



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:31 AM UTC

PDB ID : 4P8N / pdb_00004p8n
Title : Crystal structure of M. tuberculosis DprE1 in complex with the non-covalent inhibitor QN118
Authors : Neres, J.; Pojer, F.; Cole, S.T.
Deposited on : 2014-03-31
Resolution : 1.79 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

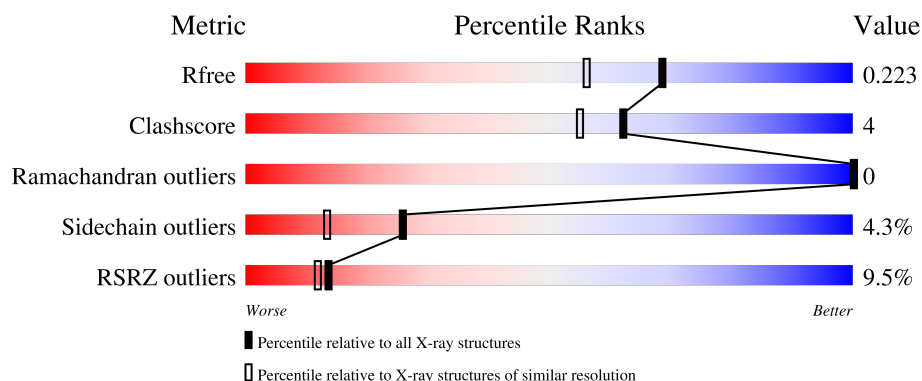
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.79 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	480	
1	B	480	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7246 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 Probable decaprenylphosphoryl-beta-D-ribose oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	454	Total	C	N	O	S	0	1	0
			3502	2230	616	645	11			
1	B	419	Total	C	N	O	S	0	2	0
			3230	2049	574	596	11			

There are 38 discrepancies between the modelled and reference sequences:

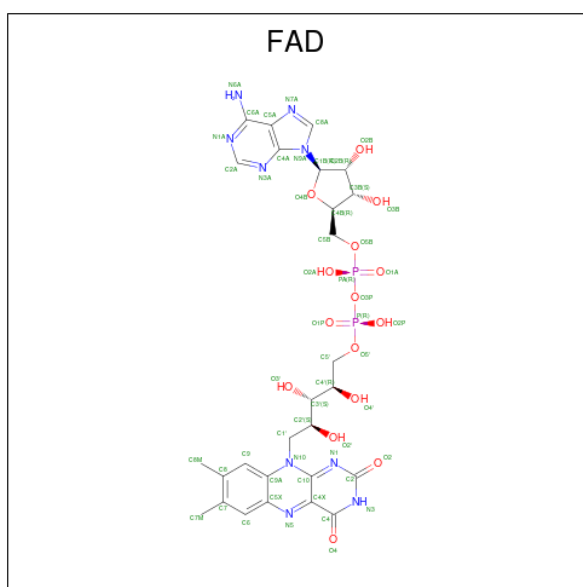
Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	GLY	-	expression tag	UNP P72056
A	-17	SER	-	expression tag	UNP P72056
A	-16	SER	-	expression tag	UNP P72056
A	-15	HIS	-	expression tag	UNP P72056
A	-14	HIS	-	expression tag	UNP P72056
A	-13	HIS	-	expression tag	UNP P72056
A	-12	HIS	-	expression tag	UNP P72056
A	-11	HIS	-	expression tag	UNP P72056
A	-10	HIS	-	expression tag	UNP P72056
A	-9	SER	-	expression tag	UNP P72056
A	-8	SER	-	expression tag	UNP P72056
A	-7	GLY	-	expression tag	UNP P72056
A	-6	LEU	-	expression tag	UNP P72056
A	-5	VAL	-	expression tag	UNP P72056
A	-4	PRO	-	expression tag	UNP P72056
A	-3	ARG	-	expression tag	UNP P72056
A	-2	GLY	-	expression tag	UNP P72056
A	-1	SER	-	expression tag	UNP P72056
A	0	HIS	-	expression tag	UNP P72056
B	-18	GLY	-	expression tag	UNP P72056
B	-17	SER	-	expression tag	UNP P72056
B	-16	SER	-	expression tag	UNP P72056
B	-15	HIS	-	expression tag	UNP P72056
B	-14	HIS	-	expression tag	UNP P72056
B	-13	HIS	-	expression tag	UNP P72056

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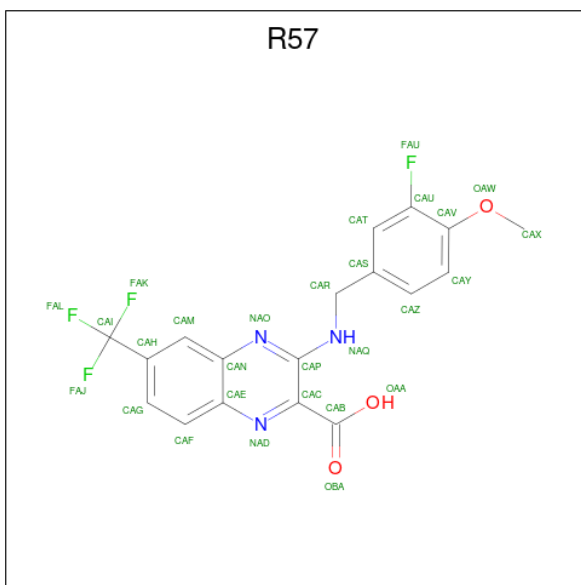
Chain	Residue	Modelled	Actual	Comment	Reference
B	-12	HIS	-	expression tag	UNP P72056
B	-11	HIS	-	expression tag	UNP P72056
B	-10	HIS	-	expression tag	UNP P72056
B	-9	SER	-	expression tag	UNP P72056
B	-8	SER	-	expression tag	UNP P72056
B	-7	GLY	-	expression tag	UNP P72056
B	-6	LEU	-	expression tag	UNP P72056
B	-5	VAL	-	expression tag	UNP P72056
B	-4	PRO	-	expression tag	UNP P72056
B	-3	ARG	-	expression tag	UNP P72056
B	-2	GLY	-	expression tag	UNP P72056
B	-1	SER	-	expression tag	UNP P72056
B	0	HIS	-	expression tag	UNP P72056

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



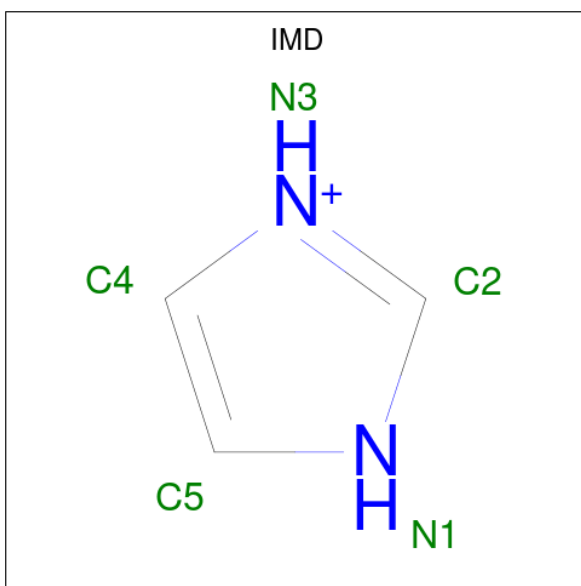
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is 3-[(3-fluoro-4-methoxybenzyl)amino]-6-(trifluoromethyl)quinoxaline-2-carboxylic acid (CCD ID: R57) (formula: $C_{18}H_{13}F_4N_3O_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	1
			56	36	8	6	6		
3	B	1	Total	C	F	N	O	0	0
			28	18	4	3	3		

- Molecule 4 is IMIDAZOLE (CCD ID: IMD) (formula: $C_3H_5N_2$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	C N	0	0
			5	3 2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	198	Total 198	O 198	0	0
5	B	121	Total 121	O 121	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	77.24Å 83.22Å 80.75Å 90.00° 103.04° 90.00°	Depositor
Resolution (Å)	41.60 – 1.79 41.60 – 1.79	Depositor EDS
% Data completeness (in resolution range)	98.3 (41.60-1.79) 92.4 (41.60-1.79)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.98 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.203 , 0.230 (Not available) , 0.223	Depositor DCC
R_{free} test set	4600 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	28.0	Xtriage
Anisotropy	0.018	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 28.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.021 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7246	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: IMD, R57, 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	0.66	1/3590 (0.0%)	0.78	2/4878 (0.0%)
1	B	0.64	1/3309 (0.0%)	0.78	0/4488
All	All	0.65	2/6899 (0.0%)	0.78	2/9366 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	45	SER	CB-OG	6.14	1.54	1.42
1	A	326	ALA	C-O	5.30	1.34	1.23

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	290	GLY	CA-C-N	5.18	124.98	119.28
1	A	290	GLY	C-N-CA	5.18	124.98	119.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3502	0	3471	28	0
1	B	3230	0	3216	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	53	0	31	0	0
2	B	53	0	31	0	0
3	A	56	0	24	7	0
3	B	28	0	12	0	0
4	B	5	0	5	0	0
5	A	198	0	0	1	0
5	B	121	0	0	2	0
All	All	7246	0	6790	56	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 4.

All (56) 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:159:ARG:HG2	1:B:159:ARG:HH11	1.08	1.12
1:B:159:ARG:HG2	1:B:159:ARG:NH1	1.86	0.89
1:A:230[B]:TRP:CD1	3:A:502[B]:R57:CAX	2.57	0.87
1:A:442:ARG:HH11	1:A:442:ARG:HG2	1.49	0.78
1:A:230[B]:TRP:CD1	3:A:502[B]:R57:OAW	2.41	0.73
1:A:230[B]:TRP:CD1	3:A:502[B]:R57:H1	2.25	0.71
1:B:372[B]:ARG:HD3	1:B:381:ILE:HG22	1.73	0.70
1:A:230[B]:TRP:CG	3:A:502[B]:R57:H1	2.29	0.68
1:A:334:GLN:HE22	1:A:387:CYS:HB3	1.58	0.68
1:A:442:ARG:HH11	1:A:442:ARG:CG	2.10	0.64
1:B:99:ASP:O	1:B:103:LYS:HG3	1.97	0.63
1:A:99:ASP:HB2	1:A:119:ARG:HB3	1.81	0.62
1:A:366:PHE:CD2	1:A:386:ILE:HD11	2.34	0.62
1:A:230[B]:TRP:NE1	3:A:502[B]:R57:OAW	2.33	0.61
1:B:240:LEU:HD22	1:B:355:GLN:HG2	1.87	0.57
1:A:347:PHE:O	1:A:350:ILE:HG22	2.05	0.56
3:A:502[B]:R57:FAU	3:A:502[B]:R57:H2	1.95	0.56
1:A:347:PHE:CE1	1:A:386:ILE:HD13	2.43	0.54
1:A:366:PHE:HD2	1:A:386:ILE:HD11	1.74	0.53
1:A:347:PHE:CZ	1:A:386:ILE:HD13	2.43	0.53
1:A:375:ALA:HB3	1:A:378:SER:HB2	1.91	0.52
1:B:33:GLU:O	1:B:37:LYS:HG2	2.11	0.50
1:B:347:PHE:O	1:B:350:ILE:HG22	2.11	0.50
1:B:374:GLN:HG2	5:B:689:HOH:O	2.12	0.48
1:A:164:THR:HG22	5:A:756:HOH:O	2.13	0.48
1:B:375:ALA:HB3	1:B:378:SER:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:LEU:O	1:A:299:LYS:HG3	2.15	0.47
1:B:230:TRP:HA	1:B:364:ASN:O	2.14	0.47
1:B:153:THR:HG21	1:B:159:ARG:HD3	1.96	0.47
1:A:200:ILE:HD11	1:A:255:GLN:OE1	2.16	0.46
1:A:231:PHE:CE1	1:A:364:ASN:HB3	2.50	0.46
3:A:502[B]:R57:CAX	3:A:502[B]:R57:FAU	2.54	0.46
1:B:116:PRO:HG2	1:B:118:THR:O	2.16	0.45
1:B:83:ILE:HG12	1:B:90:VAL:HG12	1.99	0.45
1:A:61:GLY:HA2	1:A:420:SER:HB3	1.98	0.45
1:B:334:GLN:HG2	1:B:389:ASP:OD1	2.16	0.45
1:B:72:ILE:HD12	1:B:183:ILE:HD11	1.98	0.44
1:B:231:PHE:CE1	1:B:364:ASN:HB3	2.53	0.43
1:A:116:PRO:HG2	1:A:118:THR:O	2.18	0.43
1:A:83:ILE:HG12	1:A:90:VAL:HG12	1.99	0.43
1:B:414:LEU:HG	1:B:430:MET:HE1	2.00	0.43
1:A:335:TYR:CZ	1:A:414:LEU:HD21	2.54	0.43
1:A:115:LEU:O	1:A:314:TYR:OH	2.33	0.43
1:A:230[B]:TRP:HZ2	1:A:316:PRO:HG2	1.84	0.43
1:B:455:ALA:HB1	1:B:461:LEU:HD13	2.02	0.42
1:B:60:TYR:CD2	1:B:418:LYS:HD3	2.54	0.42
1:B:100:GLN:HA	1:B:103:LYS:HE2	2.00	0.42
1:B:87:THR:O	1:B:88:LYS:HB2	2.20	0.41
1:A:92:ILE:HB	1:A:96:VAL:HG21	2.02	0.41
1:B:103:LYS:HG2	1:B:267:PHE:CE1	2.56	0.41
1:B:159:ARG:NH1	1:B:159:ARG:CG	2.66	0.41
1:B:90:VAL:O	1:B:188:THR:HA	2.21	0.41
1:A:75:THR:N	1:A:76:PRO:CD	2.84	0.41
1:A:442:ARG:CG	1:A:442:ARG:NH1	2.74	0.40
1:B:34:MET:HA	1:B:34:MET:HE2	2.04	0.40
1:B:88:LYS:HE3	5:B:648:HOH:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	451/480 (94%)	448 (99%)	3 (1%)	0	100	100
1	B	415/480 (86%)	410 (99%)	5 (1%)	0	100	100
All	All	866/960 (90%)	858 (99%)	8 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	364/384 (95%)	347 (95%)	17 (5%)	23	11
1	B	337/384 (88%)	324 (96%)	13 (4%)	28	16
All	All	701/768 (91%)	671 (96%)	30 (4%)	26	13

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	THR
1	A	77	LEU
1	A	88	LYS
1	A	92	ILE
1	A	121	VAL
1	A	164	THR
1	A	167	ASP
1	A	256	LEU
1	A	259	LYS
1	A	268	ASP
1	A	272	LEU
1	A	325	ARG
1	A	365	VAL
1	A	377	LEU
1	A	386	ILE

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Mol	Chain	Res	Type
1	A	421	ARG
1	A	442	ARG
1	B	44	GLU
1	B	92	ILE
1	B	159	ARG
1	B	253	VAL
1	B	260	LEU
1	B	334	GLN
1	B	355	GLN
1	B	396	LEU
1	B	421	ARG
1	B	422	THR
1	B	444	VAL
1	B	448	ARG
1	B	461	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	324	ASN
1	A	334	GLN
1	B	308	GLN
1	B	315	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry

6 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	-	58,58,58	1.50	9 (15%)	85,89,89	1.54	15 (17%)
3	R57	B	502	-	30,30,30	1.09	2 (6%)	42,44,44	1.57	4 (9%)
3	R57	A	502[B]	-	30,30,30	1.05	2 (6%)	42,44,44	1.55	3 (7%)
4	IMD	B	503	-	5,5,5	0.56	0	5,5,5	0.44	0
2	FAD	B	501	-	58,58,58	1.54	10 (17%)	85,89,89	1.58	19 (22%)
3	R57	A	502[A]	-	30,30,30	1.08	1 (3%)	42,44,44	1.50	4 (9%)

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	-	-	2/34/50/50	0/6/6/6
3	R57	B	502	-	-	2/17/17/17	0/3/3/3
3	R57	A	502[B]	-	-	2/17/17/17	0/3/3/3
4	IMD	B	503	-	-	-	0/1/1/1
2	FAD	B	501	-	-	2/34/50/50	0/6/6/6
3	R57	A	502[A]	-	-	0/17/17/17	0/3/3/3

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	FAD	C9A-C5X	5.42	1.49	1.41
2	A	501	FAD	C9A-C5X	5.37	1.49	1.41
2	B	501	FAD	C5A-C4A	4.52	1.47	1.39
2	A	501	FAD	C5A-C4A	4.48	1.47	1.39
2	B	501	FAD	C8-C7	3.66	1.49	1.40
3	A	502[A]	R57	CAP-NAO	3.59	1.36	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	502	R57	CAP-NAO	3.51	1.36	1.31
2	A	501	FAD	C8-C7	3.46	1.49	1.40
2	A	501	FAD	PA-O3P	3.31	1.63	1.59
3	A	502[B]	R57	CAP-NAO	3.13	1.35	1.31
2	B	501	FAD	PA-O3P	3.08	1.62	1.59
2	B	501	FAD	C5A-C6A	2.72	1.48	1.41
3	B	502	R57	CAC-NAD	2.50	1.36	1.32
2	B	501	FAD	C4X-N5	2.40	1.36	1.30
2	B	501	FAD	C5A-N7A	-2.39	1.34	1.39
2	A	501	FAD	C5A-C6A	2.37	1.47	1.41
2	A	501	FAD	C4-N3	-2.35	1.34	1.38
2	A	501	FAD	C4X-N5	2.33	1.35	1.30
2	B	501	FAD	C8A-N7A	2.29	1.36	1.31
2	B	501	FAD	C4A-N9A	-2.19	1.33	1.37
2	B	501	FAD	P-O3P	2.19	1.61	1.59
2	A	501	FAD	P-O3P	2.11	1.61	1.59
2	A	501	FAD	C5A-N7A	-2.10	1.35	1.39
3	A	502[B]	R57	CAC-NAD	2.07	1.35	1.32

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	R57	OAW-CAV-CAU	6.18	120.71	116.27
3	A	502[B]	R57	OAW-CAV-CAU	5.85	120.47	116.27
2	B	501	FAD	C5A-C4A-N3A	-5.03	119.79	126.72
3	A	502[A]	R57	OAW-CAV-CAU	4.80	119.72	116.27
2	A	501	FAD	C5A-C4A-N3A	-4.63	120.34	126.72
2	A	501	FAD	N3A-C4A-N9A	4.42	134.68	127.17
2	B	501	FAD	N3A-C4A-N9A	4.16	134.24	127.17
2	B	501	FAD	N3A-C2A-N1A	-4.05	122.45	128.58
2	A	501	FAD	C4A-N9A-C8A	4.01	109.94	105.74
2	A	501	FAD	N3A-C2A-N1A	-3.90	122.68	128.58
2	B	501	FAD	C2A-N3A-C4A	3.79	121.08	111.83
2	A	501	FAD	C2A-N3A-C4A	3.41	120.16	111.83
2	B	501	FAD	C4A-C5A-N7A	-3.24	106.88	110.58
2	B	501	FAD	C4A-N9A-C8A	3.20	109.10	105.74
2	B	501	FAD	C4-C4X-N5	3.19	122.62	118.21
2	A	501	FAD	C4-C4X-N5	3.08	122.47	118.21
3	B	502	R57	OAW-CAV-CAY	-2.97	119.30	124.30
3	A	502[B]	R57	OAW-CAV-CAY	-2.89	119.42	124.30
2	A	501	FAD	C4A-C5A-N7A	-2.84	107.34	110.58
2	B	501	FAD	O4-C4-C4X	-2.82	119.08	126.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	FAD	O4-C4-C4X	-2.69	119.43	126.53
2	B	501	FAD	C6A-C5A-N7A	2.57	137.05	132.09
2	A	501	FAD	N9A-C8A-N7A	-2.52	110.36	113.94
2	B	501	FAD	C5A-N7A-C8A	2.50	107.37	103.45
2	A	501	FAD	C5A-N7A-C8A	2.50	107.37	103.45
2	B	501	FAD	O2-C2-N1	-2.43	117.76	121.80
3	A	502[A]	R57	FAL-CAI-CAH	-2.39	107.77	112.90
3	A	502[A]	R57	OAW-CAV-CAY	-2.37	120.31	124.30
2	A	501	FAD	C2A-N1A-C6A	2.33	122.56	118.73
2	B	501	FAD	C2A-N1A-C6A	2.30	122.51	118.73
2	B	501	FAD	N9A-C8A-N7A	-2.30	110.67	113.94
2	A	501	FAD	C4X-C4-N3	2.25	118.97	113.25
2	B	501	FAD	C10-N1-C2	2.25	121.71	116.85
2	B	501	FAD	C4X-C4-N3	2.24	118.96	113.25
2	B	501	FAD	C4X-C10-N1	-2.24	119.10	124.59
2	B	501	FAD	C4X-C10-N10	2.21	119.64	116.48
2	A	501	FAD	C6A-C5A-N7A	2.18	136.29	132.09
3	A	502[A]	R57	CAC-CAP-NAQ	-2.12	117.86	121.94
2	B	501	FAD	C4-N3-C2	-2.11	121.90	125.64
2	A	501	FAD	C10-N1-C2	2.10	121.40	116.85
2	B	501	FAD	O2P-P-O3P	2.10	112.96	107.27
3	B	502	R57	CAP-NAO-CAN	2.06	120.52	116.66
2	A	501	FAD	C9A-C5X-N5	-2.06	120.27	122.45
3	A	502[B]	R57	CAT-CAU-CAV	-2.04	120.10	122.86
3	B	502	R57	CAT-CAU-CAV	-2.00	120.16	122.86

There are no chirality outliers.

All (8) torsion outliers are listed below:

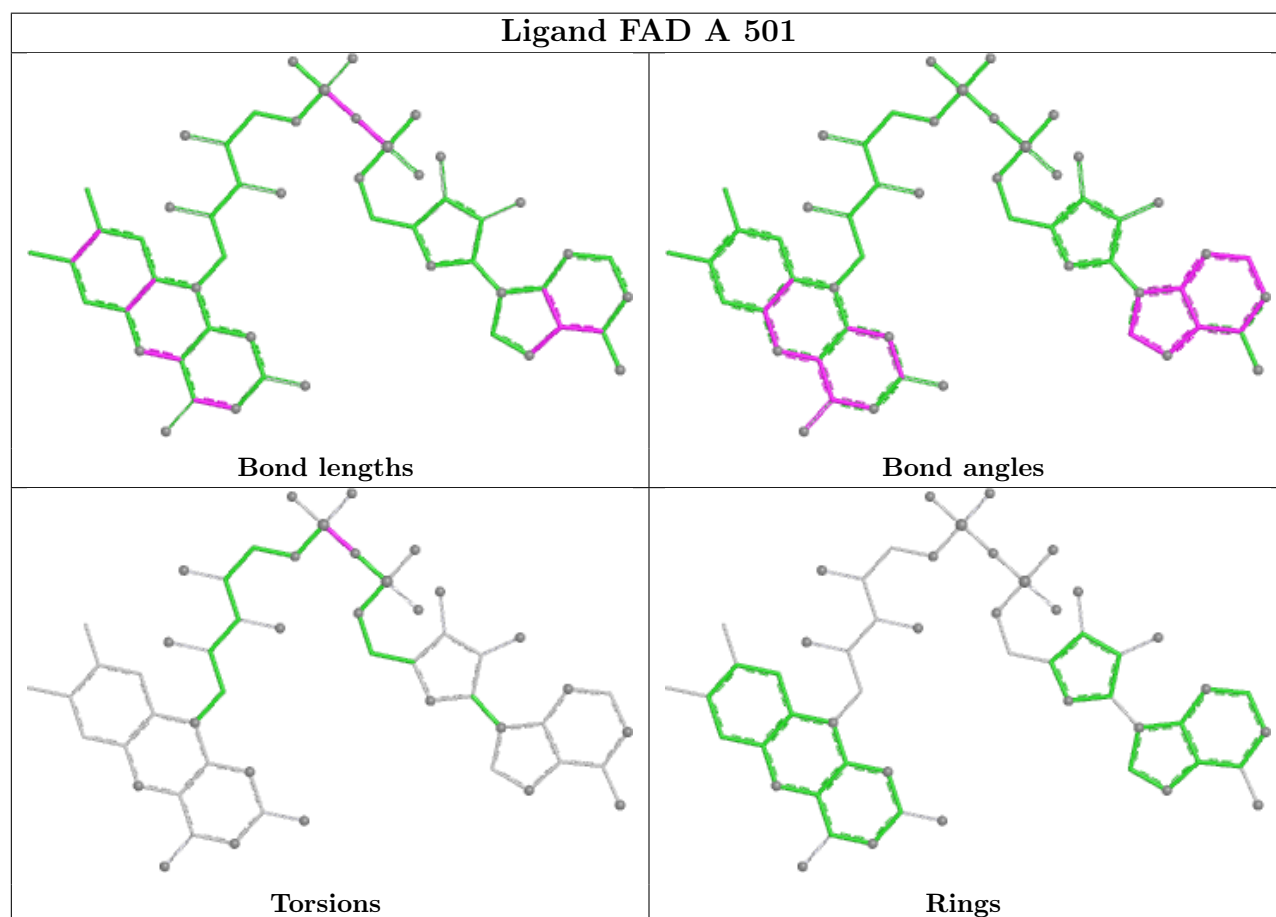
Mol	Chain	Res	Type	Atoms
3	A	502[B]	R57	CAU-CAV-OAW-CAX
3	B	502	R57	CAU-CAV-OAW-CAX
3	A	502[B]	R57	CAY-CAV-OAW-CAX
3	B	502	R57	CAY-CAV-OAW-CAX
2	A	501	FAD	PA-O3P-P-O2P
2	B	501	FAD	PA-O3P-P-O2P
2	A	501	FAD	PA-O3P-P-O1P
2	B	501	FAD	PA-O3P-P-O1P

There are no ring outliers.

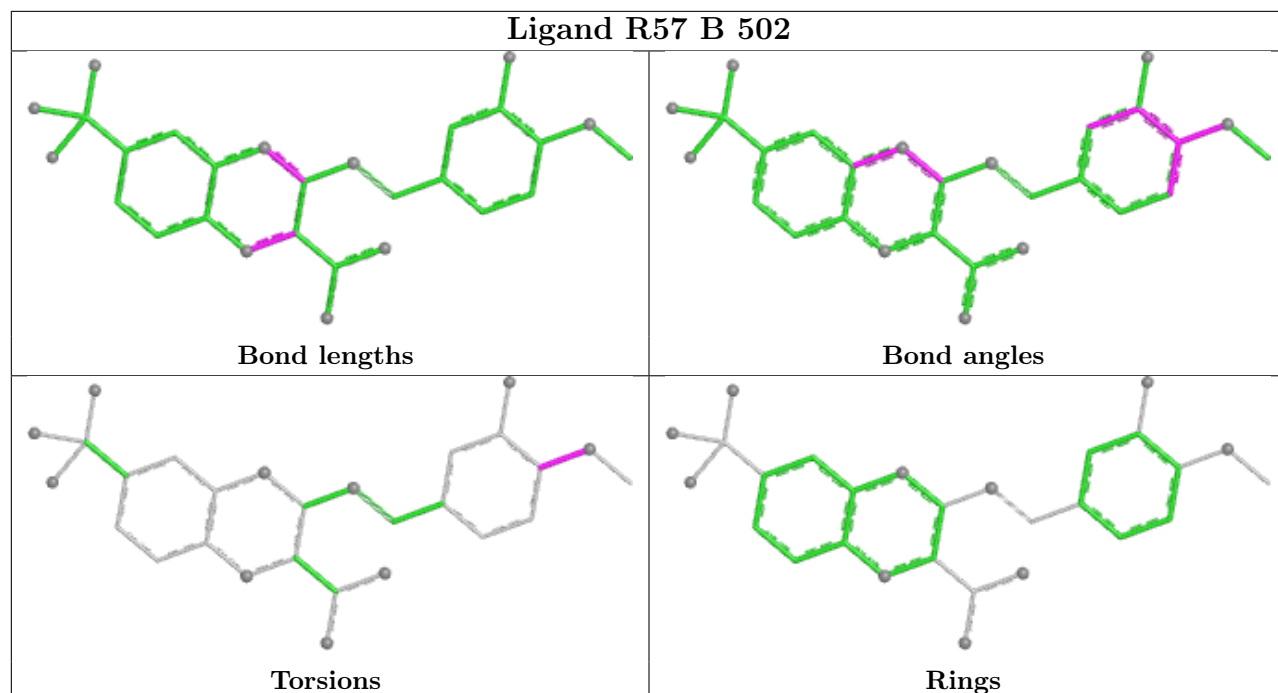
1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502[B]	R57	7	0

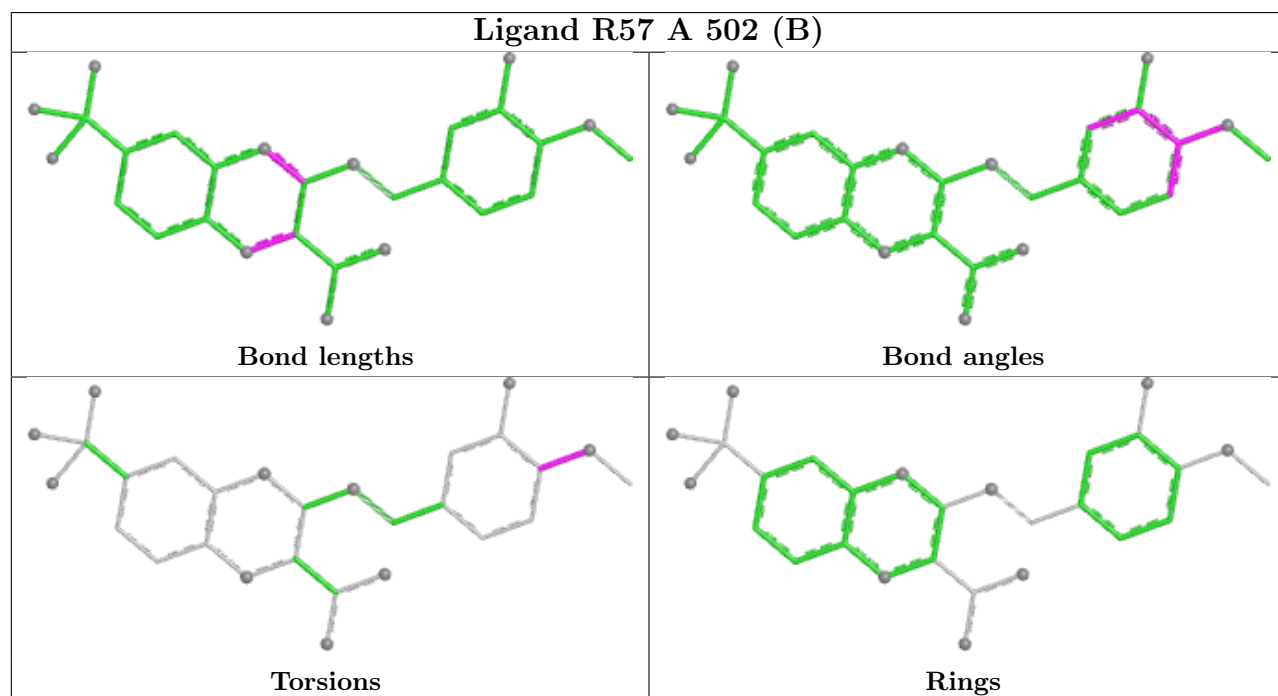
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

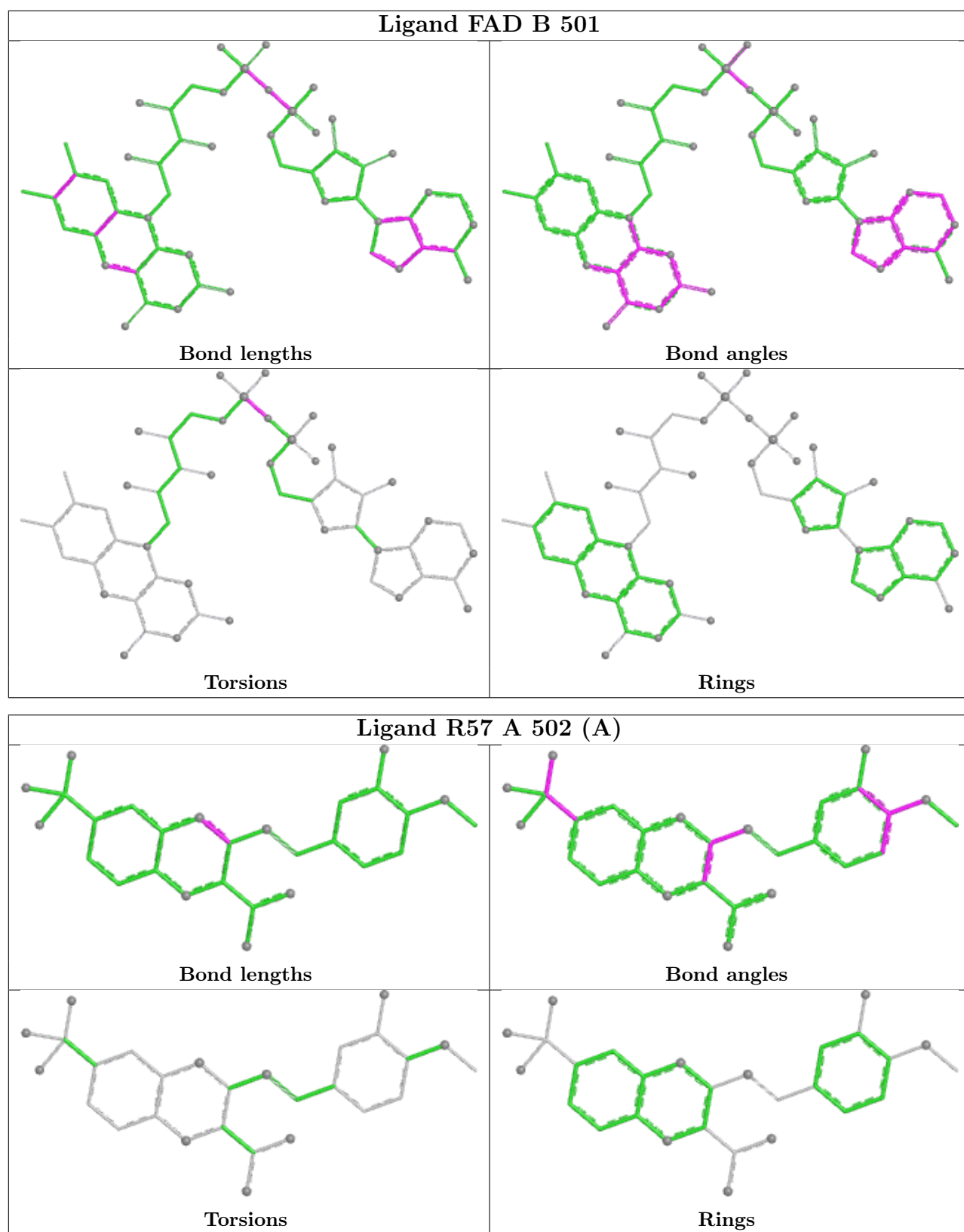


Ligand R57 B 502



Ligand R57 A 502 (B)





5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	454/480 (94%)	0.40	37 (8%) 18 16	16, 27, 53, 70	1 (0%)
1	B	419/480 (87%)	0.80	46 (10%) 10 9	14, 33, 55, 75	2 (0%)
All	All	873/960 (90%)	0.59	83 (9%) 14 12	14, 30, 55, 75	3 (0%)

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	291	PRO	6.5
1	B	292	ILE	5.0
1	B	267	PHE	4.9
1	A	230[A]	TRP	4.3
1	B	269	ALA	4.2
1	B	46	GLY	4.1
1	B	158	ILE	4.1
1	B	230	TRP	4.0
1	B	20	ALA	4.0
1	A	323	TRP	3.9
1	A	270	PRO	3.8
1	B	316	PRO	3.8
1	B	362	PHE	3.8
1	A	4	VAL	3.6
1	B	47	GLY	3.6
1	A	267	PHE	3.5
1	A	272	LEU	3.3
1	B	36	VAL	3.3
1	A	317	LEU	3.1
1	B	461	LEU	3.1
1	A	7	THR	3.1
1	A	320	PHE	3.1
1	B	315	HIS	3.0
1	B	13	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	293	GLY	2.9
1	A	47	GLY	2.9
1	B	48	GLY	2.9
1	A	325	ARG	2.8
1	B	23	VAL	2.8
1	A	119	ARG	2.8
1	B	268	ASP	2.8
1	A	6	ALA	2.7
1	A	280	PRO	2.7
1	B	8	THR	2.7
1	B	25	ASN	2.7
1	A	321	GLY	2.6
1	B	26	VAL	2.6
1	B	314	TYR	2.6
1	A	362	PHE	2.6
1	B	19	THR	2.6
1	B	16	TRP	2.6
1	B	70	LEU	2.6
1	A	269	ALA	2.6
1	B	10	ALA	2.6
1	A	9	THR	2.5
1	B	42	VAL	2.5
1	A	297	TYR	2.5
1	B	392	ILE	2.5
1	B	27	LEU	2.4
1	A	331	GLY	2.4
1	A	319	MET	2.4
1	A	46	GLY	2.4
1	A	278	VAL	2.4
1	B	262	SER	2.4
1	A	10	ALA	2.3
1	A	19	THR	2.3
1	A	324	ASN	2.3
1	A	44	GLU	2.3
1	A	316	PRO	2.3
1	B	447	LEU	2.3
1	B	420	SER	2.3
1	B	34	MET	2.3
1	A	326	ALA	2.2
1	A	461	LEU	2.2
1	B	45	SER	2.2
1	A	273	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	52	ILE	2.2
1	B	14	THR	2.2
1	B	18	ARG	2.2
1	B	360	TYR	2.2
1	B	438	ILE	2.2
1	A	25	ASN	2.1
1	B	15	GLY	2.1
1	A	8	THR	2.1
1	B	35	ILE	2.1
1	A	365	VAL	2.1
1	A	34	MET	2.1
1	A	283	LEU	2.1
1	B	29	THR	2.1
1	A	281	ASN	2.0
1	B	460	LEU	2.0
1	B	60	TYR	2.0
1	B	253	VAL	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

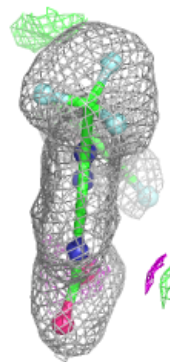
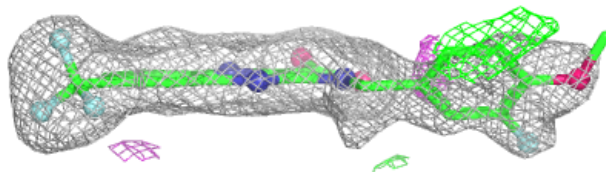
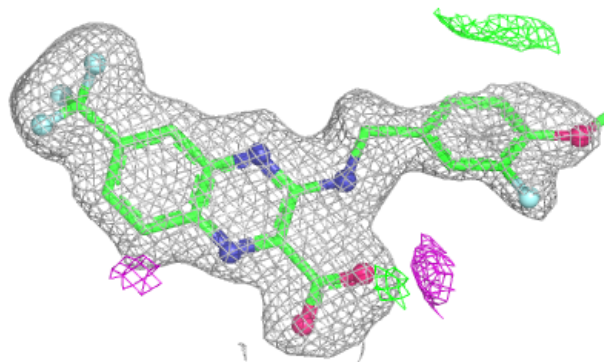
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	R57	A	502[A]	28/28	0.90	0.12	27,28,33,33	28
3	R57	A	502[B]	28/28	0.90	0.12	28,30,39,39	28
3	R57	B	502	28/28	0.91	0.13	33,37,44,44	8
4	IMD	B	503	5/5	0.94	0.08	36,36,37,37	0
2	FAD	B	501	53/53	0.97	0.06	21,27,29,31	0
2	FAD	A	501	53/53	0.98	0.06	18,20,22,22	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.

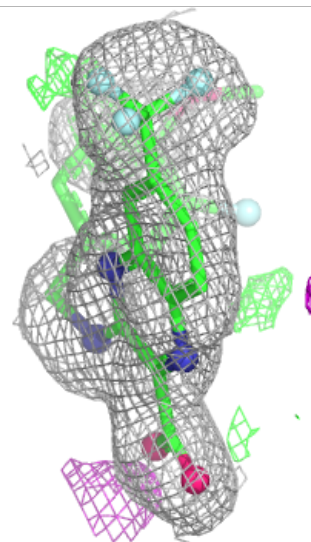
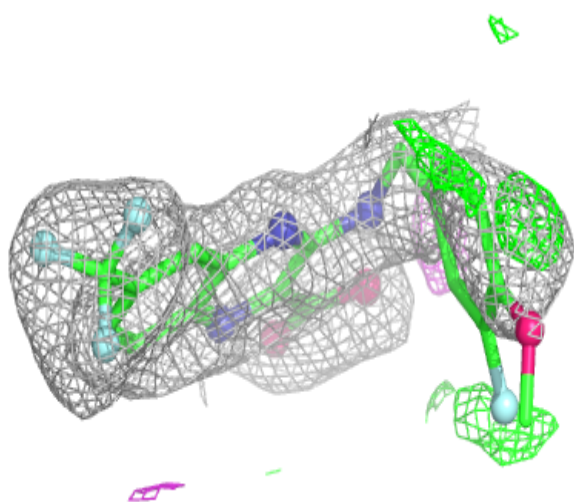
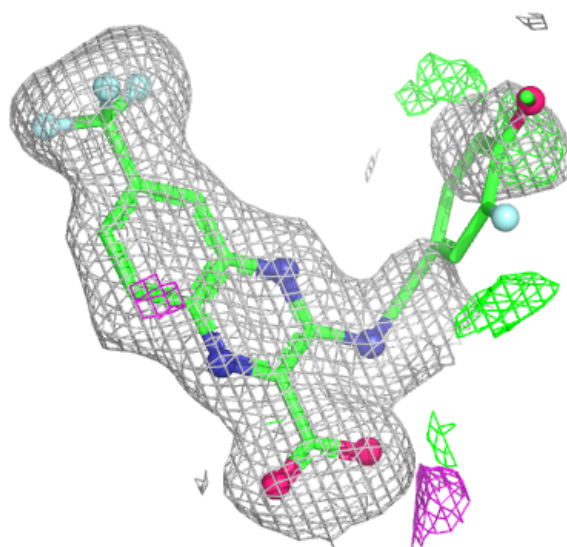
Electron density around R57 A 502 (A):

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



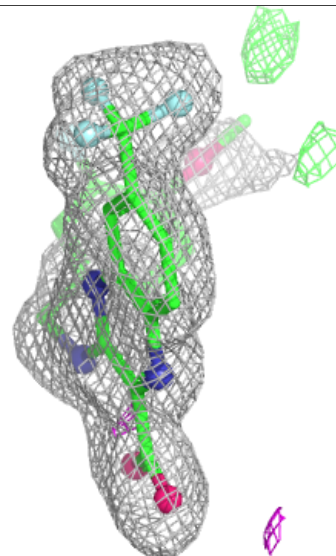
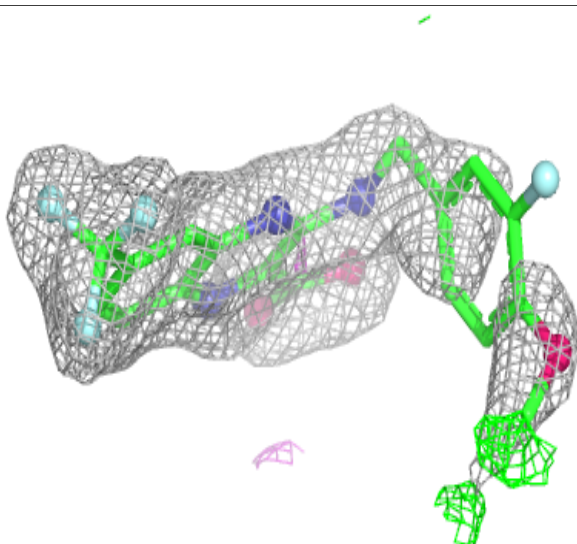
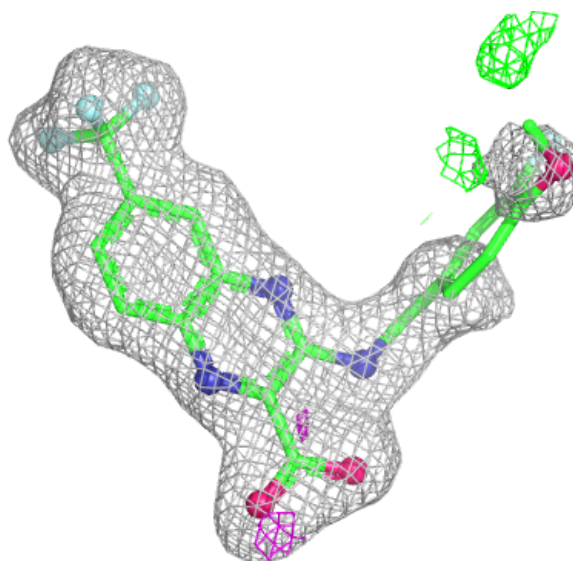
Electron density around R57 A 502 (B):

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



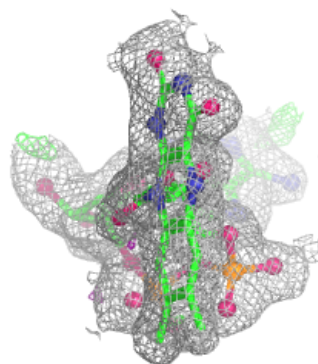
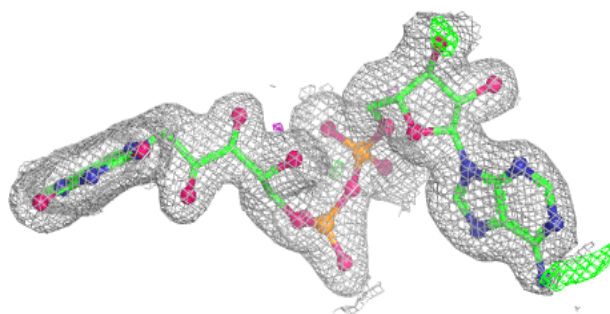
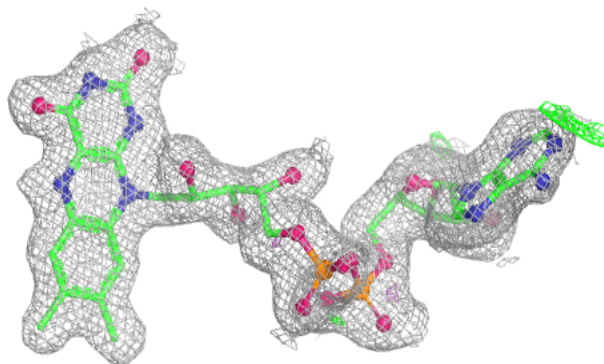
Electron density around R57 B 502:

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

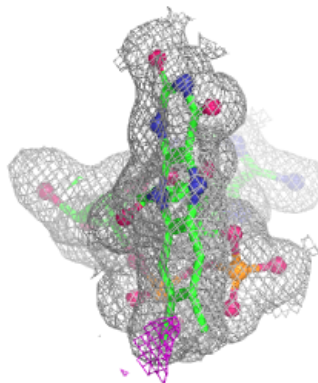
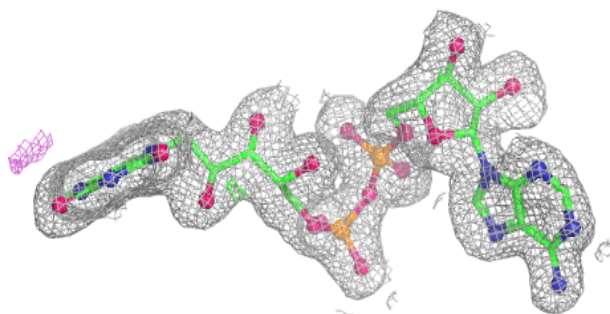
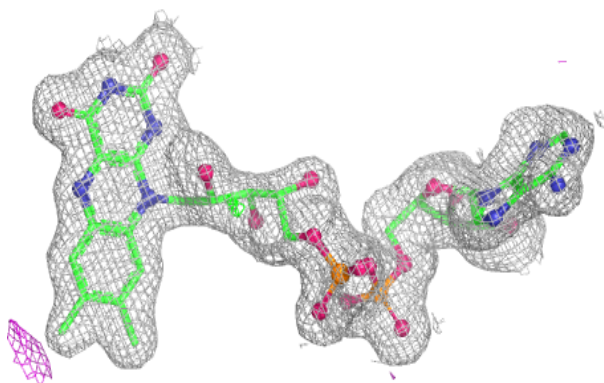


Electron density around FAD B 501:

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

**Electron density around FAD A 501:**

$2mF_o-DF_c$ (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.