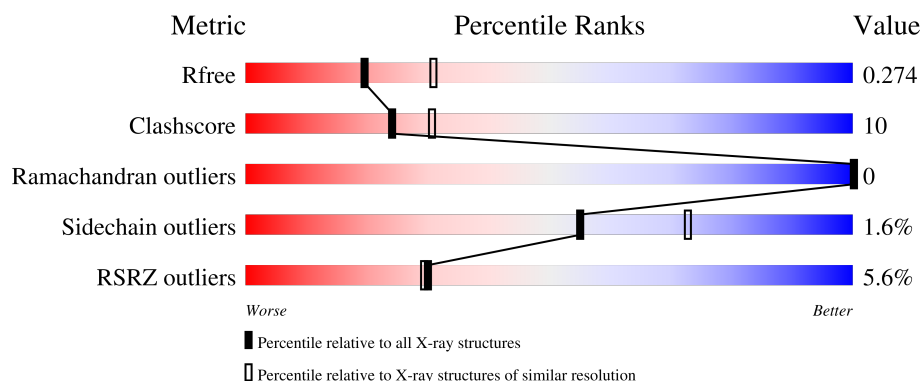
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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}	180053	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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 ≥ 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	476	3% 80% 16% ..
2	B	289	10% 78% 17% ..

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 6196 atoms, of which 62 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.

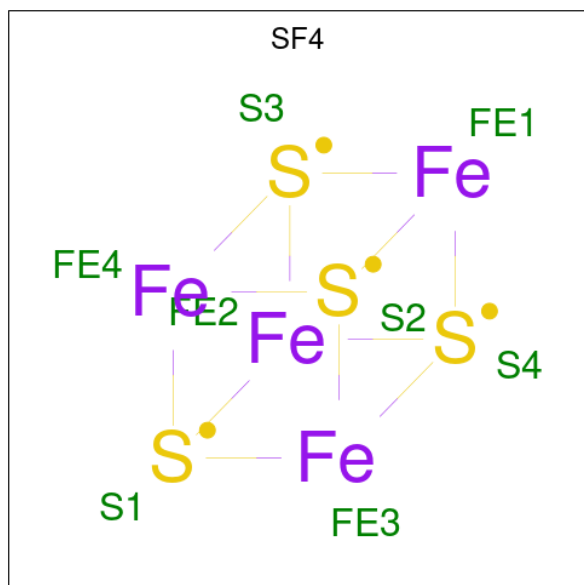
- Molecule 1 is a protein called Oxidoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	462	Total	C	N	O	S	0	0	0
			3572	2272	626	658	16			

- Molecule 2 is a protein called Ferredoxin-NADP(+) reductase subunit alpha.

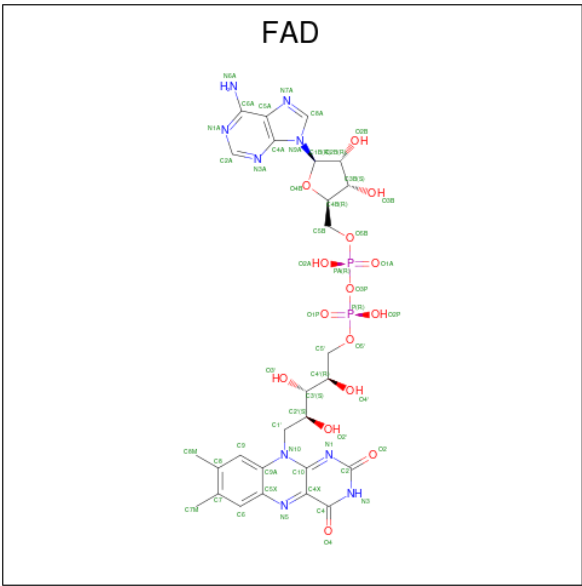
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	282	Total	C	N	O	S	0	0	0
			2250	1460	370	405	15			

- Molecule 3 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄).



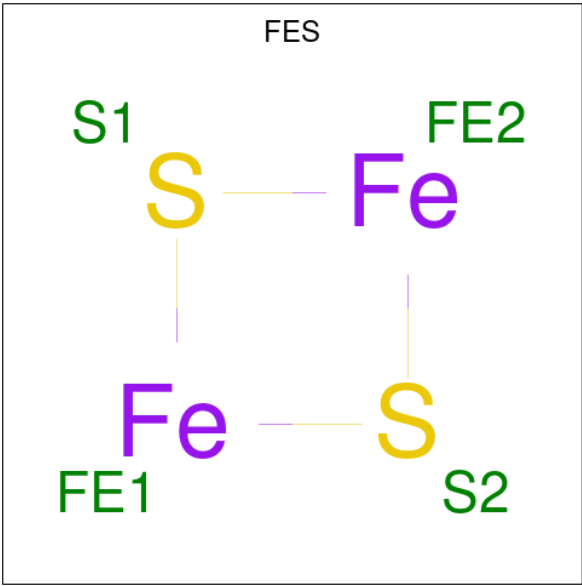
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Fe	S	0	0
			8	4	4		
3	A	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total 84	C 27	H 31	N 9	O 15	P 2	0	0
4	B	1	Total 84	C 27	H 31	N 9	O 15	P 2	0	0

- Molecule 5 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe₂S₂).



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

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total 1	Mg 1	0	0

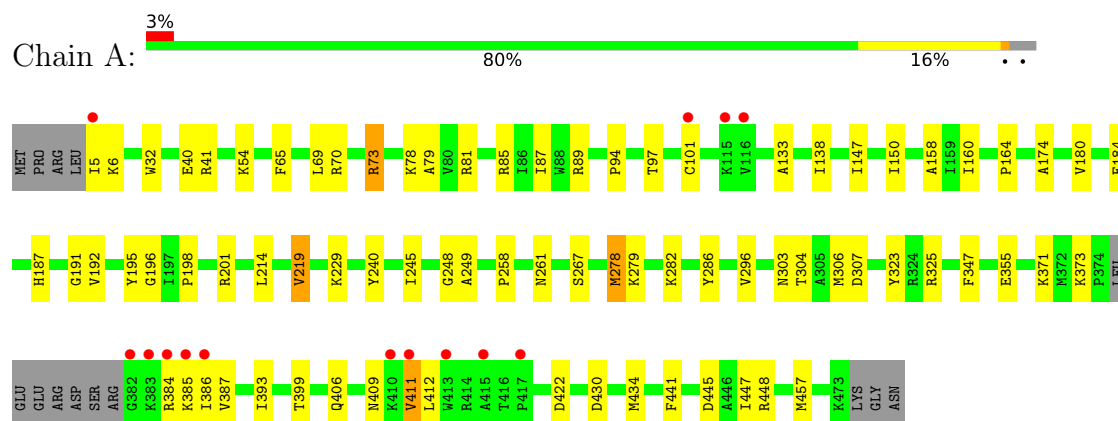
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	104	Total 104	O 104	0	0
7	B	81	Total 81	O 81	0	0

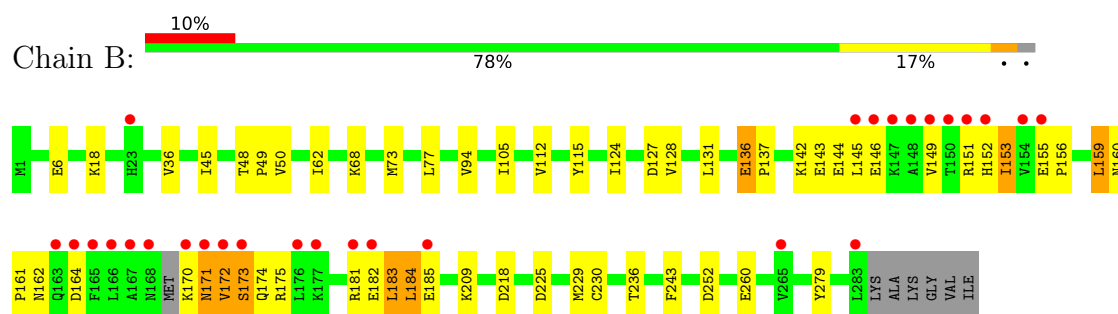
3 Residue-property plots [i](#)

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: Oxidoreductase



• Molecule 2: Ferredoxin-NADP(+) reductase subunit alpha



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.74Å 73.14Å 99.96Å 90.00° 96.73° 90.00°	Depositor
Resolution (Å)	39.00 – 2.55 39.00 – 2.55	Depositor EDS
% Data completeness (in resolution range)	98.6 (39.00-2.55) 98.6 (39.00-2.55)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.20 (at 2.54Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.183 , 0.251 0.209 , 0.274	Depositor DCC
R_{free} test set	1227 reflections (4.68%)	wwPDB-VP
Wilson B-factor (Å ²)	35.3	Xtriage
Anisotropy	1.100	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 77.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6196	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: FES, FAD, MG, SF4

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	0.91	1/3634 (0.0%)	1.38	27/4909 (0.6%)
2	B	0.90	1/2302 (0.0%)	1.41	20/3119 (0.6%)
All	All	0.91	2/5936 (0.0%)	1.39	47/8028 (0.6%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	278	MET	SD-CE	-6.01	1.64	1.79
2	B	182	GLU	CA-C	5.01	1.59	1.53

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	171	ASN	N-CA-C	-13.70	96.60	113.18
2	B	173	SER	N-CA-C	-10.85	98.76	113.30
2	B	183	LEU	N-CA-C	-10.76	99.89	113.23
2	B	152	HIS	CB-CA-C	8.17	126.46	109.38
1	A	73	ARG	CA-C-N	8.10	131.34	120.65
1	A	73	ARG	C-N-CA	8.10	131.34	120.65
2	B	184	LEU	N-CA-C	-7.71	99.07	110.48
1	A	323	TYR	CA-C-N	7.61	131.23	120.28
1	A	323	TYR	C-N-CA	7.61	131.23	120.28
2	B	182	GLU	CA-C-N	7.55	134.94	121.66
2	B	182	GLU	C-N-CA	7.55	134.94	121.66
2	B	159	LEU	CA-C-N	7.26	129.16	120.09
2	B	159	LEU	C-N-CA	7.26	129.16	120.09
1	A	267	SER	N-CA-C	-7.12	99.98	110.52
2	B	243	PHE	CA-CB-CG	6.57	120.37	113.80
2	B	183	LEU	CA-C-N	6.49	132.16	121.39
2	B	183	LEU	C-N-CA	6.49	132.16	121.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	152	HIS	N-CA-C	-6.37	99.43	109.50
1	A	430	ASP	CA-CB-CG	6.21	118.81	112.60
1	A	78	LYS	CA-C-N	6.12	128.48	120.28
1	A	78	LYS	C-N-CA	6.12	128.48	120.28
1	A	32	TRP	CA-C-N	6.01	128.25	120.44
1	A	32	TRP	C-N-CA	6.01	128.25	120.44
1	A	422	ASP	CA-C-N	5.77	128.81	120.38
1	A	422	ASP	C-N-CA	5.77	128.81	120.38
1	A	198	PRO	CA-C-N	5.76	128.58	120.28
1	A	198	PRO	C-N-CA	5.76	128.58	120.28
1	A	229	LYS	CA-C-N	5.65	127.79	120.44
1	A	229	LYS	C-N-CA	5.65	127.79	120.44
1	A	65	PHE	CA-CB-CG	5.52	119.32	113.80
1	A	411	VAL	CA-C-N	5.38	127.49	120.28
1	A	411	VAL	C-N-CA	5.38	127.49	120.28
2	B	252	ASP	CA-CB-CG	5.37	117.97	112.60
1	A	303	ASN	CA-C-N	5.36	127.46	120.28
1	A	303	ASN	C-N-CA	5.36	127.46	120.28
1	A	422	ASP	CA-CB-CG	5.35	117.95	112.60
2	B	171	ASN	CA-C-N	5.32	131.55	121.97
2	B	171	ASN	C-N-CA	5.32	131.55	121.97
1	A	457	MET	CA-C-N	5.31	125.99	119.99
1	A	457	MET	C-N-CA	5.31	125.99	119.99
1	A	180	VAL	N-CA-C	5.18	115.73	108.89
2	B	218	ASP	CA-CB-CG	5.16	117.76	112.60
2	B	153	ILE	N-CA-CB	5.04	118.67	111.82
2	B	172	VAL	CA-C-N	5.02	130.28	122.49
2	B	172	VAL	C-N-CA	5.02	130.28	122.49
1	A	174	ALA	CA-C-N	5.00	127.48	120.28
1	A	174	ALA	C-N-CA	5.00	127.48	120.28

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	3572	0	3662	58	0
2	B	2250	0	2286	65	0
3	A	16	0	0	0	0
4	A	53	31	31	12	0
4	B	53	31	31	2	0
5	B	4	0	0	1	0
6	B	1	0	0	0	0
7	A	104	0	0	0	0
7	B	81	0	0	7	0
All	All	6134	62	6010	120	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 (120) 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)
2:B:225:ASP:OD1	2:B:230:CYS:SG	1.92	1.26
2:B:160:ASN:OD1	2:B:161:PRO:HD3	1.26	1.25
2:B:142:LYS:O	2:B:146:GLU:HG3	1.36	1.21
1:A:184:GLU:OE1	4:A:503:FAD:O2B	1.64	1.12
1:A:386:ILE:HG22	1:A:387:VAL:H	1.14	1.12
2:B:115:TYR:HD1	2:B:145:LEU:HD23	1.25	1.00
2:B:145:LEU:HD13	2:B:151:ARG:HH12	1.24	1.00
1:A:191:GLY:HA2	4:A:503:FAD:O3B	1.62	0.99
1:A:386:ILE:CG2	1:A:387:VAL:H	1.74	0.99
2:B:160:ASN:OD1	2:B:161:PRO:CD	2.12	0.98
1:A:386:ILE:HG22	1:A:387:VAL:N	1.70	0.98
2:B:170:LYS:HE3	7:B:451:HOH:O	1.63	0.97
2:B:115:TYR:HD1	2:B:145:LEU:CD2	1.78	0.96
1:A:164:PRO:HD2	4:A:503:FAD:O1P	1.69	0.93
1:A:373:LYS:O	1:A:386:ILE:HG23	1.69	0.91
2:B:127:ASP:OD1	2:B:149:VAL:HG21	1.73	0.88
2:B:172:VAL:CG1	2:B:175:ARG:HG2	2.06	0.85
2:B:115:TYR:CD1	2:B:145:LEU:HD23	2.11	0.84
2:B:230:CYS:SG	5:B:302:FES:FE2	1.69	0.84
2:B:151:ARG:NH2	7:B:401:HOH:O	1.92	0.84
1:A:195:TYR:CE2	2:B:229:MET:HE1	2.13	0.83
2:B:170:LYS:CE	7:B:451:HOH:O	2.19	0.83
1:A:187:HIS:HE1	1:A:278:MET:HE1	1.42	0.82
2:B:142:LYS:HG3	2:B:153:ILE:HD11	1.62	0.81
1:A:296:VAL:HG22	1:A:399:THR:HB	1.64	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:LYS:HE2	1:A:41:ARG:HG2	1.64	0.78
2:B:145:LEU:CD1	2:B:151:ARG:HH12	1.97	0.78
2:B:172:VAL:HA	2:B:174:GLN:HG2	1.67	0.76
2:B:170:LYS:HG3	7:B:451:HOH:O	1.85	0.75
2:B:145:LEU:HD13	2:B:151:ARG:NH1	2.00	0.74
1:A:97:THR:HB	1:A:101:CYS:SG	2.29	0.73
1:A:412:LEU:HD11	4:A:503:FAD:H61A	1.54	0.72
1:A:6:LYS:HE2	1:A:41:ARG:CG	2.19	0.71
2:B:145:LEU:CD1	2:B:151:ARG:NH1	2.51	0.71
2:B:172:VAL:HG13	2:B:175:ARG:HG2	1.71	0.71
1:A:187:HIS:CE1	1:A:278:MET:HE1	2.25	0.71
2:B:172:VAL:CG1	2:B:175:ARG:CG	2.69	0.70
2:B:172:VAL:HG12	2:B:175:ARG:HG2	1.71	0.70
1:A:373:LYS:O	1:A:386:ILE:CG2	2.40	0.69
2:B:142:LYS:O	2:B:146:GLU:CG	2.29	0.68
2:B:115:TYR:CD1	2:B:145:LEU:CD2	2.70	0.68
2:B:172:VAL:HG12	2:B:175:ARG:CG	2.26	0.66
1:A:97:THR:O	1:A:101:CYS:SG	2.55	0.64
2:B:170:LYS:CG	7:B:451:HOH:O	2.40	0.64
1:A:445:ASP:OD1	4:A:503:FAD:H5'1	1.98	0.64
1:A:81:ARG:NH1	1:A:85:ARG:HH12	1.97	0.62
2:B:159:LEU:HG	2:B:161:PRO:HD2	1.82	0.62
2:B:155:GLU:HG2	2:B:155:GLU:O	2.00	0.62
1:A:304:THR:HA	4:A:503:FAD:HM73	1.83	0.61
2:B:142:LYS:HG2	2:B:146:GLU:OE2	2.00	0.60
1:A:192:VAL:HG11	4:A:503:FAD:C9	2.31	0.60
1:A:184:GLU:CD	4:A:503:FAD:HO2A	2.09	0.59
2:B:142:LYS:HG3	2:B:153:ILE:CD1	2.31	0.58
1:A:409:ASN:HB3	1:A:412:LEU:HB2	1.85	0.58
1:A:81:ARG:NH1	1:A:85:ARG:NH1	2.52	0.58
1:A:386:ILE:CG2	1:A:387:VAL:N	2.38	0.57
2:B:145:LEU:CB	2:B:151:ARG:NH1	2.68	0.57
2:B:127:ASP:OD1	2:B:149:VAL:CG2	2.49	0.57
2:B:159:LEU:HB3	2:B:162:ASN:HB2	1.88	0.56
4:B:301:FAD:O3B	7:B:402:HOH:O	2.17	0.56
1:A:249:ALA:O	1:A:406:GLN:HB3	2.06	0.56
1:A:214:LEU:HB3	1:A:219:VAL:HG22	1.89	0.55
1:A:325:ARG:HG2	1:A:386:ILE:HD12	1.87	0.55
2:B:156:PRO:HG2	2:B:173:SER:HB2	1.87	0.55
2:B:115:TYR:HD1	2:B:145:LEU:HD21	1.71	0.54
1:A:147:ILE:HA	1:A:150:ILE:HD12	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:170:LYS:C	2:B:172:VAL:N	2.61	0.54
1:A:258:PRO:HB2	1:A:355:GLU:HG3	1.90	0.53
2:B:128:VAL:O	2:B:149:VAL:HB	2.08	0.53
2:B:144:GLU:C	2:B:146:GLU:H	2.17	0.53
2:B:136:GLU:HG2	2:B:137:PRO:HD3	1.91	0.52
1:A:6:LYS:O	1:A:6:LYS:HG3	2.09	0.52
2:B:143:GLU:O	2:B:146:GLU:HB2	2.10	0.51
1:A:40:GLU:HA	1:A:70:ARG:HD2	1.91	0.51
1:A:248:GLY:HA2	1:A:445:ASP:HB2	1.93	0.50
2:B:164:ASP:O	2:B:171:ASN:HB2	2.11	0.50
2:B:145:LEU:HB3	2:B:151:ARG:NH1	2.26	0.49
1:A:411:VAL:HG12	1:A:411:VAL:O	2.12	0.49
2:B:73:MET:O	2:B:77:LEU:HB2	2.11	0.49
2:B:131:LEU:CD2	2:B:181:ARG:HH12	2.27	0.48
1:A:325:ARG:HH22	1:A:385:LYS:HD3	1.78	0.48
2:B:236:THR:HG22	7:B:426:HOH:O	2.14	0.48
1:A:384:ARG:O	1:A:385:LYS:HB2	2.12	0.48
1:A:85:ARG:O	1:A:89:ARG:HG3	2.13	0.48
2:B:115:TYR:HB2	2:B:145:LEU:HD21	1.95	0.47
1:A:69:LEU:HA	1:A:79:ALA:HB1	1.96	0.47
1:A:306:MET:HE1	1:A:347:PHE:HE2	1.81	0.46
1:A:184:GLU:OE2	4:A:503:FAD:H1B	2.15	0.46
1:A:160:ILE:HD12	1:A:245:ILE:HG12	1.97	0.45
1:A:434:MET:HE2	1:A:441:PHE:CZ	2.51	0.45
1:A:445:ASP:CG	4:A:503:FAD:H5'1	2.42	0.45
2:B:144:GLU:C	2:B:146:GLU:N	2.75	0.45
1:A:195:TYR:CZ	2:B:229:MET:CE	3.00	0.45
1:A:279:LYS:HD2	1:A:286:TYR:CZ	2.51	0.44
2:B:184:LEU:O	2:B:185:GLU:HB3	2.16	0.44
2:B:115:TYR:CD1	2:B:145:LEU:HD21	2.49	0.44
1:A:201:ARG:HD3	1:A:307:ASP:OD1	2.19	0.43
1:A:191:GLY:CA	4:A:503:FAD:O3B	2.50	0.43
2:B:127:ASP:OD1	2:B:149:VAL:HG11	2.17	0.43
2:B:172:VAL:HG12	2:B:175:ARG:HG3	2.01	0.43
1:A:279:LYS:HB3	1:A:282:LYS:HD2	2.01	0.43
1:A:371:LYS:HE3	1:A:393:ILE:HB	2.00	0.42
1:A:192:VAL:HG13	1:A:196:GLY:HA3	2.02	0.42
1:A:192:VAL:HG11	4:A:503:FAD:C8	2.49	0.42
2:B:45:ILE:HB	4:B:301:FAD:O4'	2.19	0.42
2:B:145:LEU:HB3	2:B:151:ARG:HH11	1.84	0.42
2:B:170:LYS:C	2:B:172:VAL:H	2.22	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:185:GLU:O	2:B:185:GLU:HG2	2.20	0.41
2:B:68:LYS:HD3	2:B:279:TYR:HB2	2.02	0.41
1:A:133:ALA:HA	1:A:138:ILE:HB	2.03	0.41
1:A:158:ALA:HB2	1:A:240:TYR:CD1	2.56	0.41
1:A:195:TYR:CZ	2:B:229:MET:HE1	2.54	0.41
1:A:447:ILE:HG13	1:A:448:ARG:HG3	2.02	0.41
1:A:73:ARG:H	1:A:73:ARG:HG2	1.76	0.40
1:A:261:ASN:HD22	2:B:260:GLU:CD	2.29	0.40
2:B:62:ILE:HB	2:B:112:VAL:HG11	2.03	0.40
2:B:48:THR:HA	2:B:49:PRO:HD3	2.01	0.40
1:A:87:ILE:HG21	1:A:94:PRO:HB3	2.04	0.40
2:B:6:GLU:HB2	2:B:18:LYS:HB3	2.04	0.40
2:B:50:VAL:HG21	2:B:112:VAL:HG13	2.02	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	458/476 (96%)	434 (95%)	24 (5%)	0	100	100
2	B	278/289 (96%)	257 (92%)	21 (8%)	0	100	100
All	All	736/765 (96%)	691 (94%)	45 (6%)	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	375/388 (97%)	372 (99%)	3 (1%)	73	84
2	B	244/249 (98%)	237 (97%)	7 (3%)	37	55
All	All	619/637 (97%)	609 (98%)	10 (2%)	55	73

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ILE
1	A	54	LYS
1	A	219	VAL
2	B	36	VAL
2	B	94	VAL
2	B	105	ILE
2	B	124	ILE
2	B	136	GLU
2	B	183	LEU
2	B	209	LYS

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

Mol	Chain	Res	Type
1	A	61	ASN
1	A	121	ASN
1	A	188	GLN
1	A	432	ASN
2	B	203	GLN
2	B	255	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 for Mogul analysis.

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	$\# Z > 2$
3	SF4	A	501	1	0,12,12	-	-	-		
4	FAD	A	503	-	58,58,58	1.47	8 (13%)	85,89,89	1.61	19 (22%)
4	FAD	B	301	6	58,58,58	1.47	8 (13%)	85,89,89	1.61	18 (21%)
3	SF4	A	502	1	0,12,12	-	-	-		
5	FES	B	302	2	0,4,4	-	-	-		

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
3	SF4	A	501	1	-	-	0/6/5/5
4	FAD	A	503	-	-	13/34/50/50	0/6/6/6
4	FAD	B	301	6	-	4/34/50/50	0/6/6/6
3	SF4	A	502	1	-	-	0/6/5/5
5	FES	B	302	2	-	-	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	301	FAD	C9A-C5X	5.29	1.49	1.41
4	A	503	FAD	C9A-C5X	5.24	1.49	1.41
4	A	503	FAD	C5A-C4A	4.63	1.47	1.39
4	B	301	FAD	C5A-C4A	4.60	1.47	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	503	FAD	C8-C7	3.44	1.49	1.40
4	B	301	FAD	C8-C7	3.40	1.49	1.40
4	A	503	FAD	C5A-C6A	2.75	1.48	1.41
4	B	301	FAD	C5A-C6A	2.70	1.48	1.41
4	B	301	FAD	C4-N3	-2.42	1.34	1.38
4	B	301	FAD	C8A-N7A	2.42	1.36	1.31
4	A	503	FAD	C8A-N7A	2.39	1.36	1.31
4	A	503	FAD	C4-N3	-2.38	1.34	1.38
4	A	503	FAD	C5A-N7A	-2.28	1.34	1.39
4	B	301	FAD	C5A-N7A	-2.23	1.35	1.39
4	B	301	FAD	C4X-N5	2.23	1.35	1.30
4	A	503	FAD	C4X-N5	2.20	1.35	1.30

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	301	FAD	C5A-C4A-N3A	-5.83	118.70	126.72
4	A	503	FAD	C5A-C4A-N3A	-5.81	118.72	126.72
4	A	503	FAD	N3A-C4A-N9A	4.65	135.08	127.17
4	B	301	FAD	N3A-C4A-N9A	4.64	135.05	127.17
4	A	503	FAD	C2A-N3A-C4A	3.89	121.34	111.83
4	B	301	FAD	C2A-N3A-C4A	3.89	121.32	111.83
4	B	301	FAD	N3A-C2A-N1A	-3.73	122.94	128.58
4	A	503	FAD	N3A-C2A-N1A	-3.73	122.94	128.58
4	A	503	FAD	C4A-C5A-N7A	-3.53	106.55	110.58
4	B	301	FAD	C4A-C5A-N7A	-3.51	106.56	110.58
4	A	503	FAD	C4A-N9A-C8A	2.94	108.83	105.74
4	B	301	FAD	C4A-N9A-C8A	2.89	108.78	105.74
4	A	503	FAD	C4-C4X-N5	2.83	122.11	118.21
4	B	301	FAD	C4-C4X-N5	2.82	122.10	118.21
4	A	503	FAD	C5A-N7A-C8A	2.78	107.82	103.45
4	B	301	FAD	C5A-N7A-C8A	2.74	107.75	103.45
4	B	301	FAD	C3B-C2B-C1B	2.56	106.31	101.46
4	A	503	FAD	C3B-C2B-C1B	2.53	106.26	101.46
4	A	503	FAD	C4X-C10-N1	-2.48	118.51	124.59
4	B	301	FAD	C4X-C10-N1	-2.48	118.51	124.59
4	A	503	FAD	N9A-C8A-N7A	-2.35	110.60	113.94
4	B	301	FAD	N9A-C8A-N7A	-2.31	110.66	113.94
4	B	301	FAD	O4-C4-C4X	-2.23	120.64	126.53
4	A	503	FAD	O4-C4-C4X	-2.22	120.68	126.53
4	A	503	FAD	C6A-C5A-N7A	2.19	136.31	132.09
4	B	301	FAD	C6A-C5A-N7A	2.15	136.24	132.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	301	FAD	C4X-C4-N3	2.13	118.68	113.25
4	A	503	FAD	C4X-C4-N3	2.13	118.66	113.25
4	B	301	FAD	C10-N1-C2	2.12	121.44	116.85
4	B	301	FAD	C4-N3-C2	-2.12	121.89	125.64
4	A	503	FAD	C4-N3-C2	-2.11	121.89	125.64
4	A	503	FAD	C4X-C10-N10	2.11	119.50	116.48
4	B	301	FAD	C4X-C10-N10	2.11	119.50	116.48
4	A	503	FAD	C2A-N1A-C6A	2.10	122.18	118.73
4	A	503	FAD	C10-N1-C2	2.09	121.37	116.85
4	B	301	FAD	C2A-N1A-C6A	2.08	122.14	118.73
4	A	503	FAD	O2-C2-N1	-2.01	118.47	121.80

There are no chirality outliers.

All (17) torsion outliers are listed below:

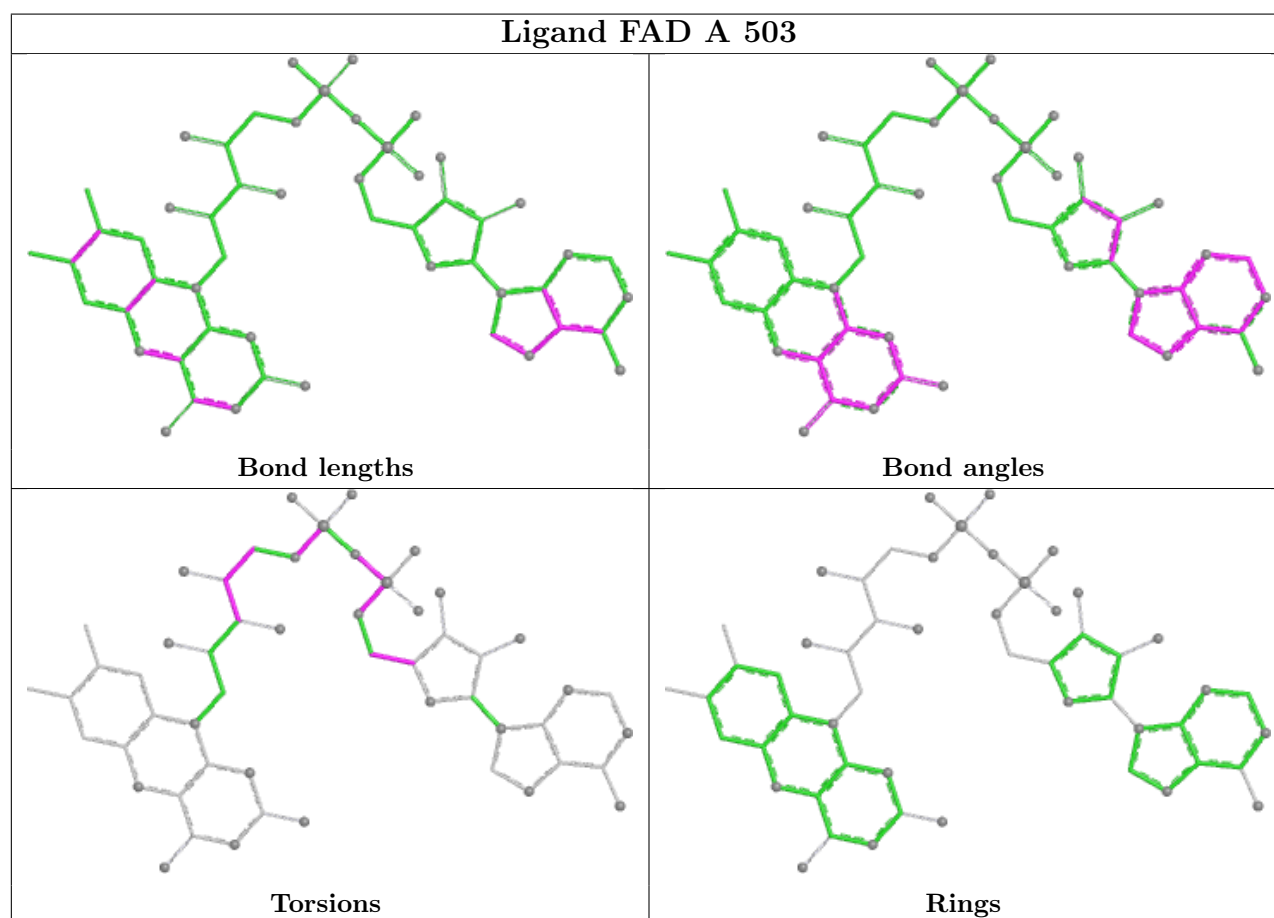
Mol	Chain	Res	Type	Atoms
4	A	503	FAD	C5B-O5B-PA-O2A
4	A	503	FAD	C5'-O5'-P-O3P
4	B	301	FAD	C5B-O5B-PA-O3P
4	B	301	FAD	C5'-O5'-P-O1P
4	B	301	FAD	C5'-O5'-P-O2P
4	A	503	FAD	O4B-C4B-C5B-O5B
4	A	503	FAD	C3B-C4B-C5B-O5B
4	A	503	FAD	C3'-C4'-C5'-O5'
4	A	503	FAD	C5B-O5B-PA-O1A
4	A	503	FAD	C5B-O5B-PA-O3P
4	A	503	FAD	C5'-O5'-P-O1P
4	B	301	FAD	C5'-O5'-P-O3P
4	A	503	FAD	C2'-C3'-C4'-O4'
4	A	503	FAD	P-O3P-PA-O1A
4	A	503	FAD	P-O3P-PA-O2A
4	A	503	FAD	O3'-C3'-C4'-C5'
4	A	503	FAD	O3'-C3'-C4'-O4'

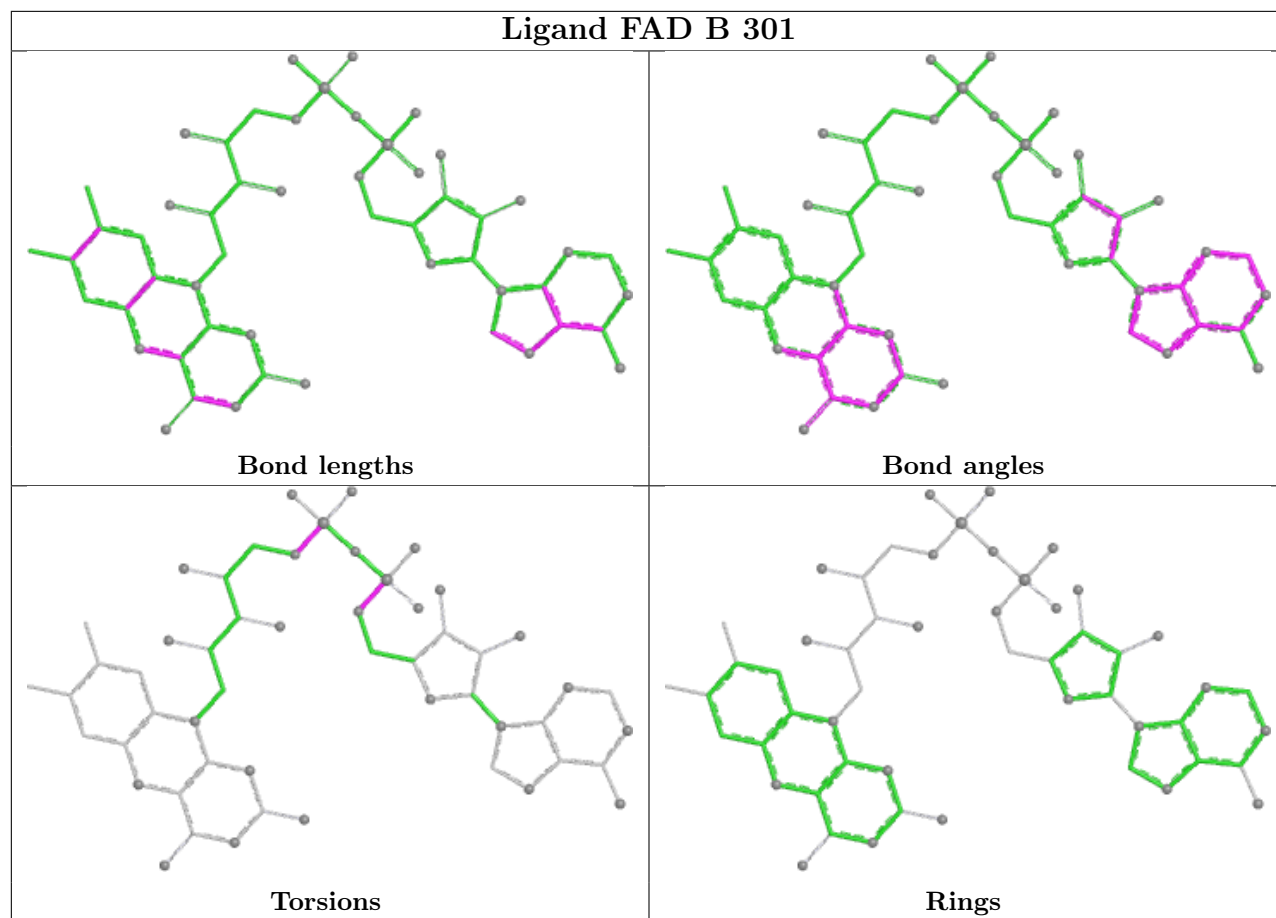
There are no ring outliers.

3 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	503	FAD	12	0
4	B	301	FAD	2	0
5	B	302	FES	1	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 [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	462/476 (97%)	0.25	14 (3%) 52 54	26, 41, 66, 114	0
2	B	282/289 (97%)	0.38	28 (9%) 13 12	27, 39, 66, 93	0
All	All	744/765 (97%)	0.30	42 (5%) 30 29	26, 40, 66, 114	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	148	ALA	4.8
2	B	181	ARG	4.6
2	B	151	ARG	4.5
2	B	152	HIS	4.5
2	B	170	LYS	4.5
2	B	165	PHE	4.3
2	B	147	LYS	3.9
2	B	145	LEU	3.8
2	B	171	ASN	3.7
2	B	172	VAL	3.7
2	B	163	GLN	3.6
2	B	150	THR	3.6
1	A	413	TRP	3.6
2	B	167	ALA	3.5
2	B	283	LEU	3.2
2	B	177	LYS	3.1
2	B	168	ASN	3.1
1	A	382	GLY	3.0
2	B	173	SER	3.0
2	B	166	LEU	3.0
1	A	386	ILE	2.9
1	A	383	LYS	2.9
2	B	146	GLU	2.8
1	A	5	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
2	B	149	VAL	2.7
2	B	155	GLU	2.7
1	A	410	LYS	2.6
1	A	101	CYS	2.6
2	B	182	GLU	2.5
2	B	154	VAL	2.5
1	A	116	VAL	2.5
1	A	384	ARG	2.4
2	B	265	VAL	2.3
1	A	385	LYS	2.2
2	B	185	GLU	2.2
1	A	411	VAL	2.2
2	B	164	ASP	2.2
1	A	415	ALA	2.1
2	B	176	LEU	2.1
1	A	417	PRO	2.1
2	B	23	HIS	2.1
1	A	115	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 oligosaccharides 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	FAD	A	503	53/53	0.81	0.14	26,35,46,53	0
4	FAD	B	301	53/53	0.87	0.11	28,37,41,44	0
6	MG	B	303	1/1	0.87	0.22	34,34,34,34	0
5	FES	B	302	4/4	0.94	0.07	35,35,36,39	0
3	SF4	A	501	8/8	0.96	0.06	37,39,40,40	0

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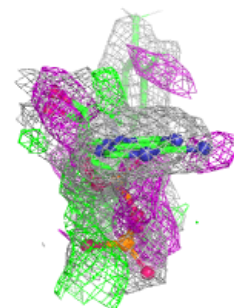
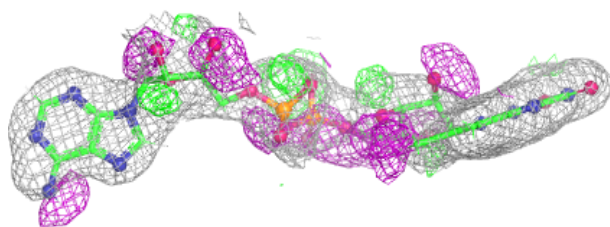
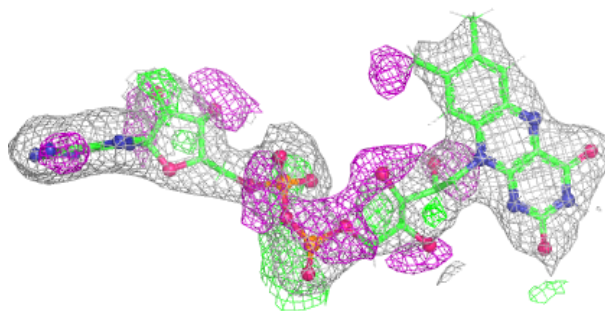
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SF4	A	502	8/8	0.96	0.05	40,45,46,53	0

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.

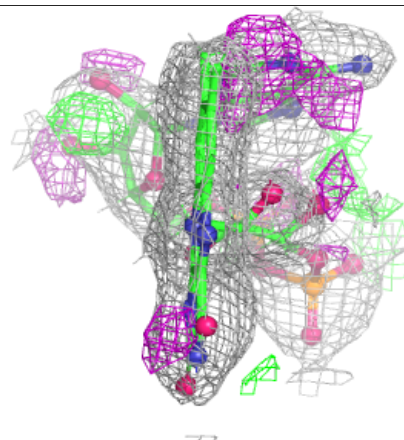
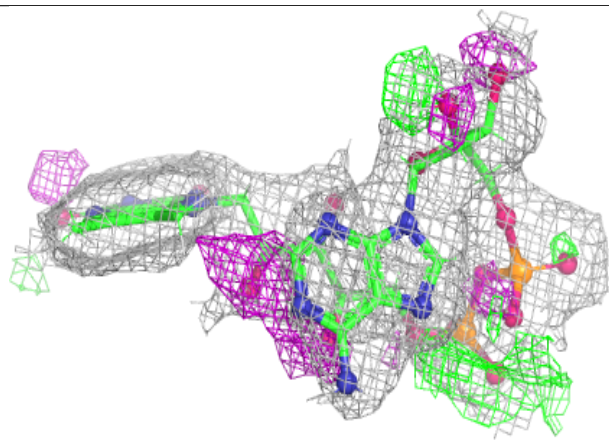
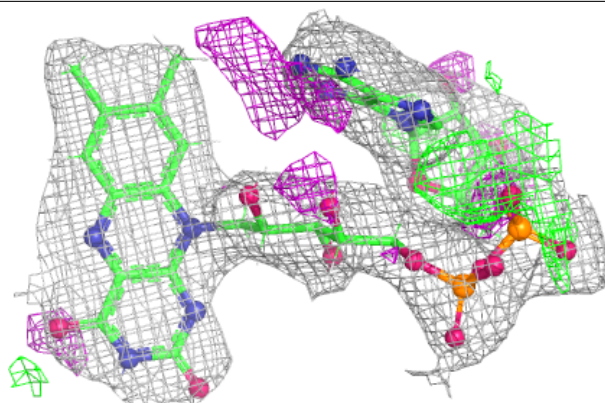
Electron density around FAD A 503:

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 301:

$2mF_o - DF_c$ (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.