



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:36 AM UTC

PDB ID : 3ZNO / pdb_00003zno
Title : IN VITRO AND IN VIVO INHIBITION OF HUMAN D-AMINO ACID OXIDASE: REGULATION OF D-SERINE CONCENTRATION IN THE BRAIN
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Deposited on : 2013-02-15
Resolution : 2.30 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)

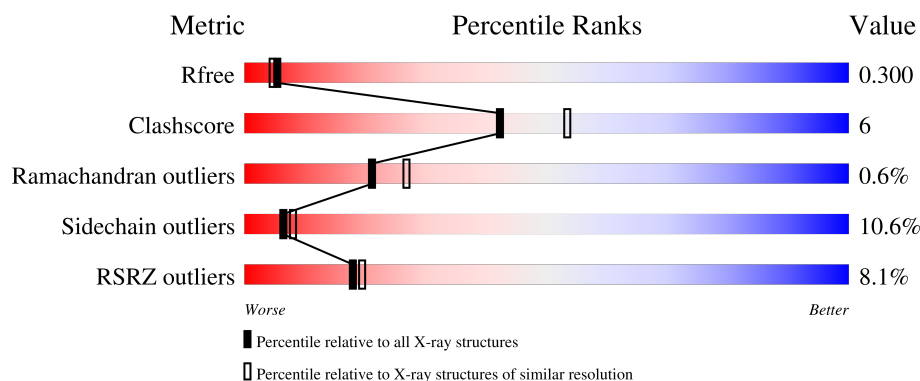
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	347	 5% 72% 22% . .
1	B	347	 11% 78% 17% . .

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

2 Entry composition [i](#)

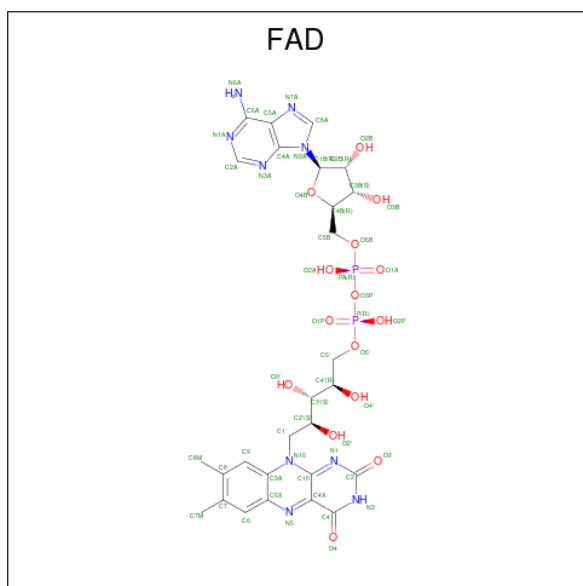
There are 5 unique types of molecules in this entry. The entry contains 5931 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 D-AMINO-ACID OXIDASE.

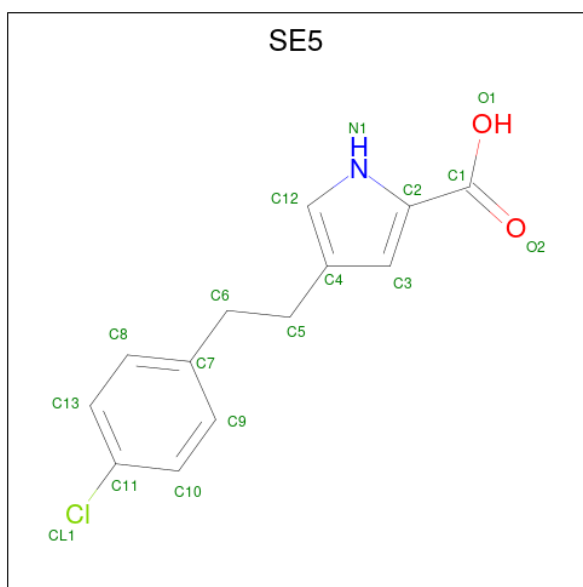
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	340	Total	C	N	O	S	0	8	0
			2789	1791	488	501	9			
1	B	340	Total	C	N	O	S	0	2	0
			2749	1761	485	494	9			

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



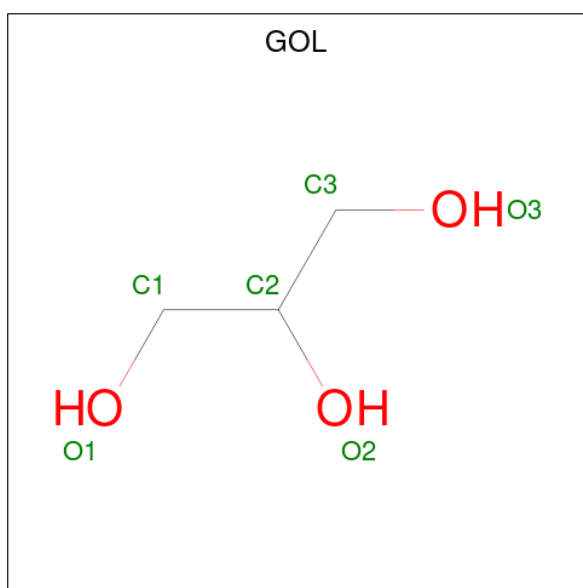
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is 4-(4-chlorophenethyl)-1H-pyrrole-2-carboxylic acid (CCD ID: SE5) (formula: $C_{13}H_{12}ClNO_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0	0
			17	13	1	1	2		
3	B	1	Total	C	Cl	N	O	0	0
			17	13	1	1	2		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		

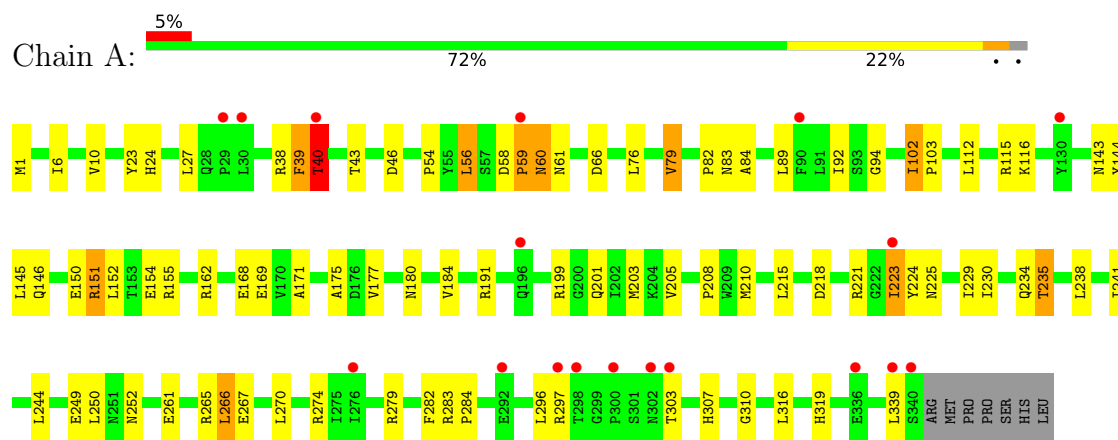
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	131	Total 139	O 139	0	8
5	B	104	Total 108	O 108	0	4

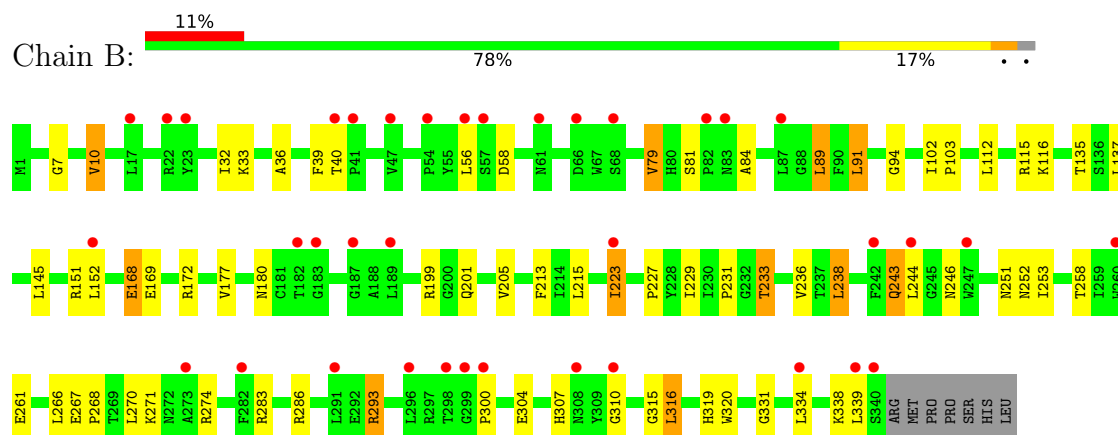
3 Residue-property plots

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: D-AMINO-ACID OXIDASE



• Molecule 1: D-AMINO-ACID OXIDASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	44.03Å 61.06Å 259.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.25 – 2.30 39.25 – 2.30	Depositor EDS
% Data completeness (in resolution range)	90.6 (39.25-2.30) 90.6 (39.25-2.30)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 2.29Å)	Xtriage
Refinement program	BUSTER 2.11.1	Depositor
R, R_{free}	0.217 , 0.289 0.226 , 0.300	Depositor DCC
R_{free} test set	1471 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	40.3	Xtriage
Anisotropy	0.392	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 71.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5931	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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: FAD, SE5, GOL

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	0.89	0/2889	1.37	22/3929 (0.6%)
1	B	0.81	0/2832	1.31	15/3852 (0.4%)
All	All	0.85	0/5721	1.34	37/7781 (0.5%)

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	60	ASN	CA-CB-CG	8.79	121.39	112.60
1	A	23	TYR	N-CA-C	8.25	124.35	113.30
1	A	59	PRO	CA-C-N	7.74	135.63	121.70
1	A	59	PRO	C-N-CA	7.74	135.63	121.70
1	B	223	ILE	N-CA-C	7.35	118.48	109.30
1	B	84	ALA	CA-C-N	6.75	129.33	120.28
1	B	84	ALA	C-N-CA	6.75	129.33	120.28
1	A	297	ARG	N-CA-C	6.65	121.61	112.90
1	B	168	GLU	CA-C-N	6.27	128.68	120.28
1	B	168	GLU	C-N-CA	6.27	128.68	120.28
1	B	227	PRO	N-CA-C	6.04	121.10	111.26
1	A	39	PHE	CA-C-N	5.96	136.34	121.80
1	A	39	PHE	C-N-CA	5.96	136.34	121.80
1	A	94	GLY	N-CA-C	5.87	119.17	110.42
1	B	213	PHE	CA-CB-CG	5.87	119.67	113.80
1	B	10	VAL	N-CA-CB	5.79	116.94	110.51
1	A	84	ALA	CA-C-N	5.71	127.93	120.28
1	A	84	ALA	C-N-CA	5.71	127.93	120.28
1	A	54	PRO	CA-C-N	5.66	128.88	120.90
1	A	54	PRO	C-N-CA	5.66	128.88	120.90
1	B	316	LEU	CA-C-N	5.57	127.74	120.28
1	B	316	LEU	C-N-CA	5.57	127.74	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	251	ASN	CA-CB-CG	5.56	118.16	112.60
1	A	151	ARG	CA-C-N	5.42	127.49	120.44
1	A	151	ARG	C-N-CA	5.42	127.49	120.44
1	B	258	THR	CA-C-N	5.35	127.39	120.70
1	B	258	THR	C-N-CA	5.35	127.39	120.70
1	A	79	VAL	N-CA-C	5.34	118.55	111.17
1	B	94	GLY	N-CA-C	5.30	117.74	110.69
1	A	230	ILE	N-CA-C	5.27	113.37	108.15
1	A	6	ILE	N-CA-CB	5.22	117.94	111.41
1	A	66	ASP	CA-CB-CG	5.21	117.81	112.60
1	A	169	GLU	CA-C-N	5.17	129.18	120.29
1	A	169	GLU	C-N-CA	5.17	129.18	120.29
1	B	320	TRP	N-CA-C	5.15	116.97	111.36
1	A	40	THR	CA-C-O	-5.07	113.22	120.16
1	A	39	PHE	CB-CA-C	5.03	120.43	110.42

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	2789	0	2748	42	0
1	B	2749	0	2706	28	0
2	A	53	0	31	4	0
2	B	53	0	31	2	0
3	A	17	0	11	5	0
3	B	17	0	11	0	0
4	A	6	0	8	3	0
5	A	139	0	0	3	0
5	B	108	0	0	2	0
All	All	5931	0	5546	71	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 6.

All (71) close contacts within the same asymmetric unit are listed below, sorted by their clash

magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:ILE:HD11	3:A:360:SE5:CL1	2.26	0.72
1:B:238:LEU:HG	5:B:524:HOH:O	1.94	0.68
1:B:102:ILE:HG13	1:B:103:PRO:HD2	1.76	0.67
1:A:223:ILE:HG23	1:A:224:TYR:CD2	2.29	0.67
1:A:59:PRO:HG2	1:A:61:ASN:H	1.60	0.66
1:A:102:ILE:HG12	1:A:103:PRO:HD2	1.78	0.66
1:A:180:ASN:HD22	1:A:307:HIS:HD2	1.41	0.66
1:A:218:ASP:HB3	1:A:221[B]:ARG:HG2	1.83	0.60
1:A:283:ARG:NE	3:A:360:SE5:O2	2.27	0.59
1:B:199:ARG:HH22	1:B:201:GLN:HE21	1.49	0.59
1:A:171:ALA:HB1	1:A:303:THR:HG21	1.85	0.58
1:B:7:GLY:HA2	2:B:350:FAD:H1B	1.87	0.57
1:A:76:LEU:O	1:A:79:VAL:HG12	2.06	0.55
4:A:1341:GOL:H32	1:B:91:LEU:HD12	1.89	0.55
1:A:184:VAL:HG11	1:A:282:PHE:HB2	1.89	0.55
1:B:199:ARG:HH22	1:B:201:GLN:NE2	2.06	0.54
1:A:210:MET:HE1	1:A:267:GLU:HG2	1.89	0.54
1:B:283:ARG:HG3	2:B:350:FAD:HM81	1.90	0.54
1:B:180:ASN:HD22	1:B:307:HIS:HD2	1.57	0.53
1:B:316:LEU:O	1:B:319:HIS:HD2	1.91	0.53
1:B:112:LEU:HB2	1:B:135:THR:HB	1.90	0.53
1:A:1:MET:HG2	1:A:27:LEU:HD13	1.90	0.53
1:A:175:ALA:O	1:A:303:THR:HG22	2.09	0.52
1:A:40:THR:O	1:A:46:ASP:OD1	2.28	0.52
1:A:229:ILE:HD12	1:A:266:LEU:HD13	1.91	0.51
1:A:146:GLN:O	1:A:150:GLU:HG2	2.09	0.51
1:A:201:GLN:HE22	1:A:252:ASN:H	1.58	0.51
1:B:201:GLN:HE22	1:B:252:ASN:H	1.56	0.51
4:A:1341:GOL:H31	1:B:91:LEU:HB2	1.94	0.50
1:B:79:VAL:HG23	1:B:89:LEU:HD22	1.92	0.50
1:B:229:ILE:HG12	5:B:524:HOH:O	2.11	0.50
1:A:199:ARG:HH22	1:A:201:GLN:NE2	2.11	0.49
1:A:115:ARG:NH1	5:A:615:HOH:O	2.46	0.48
1:A:40:THR:O	1:A:46:ASP:CG	2.57	0.48
1:A:283:ARG:HG3	2:A:350:FAD:HM81	1.96	0.47
2:A:350:FAD:C4	3:A:360:SE5:C12	2.93	0.47
1:A:39:PHE:O	1:A:43:THR:HG23	2.15	0.47
1:A:223:ILE:HD12	1:A:224:TYR:H	1.78	0.47
1:A:58:ASP:OD1	1:A:59:PRO:HD2	2.15	0.47
1:B:79:VAL:HG21	1:B:137:LEU:HD22	1.97	0.47
1:B:315:GLY:O	1:B:319:HIS:HB3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:1341:GOL:C3	1:B:91:LEU:HD12	2.46	0.46
1:B:177:VAL:HG22	1:B:304:GLU:HB2	1.97	0.46
1:A:150:GLU:O	1:A:154:GLU:HG3	2.15	0.46
1:A:235:THR:HG22	5:A:580:HOH:O	2.16	0.46
1:A:199:ARG:HH22	1:A:201:GLN:HE21	1.61	0.46
1:A:76:LEU:HD12	1:A:79:VAL:HG11	1.99	0.45
1:B:268:PRO:O	1:B:271:LYS:HD3	2.17	0.45
1:A:162:ARG:O	2:A:350:FAD:H2A	2.16	0.45
1:A:59:PRO:CG	1:A:61:ASN:H	2.29	0.45
1:B:231:PRO:HA	1:B:236:VAL:HG22	1.98	0.45
1:B:331:GLY:HA2	1:B:334:LEU:HD12	1.97	0.45
1:A:203:MET:HB3	1:A:238:LEU:HB2	1.99	0.44
1:B:36:ALA:HB3	1:B:39:PHE:CZ	2.53	0.44
1:B:293:ARG:H	1:B:293:ARG:HG3	1.71	0.43
1:B:10:VAL:HG21	1:B:316:LEU:HD13	2.00	0.43
1:A:144:TYR:OH	1:A:319:HIS:HE1	2.02	0.43
2:A:350:FAD:C4X	3:A:360:SE5:C12	2.96	0.43
1:A:56:LEU:HB2	3:A:360:SE5:CL1	2.56	0.42
1:A:10:VAL:HG21	1:A:316[A]:LEU:HD13	2.01	0.42
1:A:208:PRO:HB2	1:B:233:THR:O	2.20	0.42
1:A:1:MET:HE1	1:A:177:VAL:HG23	2.02	0.42
1:A:223:ILE:C	1:A:225:ASN:H	2.26	0.42
1:A:210:MET:CE	1:A:267:GLU:HG2	2.50	0.41
1:A:261:GLU:O	1:A:265:ARG:HD3	2.19	0.41
1:A:59:PRO:HG3	5:A:531:HOH:O	2.20	0.41
1:A:284:PRO:HB2	1:A:310:GLY:HA2	2.02	0.41
1:B:243:GLN:HE21	1:B:246:ASN:HD22	1.69	0.40
1:A:144:TYR:OH	1:A:319:HIS:CE1	2.74	0.40
1:B:267:GLU:HB3	1:B:270:LEU:HD13	2.03	0.40
1:B:286:ARG:HG3	1:B:310:GLY:HA3	2.04	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	346/347 (100%)	327 (94%)	17 (5%)	2 (1%)	21	27
1	B	340/347 (98%)	322 (95%)	16 (5%)	2 (1%)	21	27
All	All	686/694 (99%)	649 (95%)	33 (5%)	4 (1%)	21	27

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	40	THR
1	B	300	PRO
1	B	40	THR
1	A	82	PRO

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	300/299 (100%)	267 (89%)	33 (11%)	6	7
1	B	294/299 (98%)	264 (90%)	30 (10%)	7	8
All	All	594/598 (99%)	531 (89%)	63 (11%)	6	8

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24[A]	HIS
1	A	24[B]	HIS
1	A	38	ARG
1	A	56	LEU
1	A	60	ASN
1	A	83	ASN
1	A	89	LEU
1	A	92	ILE
1	A	102	ILE

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Mol	Chain	Res	Type
1	A	112	LEU
1	A	116	LYS
1	A	143	ASN
1	A	145	LEU
1	A	151	ARG
1	A	152	LEU
1	A	155	ARG
1	A	168	GLU
1	A	191	ARG
1	A	205	VAL
1	A	215	LEU
1	A	223	ILE
1	A	234	GLN
1	A	235	THR
1	A	241	ILE
1	A	244	LEU
1	A	249	GLU
1	A	250	LEU
1	A	266	LEU
1	A	270	LEU
1	A	274	ARG
1	A	279	ARG
1	A	296	LEU
1	A	339	LEU
1	B	32	ILE
1	B	33	LYS
1	B	56	LEU
1	B	58	ASP
1	B	79	VAL
1	B	81	SER
1	B	89	LEU
1	B	91	LEU
1	B	115	ARG
1	B	116	LYS
1	B	145	LEU
1	B	151	ARG
1	B	152	LEU
1	B	168	GLU
1	B	169	GLU
1	B	172	ARG
1	B	205	VAL
1	B	215	LEU

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Mol	Chain	Res	Type
1	B	223	ILE
1	B	233	THR
1	B	238	LEU
1	B	243	GLN
1	B	244	LEU
1	B	253	ILE
1	B	261	GLU
1	B	266	LEU
1	B	274	ARG
1	B	293	ARG
1	B	338	LYS
1	B	339	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	GLN
1	A	70	GLN
1	A	78	HIS
1	A	143	ASN
1	A	201	GLN
1	A	243	GLN
1	A	252	ASN
1	A	307	HIS
1	A	319	HIS
1	B	20	HIS
1	B	83	ASN
1	B	161	GLN
1	B	201	GLN
1	B	225	ASN
1	B	243	GLN
1	B	251	ASN
1	B	252	ASN
1	B	272	ASN
1	B	307	HIS
1	B	319	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 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	$\# Z > 2$
3	SE5	A	360	-	18,18,18	1.03	2 (11%)	24,24,24	1.18	3 (12%)
3	SE5	B	360	-	18,18,18	1.03	2 (11%)	24,24,24	1.23	3 (12%)
2	FAD	B	350	-	58,58,58	3.31	26 (44%)	85,89,89	2.16	28 (32%)
2	FAD	A	350	-	58,58,58	3.57	27 (46%)	85,89,89	2.62	34 (40%)
4	GOL	A	1341	-	5,5,5	1.16	0	5,5,5	1.12	1 (20%)

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	SE5	A	360	-	-	0/9/9/9	0/2/2/2
3	SE5	B	360	-	-	2/9/9/9	0/2/2/2
2	FAD	B	350	-	-	5/34/50/50	0/6/6/6
2	FAD	A	350	-	-	5/34/50/50	0/6/6/6
4	GOL	A	1341	-	-	4/4/4/4	-

All (57) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	350	FAD	C6-C7	12.93	1.57	1.39
2	B	350	FAD	C6-C7	12.38	1.56	1.39
2	A	350	FAD	C9A-C5X	10.54	1.58	1.41
2	B	350	FAD	C9A-C5X	9.47	1.56	1.41
2	B	350	FAD	P-O3P	7.94	1.68	1.59
2	A	350	FAD	C5A-N7A	-6.74	1.26	1.39
2	A	350	FAD	C9-C8	6.69	1.48	1.39
2	B	350	FAD	C9-C8	6.55	1.48	1.39
2	B	350	FAD	C4X-N5	5.94	1.43	1.30
2	A	350	FAD	C5A-C4A	5.81	1.49	1.39
2	A	350	FAD	O4-C4	5.65	1.34	1.23
2	A	350	FAD	C8M-C8	-5.55	1.40	1.51
2	A	350	FAD	C4X-N5	5.26	1.42	1.30
2	A	350	FAD	P-O3P	5.18	1.65	1.59
2	B	350	FAD	C5A-C4A	4.97	1.48	1.39
2	A	350	FAD	C4'-C3'	4.91	1.62	1.53
2	B	350	FAD	C5A-N7A	-4.18	1.31	1.39
2	A	350	FAD	O2'-C2'	4.18	1.52	1.43
2	B	350	FAD	O4-C4	4.07	1.31	1.23
2	A	350	FAD	C8A-N7A	3.98	1.39	1.31
2	B	350	FAD	C10-N1	3.95	1.41	1.33
2	A	350	FAD	C9-C9A	3.79	1.45	1.39
2	B	350	FAD	C9-C9A	3.75	1.45	1.39
2	B	350	FAD	C9A-N10	3.63	1.47	1.41
2	B	350	FAD	C8M-C8	-3.57	1.44	1.51
2	A	350	FAD	O2-C2	3.53	1.31	1.24
2	B	350	FAD	C1'-C2'	3.37	1.57	1.52
2	A	350	FAD	C1'-N10	3.34	1.55	1.47
2	B	350	FAD	O2'-C2'	3.33	1.50	1.43
2	B	350	FAD	O2-C2	3.31	1.30	1.24
2	A	350	FAD	C1'-C2'	3.04	1.56	1.52
2	A	350	FAD	C2B-C3B	-3.01	1.45	1.53
2	B	350	FAD	C8A-N7A	2.96	1.37	1.31
2	A	350	FAD	C4A-N9A	-2.90	1.31	1.37
2	A	350	FAD	O3'-C3'	2.88	1.50	1.43
2	B	350	FAD	C4A-N9A	-2.84	1.31	1.37
2	B	350	FAD	C4'-C3'	2.84	1.58	1.53
2	B	350	FAD	C4A-N3A	2.71	1.39	1.34
2	A	350	FAD	C5X-N5	2.67	1.44	1.39
2	B	350	FAD	O4B-C1B	2.57	1.47	1.42
2	A	350	FAD	O4'-C4'	2.50	1.48	1.43
2	B	350	FAD	C7M-C7	2.49	1.55	1.51
3	A	360	SE5	C2-N1	-2.48	1.34	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	350	FAD	C9A-N10	2.45	1.45	1.41
2	A	350	FAD	O5B-C5B	2.42	1.53	1.44
2	A	350	FAD	C2A-N1A	2.41	1.38	1.33
2	B	350	FAD	C8-C7	2.39	1.46	1.40
2	A	350	FAD	PA-O3P	2.38	1.62	1.59
3	B	360	SE5	C2-N1	-2.35	1.34	1.37
3	B	360	SE5	C11-CL1	2.33	1.79	1.74
3	A	360	SE5	C11-CL1	2.30	1.79	1.74
2	B	350	FAD	C10-N10	2.29	1.42	1.37
2	A	350	FAD	C10-N1	2.19	1.37	1.33
2	B	350	FAD	C2A-N1A	2.08	1.37	1.33
2	B	350	FAD	P-O5'	2.08	1.67	1.59
2	B	350	FAD	O4B-C4B	2.05	1.49	1.45
2	A	350	FAD	P-O1P	2.03	1.57	1.50

All (69) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	350	FAD	O4'-C4'-C5'	-8.93	90.29	109.99
2	B	350	FAD	O4'-C4'-C5'	-6.76	95.08	109.99
2	A	350	FAD	C4A-C5A-N7A	-6.59	103.05	110.58
2	A	350	FAD	O3P-PA-O1A	-6.03	92.58	110.70
2	A	350	FAD	C9A-C5X-N5	-5.79	116.31	122.45
2	A	350	FAD	C5A-C4A-N9A	5.70	112.03	105.81
2	A	350	FAD	C5A-C4A-N3A	-5.62	118.98	126.72
2	B	350	FAD	C9A-C5X-N5	-5.43	116.70	122.45
2	B	350	FAD	O2P-P-O3P	5.41	121.88	107.27
2	B	350	FAD	C4A-C5A-N7A	-5.34	104.48	110.58
2	B	350	FAD	C5A-C4A-N3A	-5.09	119.70	126.72
2	A	350	FAD	C5A-N7A-C8A	4.90	111.15	103.45
2	A	350	FAD	O4B-C1B-N9A	-4.53	99.38	108.09
2	A	350	FAD	O2P-P-O3P	4.27	118.80	107.27
2	A	350	FAD	O3'-C3'-C4'	4.19	118.45	108.93
2	A	350	FAD	C8M-C8-C9	-4.14	112.28	119.57
2	A	350	FAD	C4X-C10-N10	3.99	122.19	116.48
2	B	350	FAD	C5A-N7A-C8A	3.81	109.44	103.45
2	B	350	FAD	C5A-C4A-N9A	3.67	109.81	105.81
2	A	350	FAD	N6A-C6A-N1A	-3.31	111.00	118.38
2	A	350	FAD	C6A-C5A-N7A	3.30	138.46	132.09
2	B	350	FAD	C4X-C4-N3	3.28	121.61	113.25
2	B	350	FAD	O3P-PA-O1A	-3.18	101.13	110.70
2	A	350	FAD	C5A-C6A-N6A	3.13	131.03	123.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	350	FAD	O2P-P-O5'	-3.09	93.58	107.57
3	B	360	SE5	C3-C2-C1	-3.03	125.54	132.53
2	B	350	FAD	C5X-N5-C4X	2.99	122.92	118.09
2	B	350	FAD	C6A-C5A-N7A	2.98	137.84	132.09
2	B	350	FAD	O2P-P-O5'	-2.93	94.29	107.57
2	A	350	FAD	C4B-O4B-C1B	-2.87	103.13	109.47
2	A	350	FAD	C9-C8-C7	2.84	123.86	119.69
2	B	350	FAD	C4B-O4B-C1B	-2.82	103.23	109.47
2	B	350	FAD	C2A-N3A-C4A	2.82	118.71	111.83
2	A	350	FAD	O2A-PA-O3P	2.80	114.85	107.27
2	A	350	FAD	C5X-N5-C4X	2.77	122.57	118.09
2	A	350	FAD	O5'-P-O1P	-2.74	98.06	108.94
2	B	350	FAD	O2B-C2B-C3B	2.73	120.56	111.82
2	B	350	FAD	C4-N3-C2	-2.68	120.88	125.64
2	B	350	FAD	N3A-C2A-N1A	-2.67	124.54	128.58
2	A	350	FAD	C10-C4X-N5	-2.63	119.44	124.81
2	A	350	FAD	C4-C4X-N5	2.61	121.81	118.21
3	A	360	SE5	O1-C1-C2	2.59	119.86	114.27
2	A	350	FAD	C2A-N3A-C4A	2.59	118.15	111.83
2	B	350	FAD	O2P-P-O1P	2.58	124.47	112.44
2	A	350	FAD	C6-C5X-N5	2.58	122.72	118.44
2	A	350	FAD	N9A-C8A-N7A	-2.51	110.38	113.94
2	B	350	FAD	C4-C4X-N5	2.49	121.64	118.21
3	A	360	SE5	C3-C2-C1	-2.49	126.80	132.53
2	B	350	FAD	N9A-C8A-N7A	-2.46	110.44	113.94
2	A	350	FAD	C4X-C4-N3	2.44	119.47	113.25
2	B	350	FAD	O2B-C2B-C1B	2.39	118.32	110.10
2	B	350	FAD	O3'-C3'-C2'	-2.32	103.65	108.93
2	A	350	FAD	C4A-N9A-C8A	-2.27	103.35	105.74
2	B	350	FAD	O4-C4-N3	-2.24	115.91	120.11
2	B	350	FAD	C4X-C10-N1	-2.24	119.11	124.59
2	B	350	FAD	C6-C5X-N5	2.23	122.14	118.44
2	A	350	FAD	C4'-C3'-C2'	-2.22	109.88	113.57
3	A	360	SE5	C5-C4-C3	2.21	127.39	123.52
2	B	350	FAD	O4B-C4B-C5B	-2.21	102.25	109.33
2	A	350	FAD	O2B-C2B-C1B	2.21	117.72	110.10
2	B	350	FAD	C8M-C8-C9	-2.20	115.69	119.57
3	B	360	SE5	C1-C2-N1	2.20	125.95	120.56
2	A	350	FAD	C4-N3-C2	-2.18	121.77	125.64
2	A	350	FAD	O4B-C1B-C2B	-2.18	101.95	106.62
2	A	350	FAD	C5'-C4'-C3'	-2.18	108.11	112.22
3	B	360	SE5	O1-C1-C2	2.13	118.87	114.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	350	FAD	N6A-C6A-N1A	-2.12	113.66	118.38
2	A	350	FAD	O2P-P-O1P	2.09	122.15	112.44
4	A	1341	GOL	O3-C3-C2	2.05	119.62	110.38

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	350	FAD	P-O3P-PA-O5B
2	B	350	FAD	C5'-O5'-P-O1P
2	B	350	FAD	C5'-O5'-P-O3P
4	A	1341	GOL	C1-C2-C3-O3
4	A	1341	GOL	O2-C2-C3-O3
2	B	350	FAD	O3'-C3'-C4'-C5'
4	A	1341	GOL	O1-C1-C2-C3
4	A	1341	GOL	O1-C1-C2-O2
2	A	350	FAD	C2'-C3'-C4'-C5'
2	A	350	FAD	PA-O3P-P-O1P
2	B	350	FAD	PA-O3P-P-O2P
3	B	360	SE5	C5-C6-C7-C8
3	B	360	SE5	C5-C6-C7-C9
2	A	350	FAD	PA-O3P-P-O2P
2	A	350	FAD	C1'-C2'-C3'-O3'
2	B	350	FAD	C2'-C3'-C4'-C5'

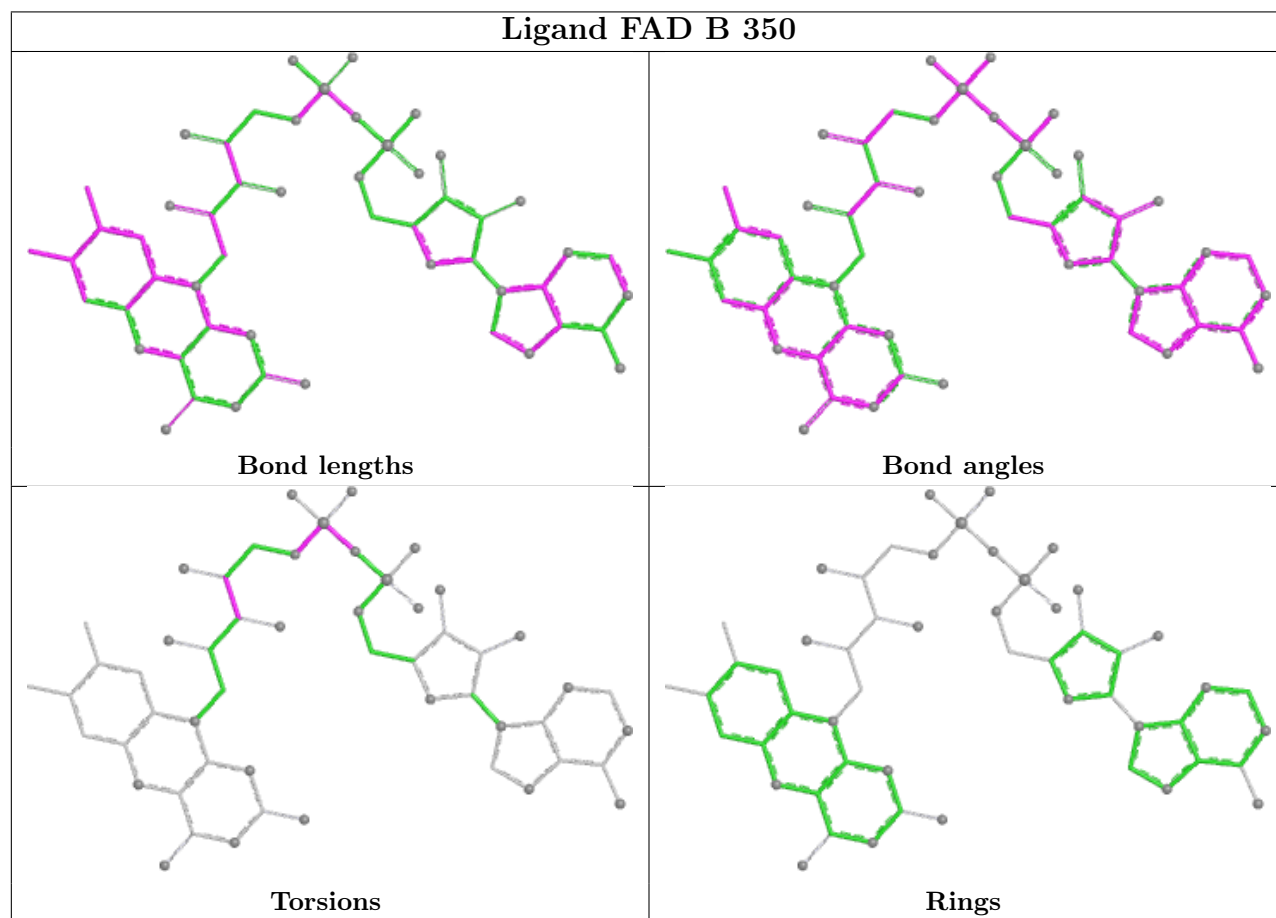
There are no ring outliers.

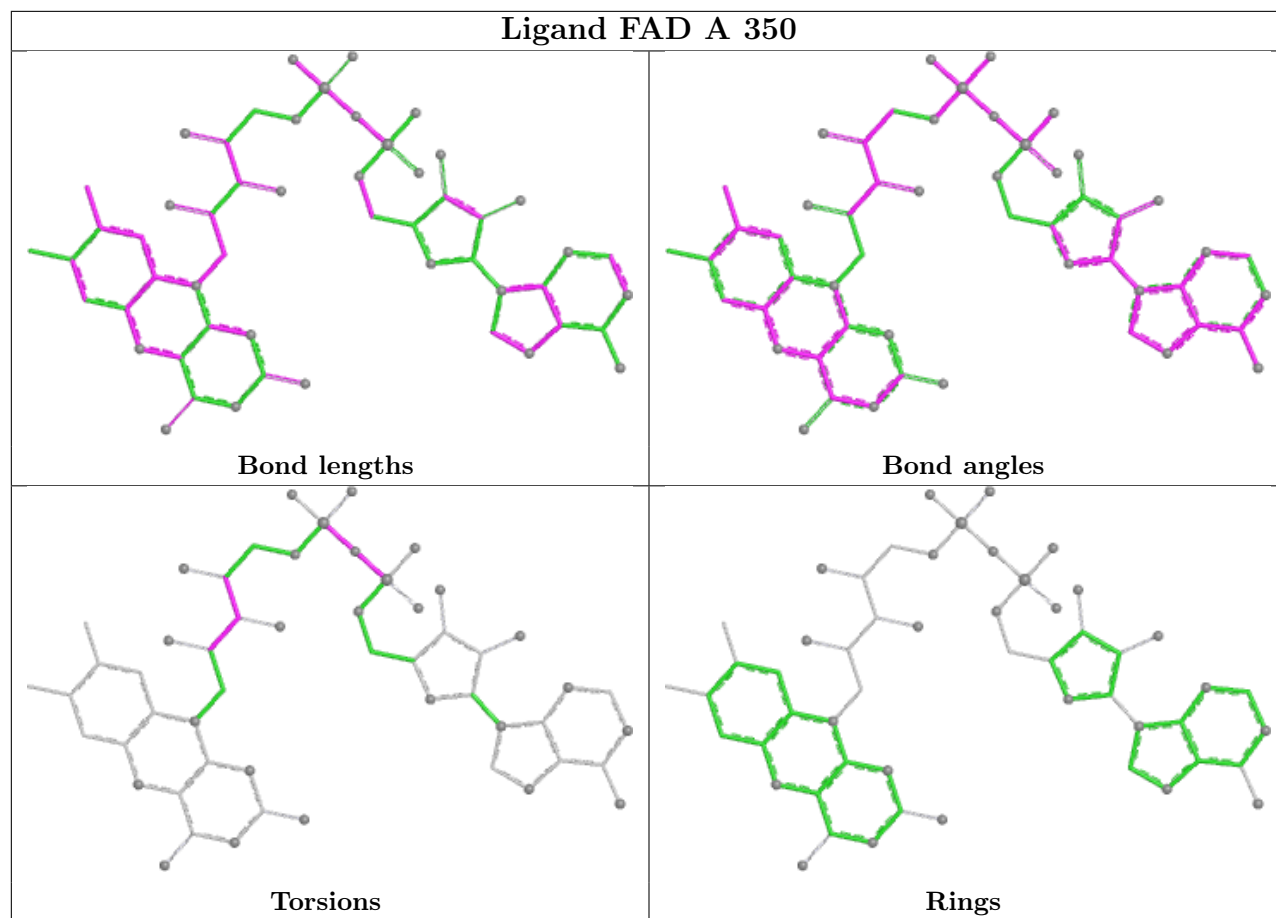
4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	360	SE5	5	0
2	B	350	FAD	2	0
2	A	350	FAD	4	0
4	A	1341	GOL	3	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, 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	340/347 (97%)	0.48	18 (5%)	32	34	16, 39, 69, 118	8 (2%)
1	B	340/347 (97%)	0.95	37 (10%)	10	12	23, 56, 98, 141	2 (0%)
All	All	680/694 (97%)	0.72	55 (8%)	18	19	16, 48, 89, 141	10 (1%)

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	90[A]	PHE	5.5
1	A	339	LEU	4.9
1	B	66	ASP	4.6
1	B	339	LEU	4.3
1	A	340	SER	4.1
1	A	298	THR	3.2
1	A	40	THR	3.2
1	B	300	PRO	3.1
1	B	340	SER	3.1
1	A	300	PRO	3.0
1	A	30	LEU	3.0
1	B	299	GLY	3.0
1	B	82	PRO	3.0
1	A	302	ASN	2.9
1	B	56	LEU	2.9
1	B	310	GLY	2.9
1	A	130	TYR	2.8
1	B	54	PRO	2.8
1	B	41	PRO	2.8
1	B	57	SER	2.8
1	A	223	ILE	2.8
1	A	292[A]	GLU	2.8
1	B	260	TRP	2.7
1	B	68	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	244	LEU	2.7
1	B	291	LEU	2.7
1	A	29	PRO	2.6
1	B	298	THR	2.6
1	B	223	ILE	2.6
1	A	59	PRO	2.5
1	B	40	THR	2.5
1	B	152	LEU	2.5
1	B	187	GLY	2.5
1	A	303	THR	2.4
1	B	183	GLY	2.4
1	B	61	ASN	2.3
1	A	336	GLU	2.3
1	B	273	ALA	2.3
1	B	83	ASN	2.3
1	B	334	LEU	2.2
1	A	297	ARG	2.2
1	B	182	THR	2.2
1	B	247	TRP	2.2
1	B	22	ARG	2.2
1	B	47	VAL	2.2
1	B	282	PHE	2.2
1	B	23	TYR	2.1
1	B	242	PHE	2.1
1	B	296	LEU	2.1
1	A	196	GLN	2.1
1	B	87	LEU	2.1
1	B	189	LEU	2.1
1	B	17	LEU	2.0
1	A	276	ILE	2.0
1	B	308	ASN	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 ⓘ

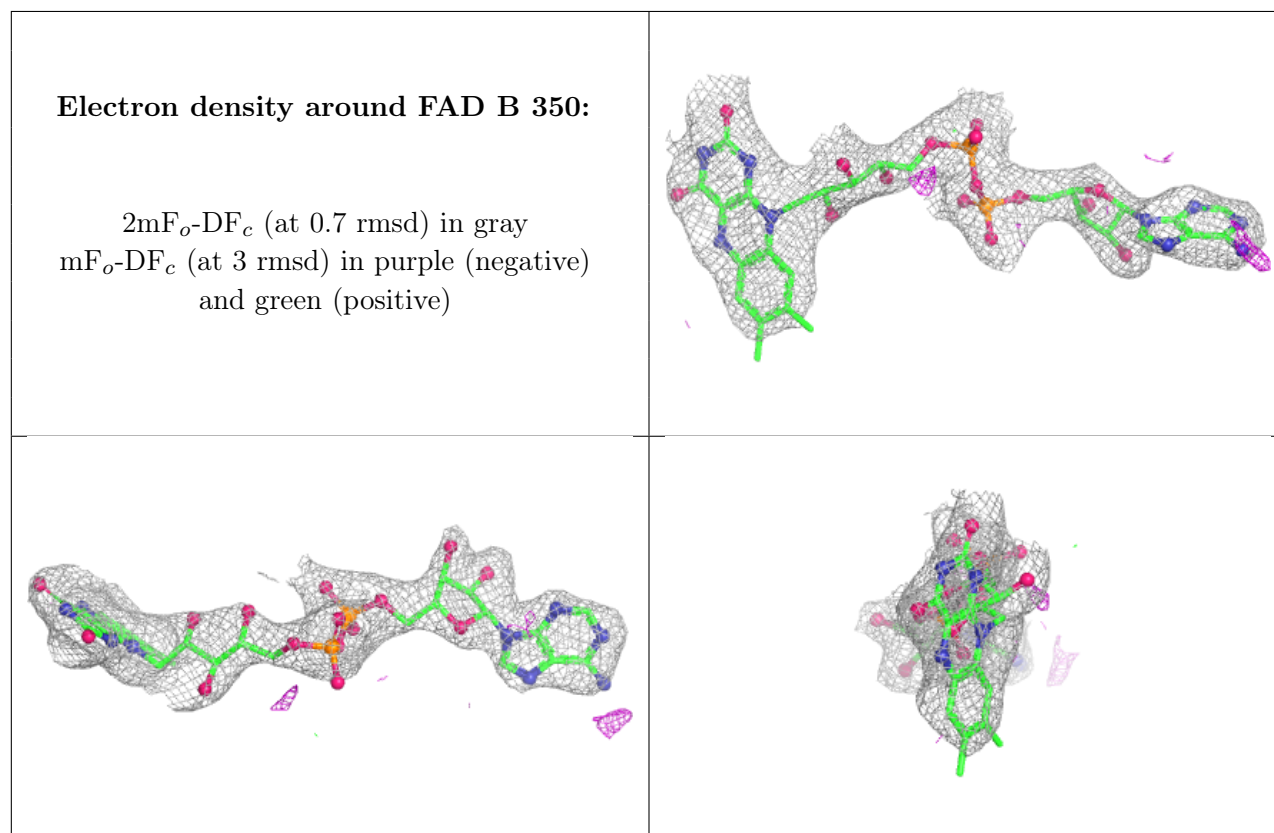
There are no oligosaccharides in this entry.

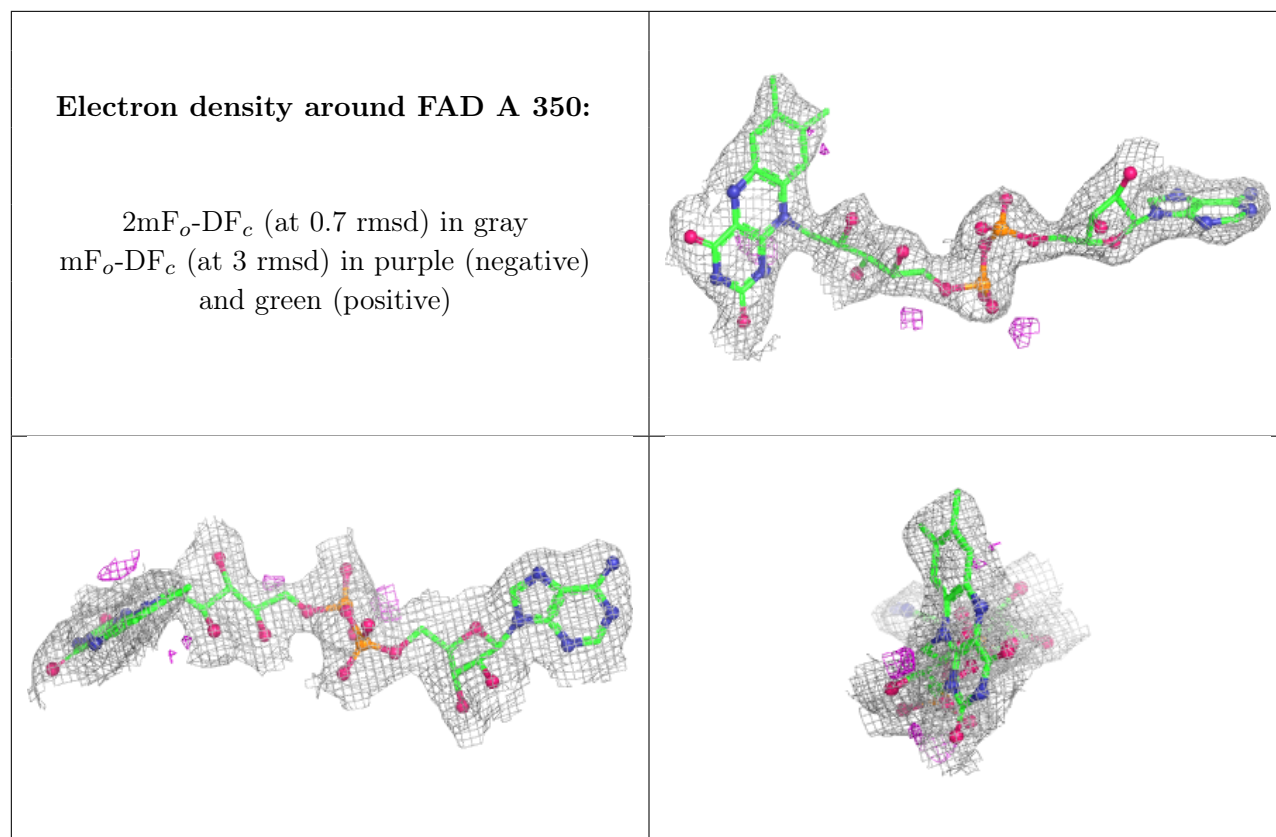
6.4 Ligands ⓘ

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($\text{\AA}^2$)	Q<0.9
4	GOL	A	1341	6/6	0.74	0.18	56,56,56,56	0
3	SE5	B	360	17/17	0.84	0.14	54,59,67,70	0
3	SE5	A	360	17/17	0.87	0.13	30,36,53,58	0
2	FAD	B	350	53/53	0.94	0.09	39,47,56,59	0
2	FAD	A	350	53/53	0.95	0.08	23,30,36,37	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.