



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:22 AM UTC

PDB ID : 4BO2 / pdb_00004bo2
Title : Crystal structure of 3-oxoacyl-(acyl-carrier-protein) reductase (FabG) from *Pseudomonas aeruginosa* in complex with 1-(1-ethylbenzimidazol-2-yl)-3-(2-methoxyphenyl)urea at 1.9Å resolution
Authors : Cukier, C.D.; Schnell, R.; Lindqvist, Y.; Schneider, G.
Deposited on : 2013-05-18
Resolution : 1.90 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

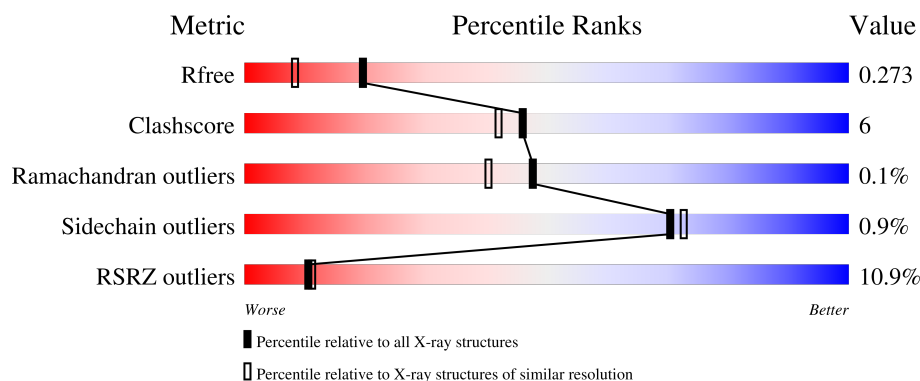
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	269	11% 79% 9% 12%
1	B	269	13% 79% 8% 13%
1	C	269	7% 83% 7% 11%
1	D	269	7% 86% 5% 9%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	36K	A	1248[A]	-	-	X	-
2	36K	A	1248[B]	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7250 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 3-OXOACYL-[ACYL-CARRIER-PROTEIN] REDUCTASE FABG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	238	Total	C	N	O	S	0	3	0
			1740	1089	308	337	6			
1	B	235	Total	C	N	O	S	0	2	0
			1704	1066	304	329	5			
1	C	240	Total	C	N	O	S	0	1	0
			1745	1092	310	337	6			
1	D	245	Total	C	N	O	S	0	1	0
			1789	1117	319	346	7			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	expression tag	UNP O54438
A	-20	HIS	-	expression tag	UNP O54438
A	-19	HIS	-	expression tag	UNP O54438
A	-18	HIS	-	expression tag	UNP O54438
A	-17	HIS	-	expression tag	UNP O54438
A	-16	HIS	-	expression tag	UNP O54438
A	-15	HIS	-	expression tag	UNP O54438
A	-14	SER	-	expression tag	UNP O54438
A	-13	SER	-	expression tag	UNP O54438
A	-12	GLY	-	expression tag	UNP O54438
A	-11	VAL	-	expression tag	UNP O54438
A	-10	ASP	-	expression tag	UNP O54438
A	-9	LEU	-	expression tag	UNP O54438
A	-8	GLY	-	expression tag	UNP O54438
A	-7	THR	-	expression tag	UNP O54438
A	-6	GLU	-	expression tag	UNP O54438
A	-5	ASN	-	expression tag	UNP O54438
A	-4	LEU	-	expression tag	UNP O54438
A	-3	TYR	-	expression tag	UNP O54438
A	-2	PHE	-	expression tag	UNP O54438

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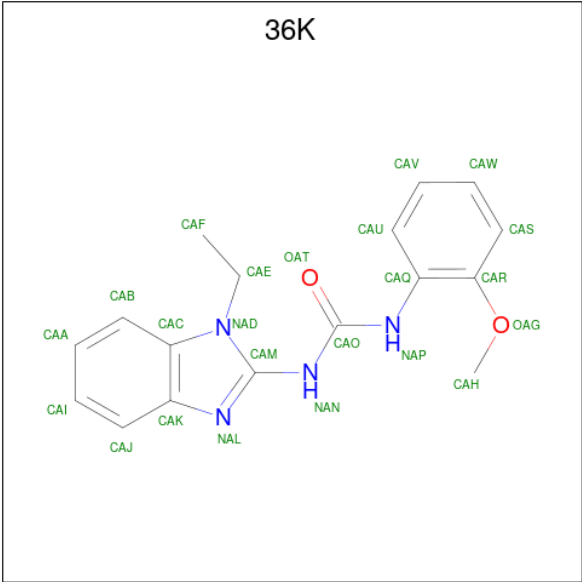
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLN	-	expression tag	UNP O54438
A	0	SER	-	expression tag	UNP O54438
B	-21	MET	-	expression tag	UNP O54438
B	-20	HIS	-	expression tag	UNP O54438
B	-19	HIS	-	expression tag	UNP O54438
B	-18	HIS	-	expression tag	UNP O54438
B	-17	HIS	-	expression tag	UNP O54438
B	-16	HIS	-	expression tag	UNP O54438
B	-15	HIS	-	expression tag	UNP O54438
B	-14	SER	-	expression tag	UNP O54438
B	-13	SER	-	expression tag	UNP O54438
B	-12	GLY	-	expression tag	UNP O54438
B	-11	VAL	-	expression tag	UNP O54438
B	-10	ASP	-	expression tag	UNP O54438
B	-9	LEU	-	expression tag	UNP O54438
B	-8	GLY	-	expression tag	UNP O54438
B	-7	THR	-	expression tag	UNP O54438
B	-6	GLU	-	expression tag	UNP O54438
B	-5	ASN	-	expression tag	UNP O54438
B	-4	LEU	-	expression tag	UNP O54438
B	-3	TYR	-	expression tag	UNP O54438
B	-2	PHE	-	expression tag	UNP O54438
B	-1	GLN	-	expression tag	UNP O54438
B	0	SER	-	expression tag	UNP O54438
C	-21	MET	-	expression tag	UNP O54438
C	-20	HIS	-	expression tag	UNP O54438
C	-19	HIS	-	expression tag	UNP O54438
C	-18	HIS	-	expression tag	UNP O54438
C	-17	HIS	-	expression tag	UNP O54438
C	-16	HIS	-	expression tag	UNP O54438
C	-15	HIS	-	expression tag	UNP O54438
C	-14	SER	-	expression tag	UNP O54438
C	-13	SER	-	expression tag	UNP O54438
C	-12	GLY	-	expression tag	UNP O54438
C	-11	VAL	-	expression tag	UNP O54438
C	-10	ASP	-	expression tag	UNP O54438
C	-9	LEU	-	expression tag	UNP O54438
C	-8	GLY	-	expression tag	UNP O54438
C	-7	THR	-	expression tag	UNP O54438
C	-6	GLU	-	expression tag	UNP O54438
C	-5	ASN	-	expression tag	UNP O54438
C	-4	LEU	-	expression tag	UNP O54438

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	TYR	-	expression tag	UNP O54438
C	-2	PHE	-	expression tag	UNP O54438
C	-1	GLN	-	expression tag	UNP O54438
C	0	SER	-	expression tag	UNP O54438
D	-21	MET	-	expression tag	UNP O54438
D	-20	HIS	-	expression tag	UNP O54438
D	-19	HIS	-	expression tag	UNP O54438
D	-18	HIS	-	expression tag	UNP O54438
D	-17	HIS	-	expression tag	UNP O54438
D	-16	HIS	-	expression tag	UNP O54438
D	-15	HIS	-	expression tag	UNP O54438
D	-14	SER	-	expression tag	UNP O54438
D	-13	SER	-	expression tag	UNP O54438
D	-12	GLY	-	expression tag	UNP O54438
D	-11	VAL	-	expression tag	UNP O54438
D	-10	ASP	-	expression tag	UNP O54438
D	-9	LEU	-	expression tag	UNP O54438
D	-8	GLY	-	expression tag	UNP O54438
D	-7	THR	-	expression tag	UNP O54438
D	-6	GLU	-	expression tag	UNP O54438
D	-5	ASN	-	expression tag	UNP O54438
D	-4	LEU	-	expression tag	UNP O54438
D	-3	TYR	-	expression tag	UNP O54438
D	-2	PHE	-	expression tag	UNP O54438
D	-1	GLN	-	expression tag	UNP O54438
D	0	SER	-	expression tag	UNP O54438

- Molecule 2 is 1-(1-ethylbenzimidazol-2-yl)-3-(2-methoxyphenyl)urea (CCD ID: 36K) (formula: C₁₇H₁₈N₄O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	1
			46	34	8	4		
2	C	1	Total	C	N	O	0	0
			23	17	4	2		

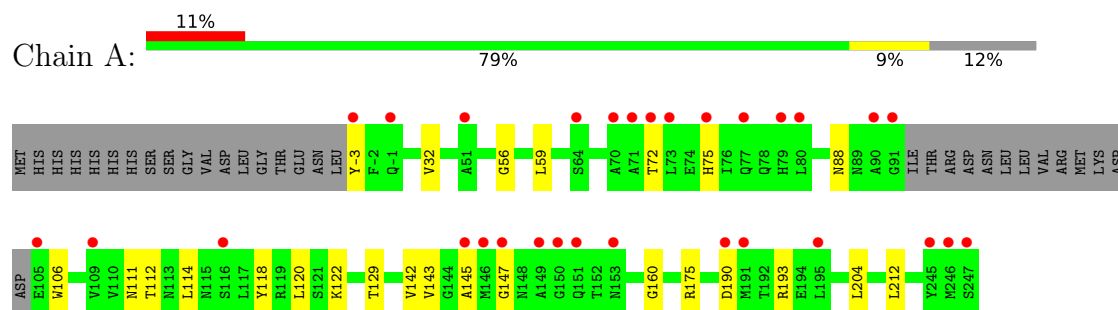
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	55	Total	O	0	0
			55	55		
3	B	48	Total	O	0	0
			48	48		
3	C	43	Total	O	0	0
			43	43		
3	D	57	Total	O	0	0
			57	57		

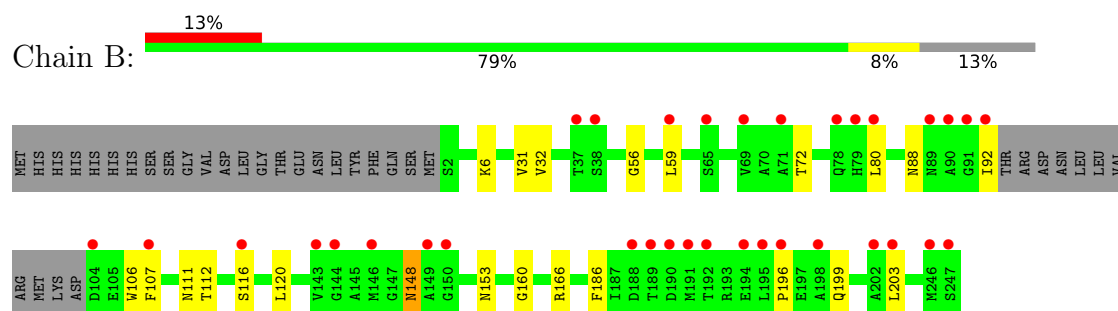
3 Residue-property plots [i](#)

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.

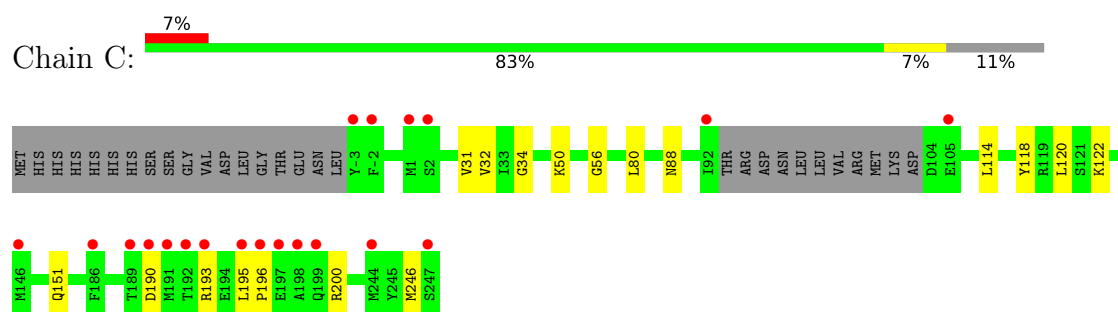
• Molecule 1: 3-OXOACYL-[ACYL-CARRIER-PROTEIN] REDUCTASE FABG



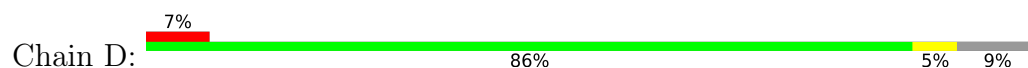
• Molecule 1: 3-OXOACYL-[ACYL-CARRIER-PROTEIN] REDUCTASE FABG

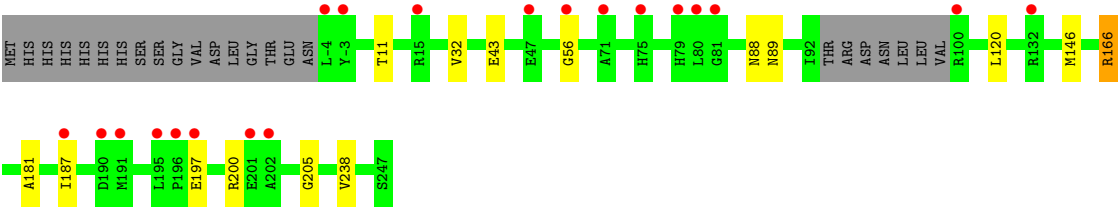


• Molecule 1: 3-OXOACYL-[ACYL-CARRIER-PROTEIN] REDUCTASE FABG



• Molecule 1: 3-OXOACYL-[ACYL-CARRIER-PROTEIN] REDUCTASE FABG





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.55Å 108.92Å 149.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.56 – 1.90 38.56 – 1.90	Depositor EDS
% Data completeness (in resolution range)	87.3 (38.56-1.90) 87.3 (38.56-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.235 , 0.269 0.240 , 0.273	Depositor DCC
R_{free} test set	3401 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	19.1	Xtriage
Anisotropy	0.415	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 37.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7250	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.35% 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: 36K

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	1.10	0/1768	1.01	0/2392
1	B	1.17	0/1727	1.03	3/2338 (0.1%)
1	C	1.13	1/1767 (0.1%)	1.03	0/2391
1	D	1.15	3/1811 (0.2%)	1.05	1/2448 (0.0%)
All	All	1.14	4/7073 (0.1%)	1.03	4/9569 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	181	ALA	N-CA	-5.52	1.38	1.45
1	D	187	ILE	CA-C	5.40	1.59	1.52
1	C	34	GLY	CA-C	-5.20	1.48	1.52
1	D	238	VAL	N-CA	5.03	1.50	1.46

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	146	MET	N-CA-C	9.52	122.86	111.82
1	B	148	ASN	N-CA-C	6.88	118.43	111.07
1	B	6	LYS	CB-CG-CD	5.58	124.14	111.30
1	B	6	LYS	CA-CB-CG	-5.12	103.85	114.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	147	GLY	Peptide

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	1740	0	1775	21	0
1	B	1704	0	1748	19	0
1	C	1745	0	1780	12	0
1	D	1789	0	1823	6	0
2	A	46	0	36	31	0
2	C	23	0	18	5	0
3	A	55	0	0	3	0
3	B	48	0	0	1	0
3	C	43	0	0	0	0
3	D	57	0	0	0	0
All	All	7250	0	7180	82	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 (82) 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:A:1248[B]:36K:HAP	2:A:1248[B]:36K:CAH	1.29	1.43
2:A:1248[B]:36K:CAF	2:A:1248[B]:36K:HAB	1.64	1.27
2:A:1248[B]:36K:HAN	2:A:1248[B]:36K:CAU	1.34	1.21
2:A:1248[A]:36K:HAN	2:A:1248[A]:36K:CAU	1.43	1.14
1:A:59[B]:LEU:HD21	1:A:72:THR:HG23	1.24	1.10
2:A:1248[B]:36K:HAU	2:A:1248[B]:36K:NAN	1.37	1.08
1:B:59[B]:LEU:HD21	1:B:72:THR:CG2	1.83	1.06
2:A:1248[B]:36K:HAB	2:A:1248[B]:36K:HAF3	1.36	1.06
2:A:1248[B]:36K:HAB	2:A:1248[B]:36K:HAF2	1.35	1.06
2:A:1248[B]:36K:HAP	2:A:1248[B]:36K:HAH2	0.92	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1248[B]:36K:HAH2	2:A:1248[B]:36K:NAP	1.58	1.03
1:B:59[B]:LEU:CD2	1:B:72:THR:HG23	1.88	1.02
2:A:1248[A]:36K:HAU	2:A:1248[A]:36K:NAN	1.51	1.02
2:A:1248[B]:36K:HAP	2:A:1248[B]:36K:HAH3	1.25	1.00
2:A:1248[A]:36K:HAF2	2:A:1248[A]:36K:OAT	1.62	1.00
2:A:1248[B]:36K:CAH	2:A:1248[B]:36K:NAP	2.00	0.99
1:A:59[B]:LEU:CD2	1:A:72:THR:HG23	1.92	0.99
2:A:1248[B]:36K:CAF	2:A:1248[B]:36K:CAB	2.35	0.90
1:B:59[B]:LEU:HD21	1:B:72:THR:HG23	0.93	0.88
2:A:1248[A]:36K:HAN	2:A:1248[A]:36K:HAU	0.66	0.83
2:A:1248[B]:36K:HAF2	2:A:1248[B]:36K:CAB	2.05	0.81
2:A:1248[A]:36K:OAT	2:A:1248[A]:36K:CAF	2.30	0.80
2:A:1248[B]:36K:NAP	2:A:1248[B]:36K:HAH3	1.87	0.78
2:A:1248[B]:36K:HAF3	2:A:1248[B]:36K:CAB	2.06	0.77
1:D:197:GLU:OE2	1:D:200:ARG:NH1	2.19	0.76
2:A:1248[B]:36K:HAN	2:A:1248[B]:36K:HAU	0.59	0.75
2:A:1248[A]:36K:OAT	2:A:1248[A]:36K:CAE	2.43	0.67
2:A:1248[A]:36K:HAH1	1:B:107:PHE:CE1	2.34	0.63
2:A:1248[B]:36K:CAU	2:A:1248[B]:36K:NAN	2.08	0.62
1:A:160:GLY:HA2	2:A:1248[B]:36K:HAE2	1.82	0.62
1:C:193:ARG:O	1:C:200:ARG:NH2	2.33	0.61
1:C:151:GLN:HG2	1:C:246:MET:HE3	1.82	0.60
1:A:-3:TYR:CD2	3:A:2001:HOH:O	2.52	0.60
1:A:204:LEU:HD21	1:A:212:LEU:HD12	1.84	0.59
1:B:92:ILE:HG13	1:B:112:THR:HG22	1.82	0.59
2:A:1248[B]:36K:OAT	2:A:1248[B]:36K:NAL	2.26	0.58
1:A:118:TYR:OH	1:A:122:LYS:HE3	2.05	0.57
1:C:195:LEU:HG	1:C:196:PRO:HD2	1.85	0.57
1:B:148:ASN:ND2	1:D:205:GLY:O	2.26	0.55
2:A:1248[B]:36K:HAE1	1:B:160:GLY:HA2	1.90	0.54
1:A:114:LEU:HD13	2:A:1248[A]:36K:HAH2	1.90	0.52
1:A:112:THR:HG23	3:A:2018:HOH:O	2.09	0.52
1:A:190:ASP:OD1	1:A:190:ASP:N	2.43	0.52
1:A:112:THR:HA	3:A:2018:HOH:O	2.10	0.52
1:A:204:LEU:HD21	1:A:212:LEU:CD1	2.40	0.52
1:B:106:TRP:HB3	1:B:153:ASN:HD22	1.74	0.51
1:C:118:TYR:OH	1:C:122:LYS:HE2	2.10	0.51
1:C:190:ASP:N	1:C:190:ASP:OD1	2.44	0.50
1:C:114:LEU:HD12	2:C:1248:36K:NAN	2.27	0.49
1:A:32:VAL:O	1:A:56:GLY:HA3	2.13	0.49
1:D:32:VAL:O	1:D:56:GLY:HA3	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1248[B]:36K:HAH1	1:B:111:ASN:HB2	1.96	0.48
1:B:31:VAL:HG11	1:B:80:LEU:HD22	1.96	0.48
1:B:196:PRO:HD2	1:B:199:GLN:HG3	1.94	0.48
1:A:106:TRP:CZ3	2:A:1248[B]:36K:HAI	2.50	0.47
1:C:32:VAL:O	1:C:56:GLY:HA3	2.14	0.47
1:A:160:GLY:CA	2:A:1248[B]:36K:HAE2	2.43	0.46
1:A:88:ASN:OD1	1:A:120:LEU:HD23	2.15	0.46
1:B:116:SER:HB2	3:B:2015:HOH:O	2.15	0.46
1:A:111:ASN:OD1	1:B:111:ASN:ND2	2.48	0.46
1:B:59[B]:LEU:CD2	1:B:72:THR:CG2	2.69	0.46
1:C:88:ASN:OD1	1:C:120:LEU:HD23	2.16	0.45
1:B:92:ILE:HG13	1:B:112:THR:CG2	2.47	0.45
2:C:1248:36K:NAL	2:C:1248:36K:NAP	2.61	0.45
1:B:88:ASN:OD1	1:B:120:LEU:HD23	2.17	0.44
1:D:88:ASN:OD1	1:D:120:LEU:HD23	2.18	0.43
1:B:32:VAL:O	1:B:56:GLY:HA3	2.19	0.43
1:B:186:PHE:CD2	1:B:203:LEU:HD11	2.53	0.43
1:D:11:THR:O	1:D:89:ASN:HB3	2.19	0.43
1:A:142:VAL:HG12	1:A:143:VAL:O	2.19	0.43
2:C:1248:36K:OAT	2:C:1248:36K:HAU	2.18	0.42
1:A:59[B]:LEU:HD23	1:A:72:THR:HG23	1.92	0.42
1:A:118:TYR:CZ	1:A:122:LYS:HE3	2.54	0.42
1:C:31:VAL:HG11	1:C:80:LEU:HD22	2.01	0.42
1:B:106:TRP:HB3	1:B:153:ASN:ND2	2.34	0.42
1:C:114:LEU:HD12	2:C:1248:36K:CAO	2.50	0.41
2:A:1248[A]:36K:OAT	2:A:1248[A]:36K:HAE1	2.19	0.41
1:C:114:LEU:CD1	2:C:1248:36K:CAM	2.99	0.41
1:A:129:THR:HG23	1:A:175:ARG:HD3	2.02	0.41
1:A:122:LYS:HA	1:A:122:LYS:HD3	1.76	0.41
1:D:166:ARG:HH11	1:D:166:ARG:HD2	1.73	0.40
1:C:118:TYR:CZ	1:C:122:LYS:HE2	2.56	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	237/269 (88%)	232 (98%)	4 (2%)	1 (0%)	30	22
1	B	233/269 (87%)	230 (99%)	3 (1%)	0	100	100
1	C	237/269 (88%)	235 (99%)	2 (1%)	0	100	100
1	D	242/269 (90%)	239 (99%)	3 (1%)	0	100	100
All	All	949/1076 (88%)	936 (99%)	12 (1%)	1 (0%)	48	40

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	145	ALA

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	175/201 (87%)	173 (99%)	2 (1%)	65	67
1	B	171/201 (85%)	170 (99%)	1 (1%)	78	81
1	C	175/201 (87%)	174 (99%)	1 (1%)	78	81
1	D	180/201 (90%)	178 (99%)	2 (1%)	65	67
All	All	701/804 (87%)	695 (99%)	6 (1%)	70	73

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	HIS
1	A	193	ARG
1	B	166	ARG
1	C	50	LYS
1	D	43	GLU
1	D	166	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	GLN
1	A	77	GLN
1	A	82	GLN
1	A	89	ASN
1	A	151	GLN
1	A	153	ASN
1	A	199	GLN
1	B	19	GLN
1	B	77	GLN
1	B	78	GLN
1	B	153	ASN
1	B	199	GLN
1	C	77	GLN
1	C	82	GLN
1	C	111	ASN
1	C	115	ASN
1	C	148	ASN
1	C	214	GLN
1	D	19	GLN
1	D	111	ASN
1	D	115	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	36K	A	1248[B]	-	24,25,25	2.39	8 (33%)	30,34,34	1.65	5 (16%)
2	36K	A	1248[A]	-	24,25,25	2.55	7 (29%)	30,34,34	2.24	9 (30%)
2	36K	C	1248	-	24,25,25	1.90	7 (29%)	30,34,34	1.59	6 (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
2	36K	A	1248[B]	-	-	6/12/12/12	0/3/3/3
2	36K	A	1248[A]	-	-	5/12/12/12	0/3/3/3
2	36K	C	1248	-	-	0/12/12/12	0/3/3/3

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1248[B]	36K	CAJ-CAK	-5.92	1.30	1.40
2	A	1248[A]	36K	CAB-CAC	-5.81	1.30	1.39
2	A	1248[A]	36K	CAJ-CAK	-5.71	1.31	1.40
2	C	1248	36K	CAJ-CAK	-5.31	1.31	1.40
2	A	1248[A]	36K	CAK-NAL	-5.07	1.29	1.39
2	A	1248[B]	36K	CAB-CAC	-4.75	1.32	1.39
2	A	1248[B]	36K	CAC-NAD	-4.69	1.30	1.39
2	A	1248[A]	36K	CAQ-NAP	-4.00	1.33	1.41
2	A	1248[B]	36K	CAC-CAK	-3.82	1.34	1.40
2	A	1248[B]	36K	CAK-NAL	-3.81	1.32	1.39
2	A	1248[A]	36K	CAC-CAK	-3.78	1.34	1.40
2	A	1248[B]	36K	CAQ-NAP	-3.65	1.34	1.41
2	C	1248	36K	CAQ-NAP	-3.61	1.34	1.41
2	A	1248[A]	36K	CAC-NAD	-3.24	1.33	1.39
2	C	1248	36K	CAC-CAK	-3.05	1.35	1.40
2	C	1248	36K	CAI-CAJ	2.55	1.43	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1248	36K	CAO-NAP	2.48	1.42	1.37
2	A	1248[A]	36K	CAI-CAJ	-2.24	1.35	1.38
2	C	1248	36K	CAO-NAN	-2.21	1.34	1.39
2	C	1248	36K	CAM-NAL	2.15	1.35	1.32
2	A	1248[B]	36K	CAO-NAN	-2.10	1.34	1.39
2	A	1248[B]	36K	CAO-NAP	-2.07	1.32	1.37

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1248[A]	36K	CAH-OAG-CAR	-6.82	107.51	117.51
2	A	1248[A]	36K	CAK-CAC-NAD	5.68	110.04	105.55
2	A	1248[B]	36K	CAM-NAN-CAO	-5.12	115.94	125.72
2	A	1248[B]	36K	CAQ-NAP-CAO	-4.43	115.01	125.34
2	C	1248	36K	CAE-NAD-CAC	-3.96	116.17	124.49
2	C	1248	36K	CAH-OAG-CAR	3.62	122.83	117.51
2	A	1248[A]	36K	CAJ-CAK-CAC	3.60	124.27	120.24
2	A	1248[B]	36K	CAH-OAG-CAR	-3.60	112.24	117.51
2	A	1248[A]	36K	CAQ-NAP-CAO	-3.12	118.09	125.34
2	A	1248[A]	36K	CAA-CAI-CAJ	-2.93	116.62	120.24
2	C	1248	36K	CAI-CAA-CAB	-2.91	116.64	120.24
2	A	1248[A]	36K	CAB-CAC-CAK	-2.63	119.20	122.23
2	C	1248	36K	CAB-CAC-CAK	2.58	125.20	122.23
2	C	1248	36K	NAP-CAO-NAN	2.44	119.51	114.22
2	A	1248[A]	36K	NAN-CAM-NAL	-2.42	118.32	127.16
2	A	1248[B]	36K	CAE-NAD-CAC	-2.39	119.48	124.49
2	C	1248	36K	OAT-CAO-NAN	-2.37	115.07	120.69
2	A	1248[A]	36K	CAE-NAD-CAC	-2.33	119.59	124.49
2	A	1248[A]	36K	CAM-NAN-CAO	-2.08	121.74	125.72
2	A	1248[B]	36K	CAK-CAC-NAD	2.02	107.15	105.55

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1248[A]	36K	NAD-CAM-NAN-CAO
2	A	1248[A]	36K	CAQ-CAR-OAG-CAH
2	A	1248[B]	36K	CAF-CAE-NAD-CAC
2	A	1248[B]	36K	CAF-CAE-NAD-CAM
2	A	1248[B]	36K	CAQ-CAR-OAG-CAH
2	A	1248[A]	36K	CAS-CAR-OAG-CAH
2	A	1248[B]	36K	CAS-CAR-OAG-CAH

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Mol	Chain	Res	Type	Atoms
2	A	1248[A]	36K	OAT-CAO-NAP-CAQ
2	A	1248[B]	36K	OAT-CAO-NAP-CAQ
2	A	1248[A]	36K	NAN-CAO-NAP-CAQ
2	A	1248[B]	36K	NAN-CAO-NAP-CAQ

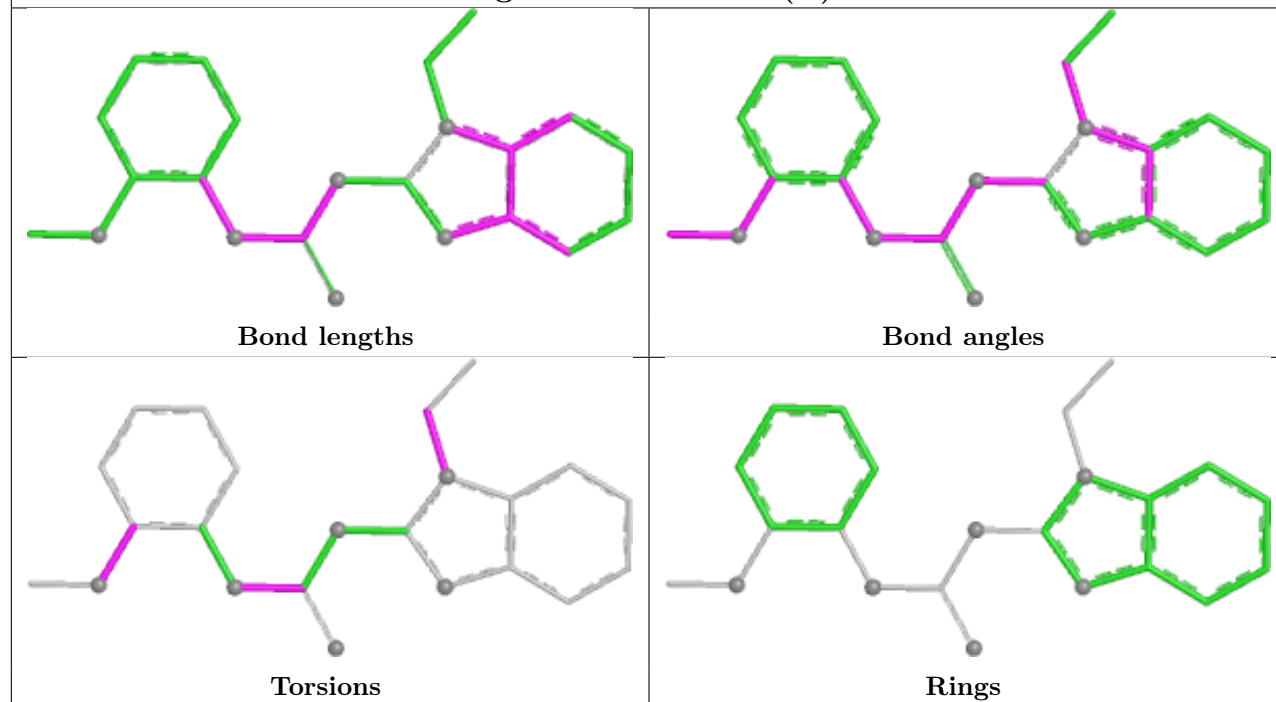
There are no ring outliers.

3 monomers are involved in 36 short contacts:

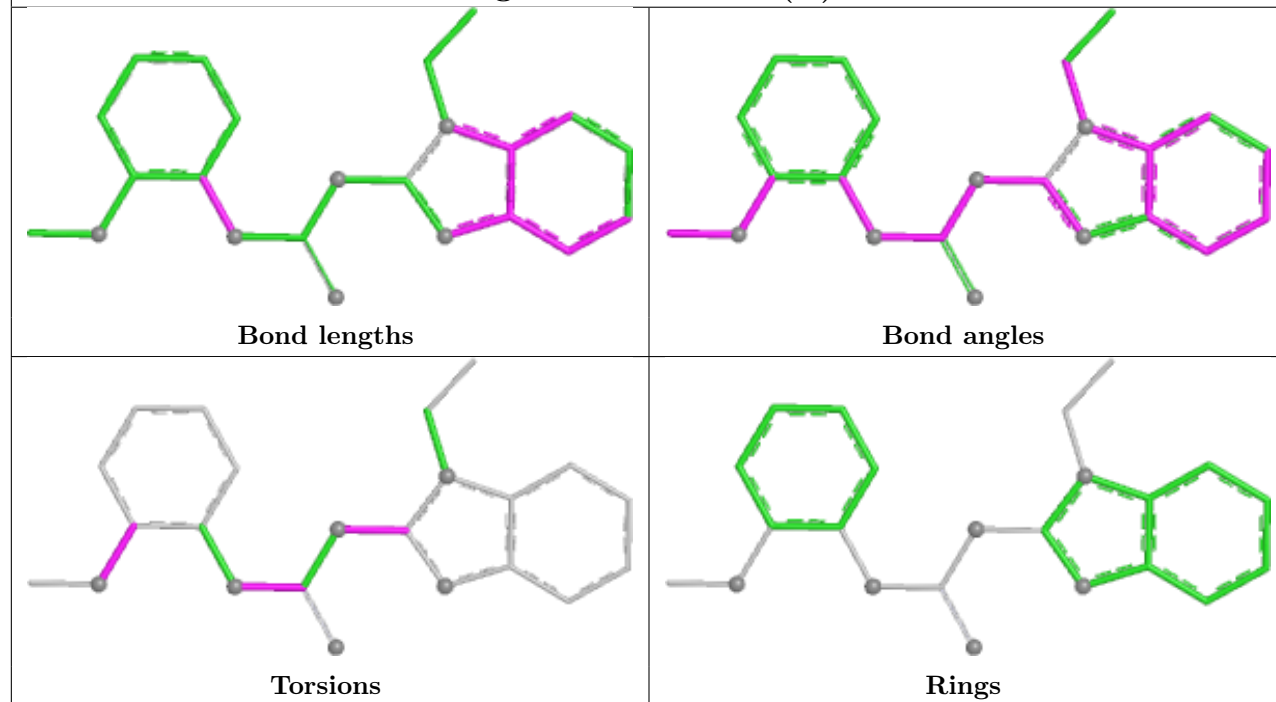
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1248[B]	36K	22	0
2	A	1248[A]	36K	9	0
2	C	1248	36K	5	0

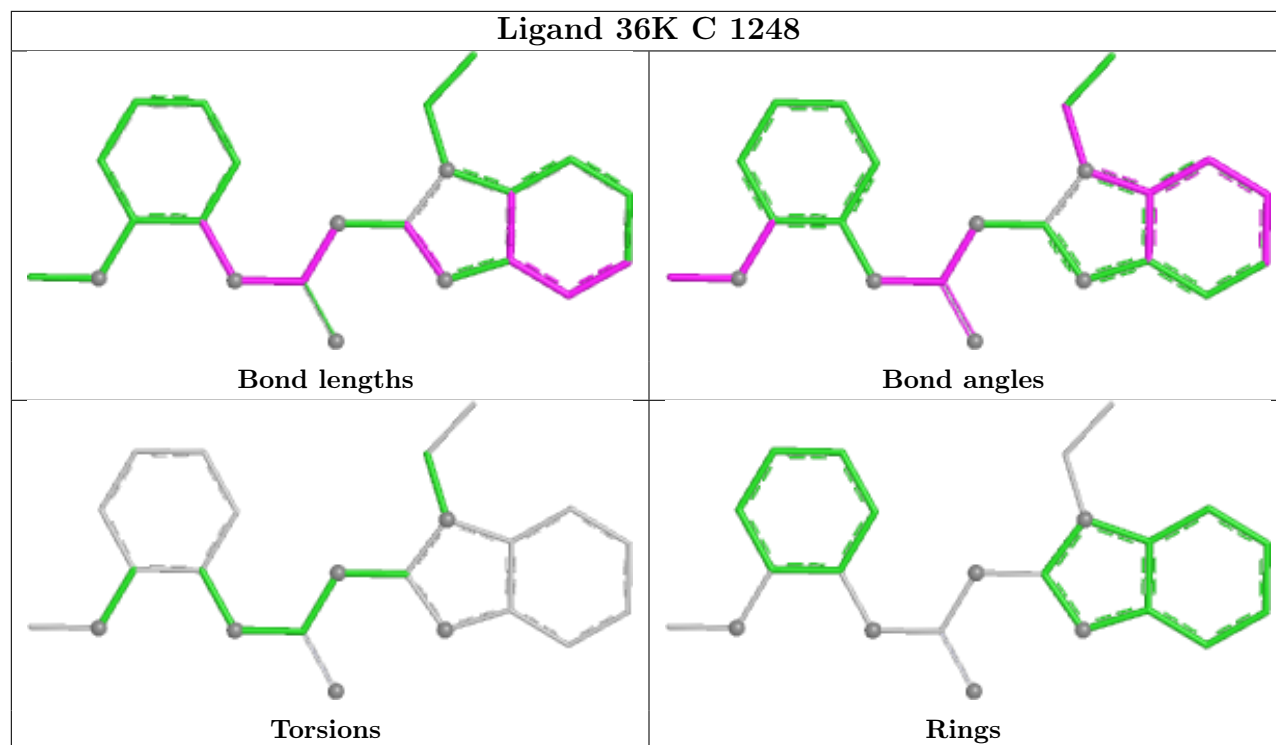
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 36K A 1248 (B)



Ligand 36K A 1248 (A)





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	238/269 (88%)	0.88	30 (12%) 8 8	16, 36, 69, 91	3 (1%)
1	B	235/269 (87%)	0.85	34 (14%) 6 6	20, 34, 72, 105	2 (0%)
1	C	240/269 (89%)	0.72	20 (8%) 17 18	18, 32, 78, 114	1 (0%)
1	D	245/269 (91%)	0.79	20 (8%) 17 18	19, 34, 66, 92	1 (0%)
All	All	958/1076 (89%)	0.81	104 (10%) 10 11	16, 34, 71, 114	7 (0%)

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	92	ILE	7.9
1	B	92	ILE	6.9
1	C	195	LEU	6.0
1	D	-4	LEU	5.3
1	A	149	ALA	5.2
1	B	90	ALA	4.7
1	C	-3	TYR	4.5
1	C	192	THR	4.5
1	A	80	LEU	4.5
1	C	196	PRO	4.4
1	C	191	MET	4.2
1	A	150	GLY	4.1
1	B	191	MET	4.1
1	C	189	THR	4.0
1	A	-3	TYR	4.0
1	B	192	THR	3.7
1	B	80	LEU	3.6
1	B	189	THR	3.6
1	D	187	ILE	3.6
1	A	191	MET	3.5
1	A	70	ALA	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	79	HIS	3.5
1	A	146	MET	3.4
1	B	146	MET	3.3
1	C	198	ALA	3.3
1	A	91	GLY	3.2
1	D	197	GLU	3.2
1	C	190	ASP	3.2
1	D	190	ASP	3.2
1	B	37	THR	3.1
1	B	143	VAL	3.1
1	B	91	GLY	3.1
1	C	105	GLU	3.1
1	B	195	LEU	3.1
1	D	-3	TYR	3.1
1	B	71	ALA	3.0
1	A	246	MET	3.0
1	A	245	TYR	3.0
1	B	190	ASP	3.0
1	D	132	ARG	2.9
1	B	149	ALA	2.9
1	A	151	GLN	2.9
1	D	191	MET	2.9
1	B	59[A]	LEU	2.8
1	A	64	SER	2.8
1	B	65	SER	2.8
1	A	153	ASN	2.8
1	D	47	GLU	2.8
1	C	193	ARG	2.7
1	D	71	ALA	2.7
1	A	75	HIS	2.7
1	B	144	GLY	2.7
1	A	195	LEU	2.7
1	B	194	GLU	2.6
1	B	188	ASP	2.6
1	C	-2	PHE	2.6
1	A	116	SER	2.6
1	A	77	GLN	2.5
1	D	56	GLY	2.5
1	B	202	ALA	2.5
1	D	202	ALA	2.5
1	A	79	HIS	2.5
1	A	73	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	80	LEU	2.5
1	C	146	MET	2.5
1	D	79	HIS	2.5
1	A	71	ALA	2.5
1	C	197	GLU	2.4
1	D	196	PRO	2.4
1	A	247	SER	2.4
1	C	244	MET	2.4
1	C	1	MET	2.4
1	C	199	GLN	2.3
1	C	247	SER	2.3
1	D	201	GLU	2.3
1	A	-1	GLN	2.3
1	A	109	VAL	2.3
1	A	72	THR	2.3
1	B	107	PHE	2.2
1	B	196	PRO	2.2
1	B	198	ALA	2.2
1	B	104	ASP	2.2
1	B	69	VAL	2.2
1	D	75	HIS	2.2
1	B	247	SER	2.2
1	B	116	SER	2.2
1	B	203	LEU	2.2
1	D	195	LEU	2.2
1	A	145	ALA	2.2
1	B	246	MET	2.1
1	A	105	GLU	2.1
1	B	150	GLY	2.1
1	D	15	ARG	2.1
1	A	51	ALA	2.1
1	A	190	ASP	2.1
1	C	2	SER	2.1
1	A	147	GLY	2.1
1	D	81	GLY	2.1
1	C	186	PHE	2.0
1	B	38	SER	2.0
1	B	89	ASN	2.0
1	A	90	ALA	2.0
1	B	78	GLN	2.0
1	D	100	ARG	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](#)

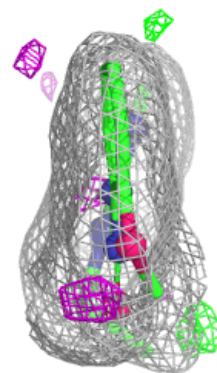
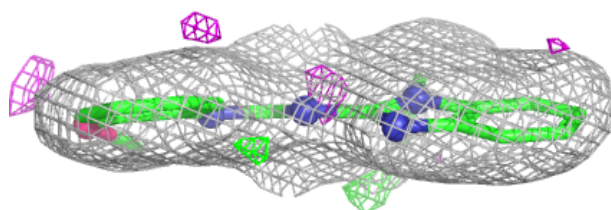
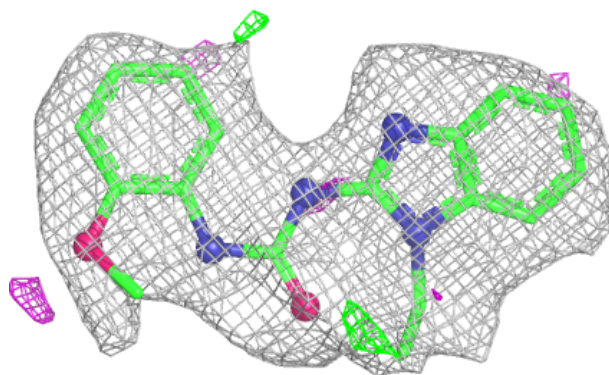
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	36K	A	1248[A]	23/23	0.86	0.12	22,34,44,44	23
2	36K	A	1248[B]	23/23	0.86	0.12	37,42,50,53	23
2	36K	C	1248	23/23	0.95	0.07	19,26,31,36	0

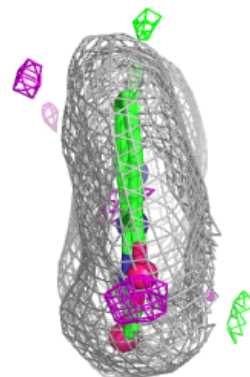
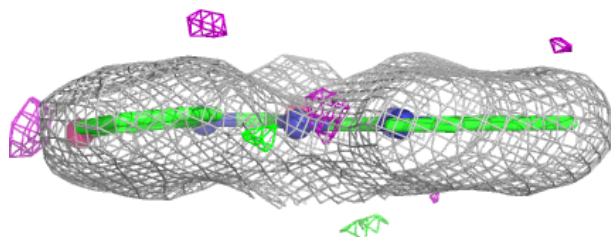
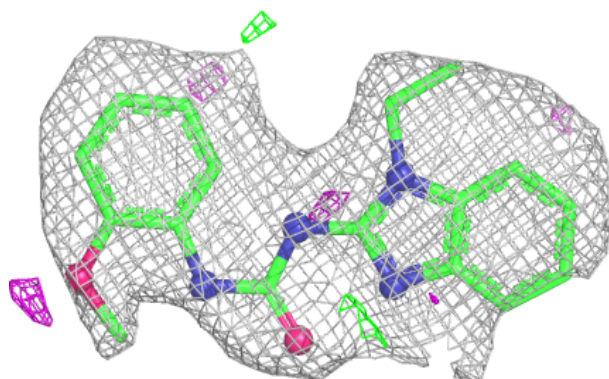
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

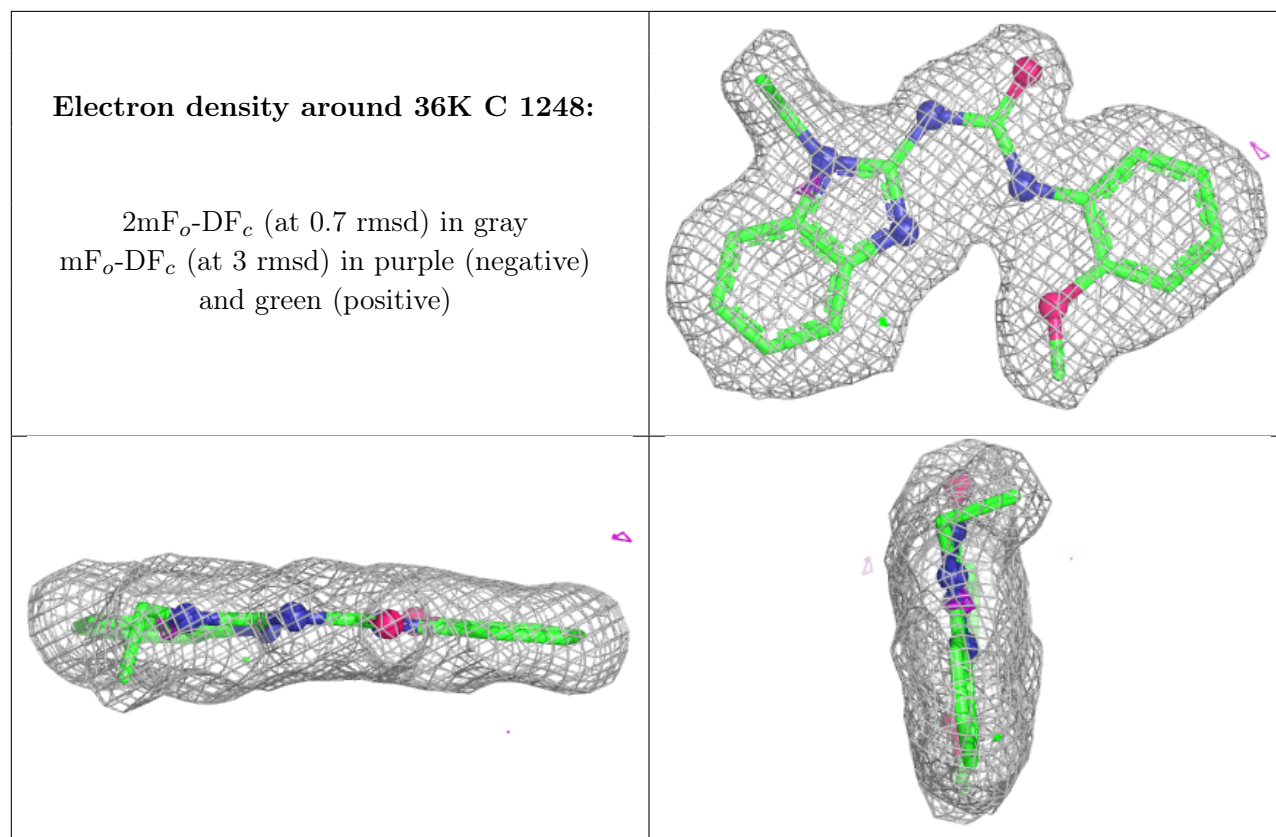
Electron density around 36K A 1248 (A):

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

**Electron density around 36K A 1248 (B):**

$2mF_o-DF_c$ (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.