



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:33 AM UTC

PDB ID : 4BNX / pdb_00004bnx
Title : Crystal structure of 3-oxoacyl-(acyl-carrier-protein) reductase (FabG) from *Pseudomonas aeruginosa* in complex with 6-(4-(2-chloroanilino)-1H-quinazolin-2-ylidene)cyclohexa-2,4-dien-1-one at 2.3Å resolution
Authors : Cukier, C.D.; Schnell, R.; Lindqvist, Y.; Schneider, G.
Deposited on : 2013-05-17
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)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

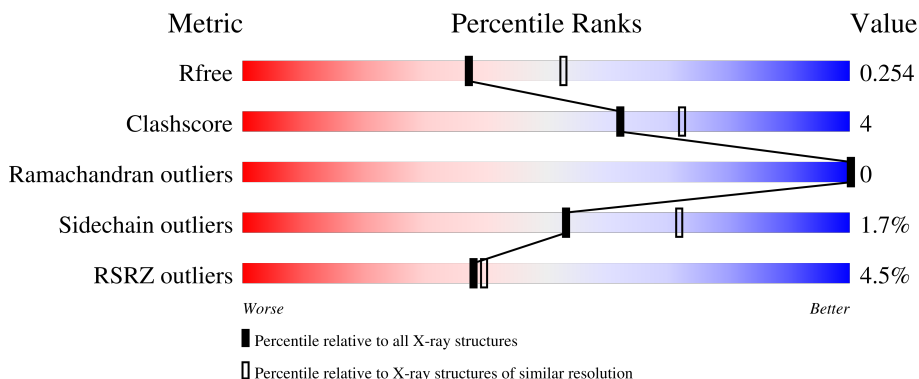
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	269	4% 81% 7% 11%
1	B	269	5% 81% 9% 10%
1	C	269	3% 79% 11% 10%
1	D	269	4% 81% 8% 11%

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	O74	A	1247	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7192 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	239	Total	C	N	O	S	0	4	0
			1755	1094	315	339	7			
1	B	241	Total	C	N	O	S	0	3	0
			1771	1110	313	341	7			
1	C	242	Total	C	N	O	S	0	2	0
			1774	1110	315	343	6			
1	D	240	Total	C	N	O	S	0	3	0
			1753	1093	315	338	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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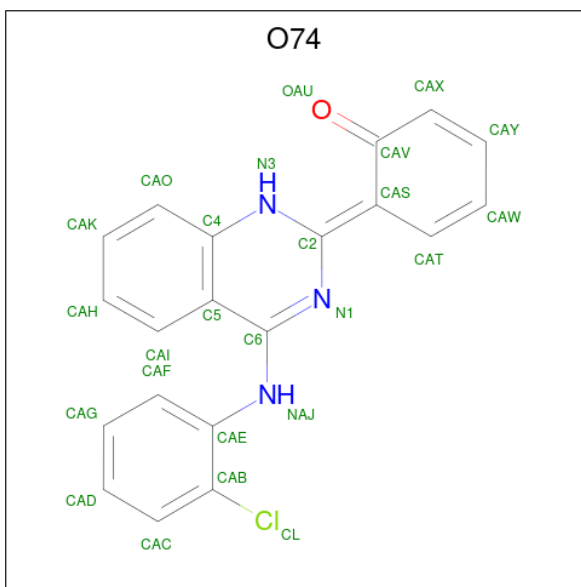
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLN	-	expression tag	UNP O54438
A	0	SER	-	expression tag	UNP O54438
B	-21	MET	-	expression tag	UNP O54438
B	-20	HIS	-	expression tag	UNP O54438
B	-19	HIS	-	expression tag	UNP O54438
B	-18	HIS	-	expression tag	UNP O54438
B	-17	HIS	-	expression tag	UNP 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 6-[4-(2-chloroanilino)-1H-quinazolin-2-ylidene]cyclohexa-2,4-dien-1-one (CCD ID: O74) (formula: C₂₀H₁₄ClN₃O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	0	0
			25	20	1	3	1		
2	C	1	Total	C	Cl	N	O	0	0
			25	20	1	3	1		

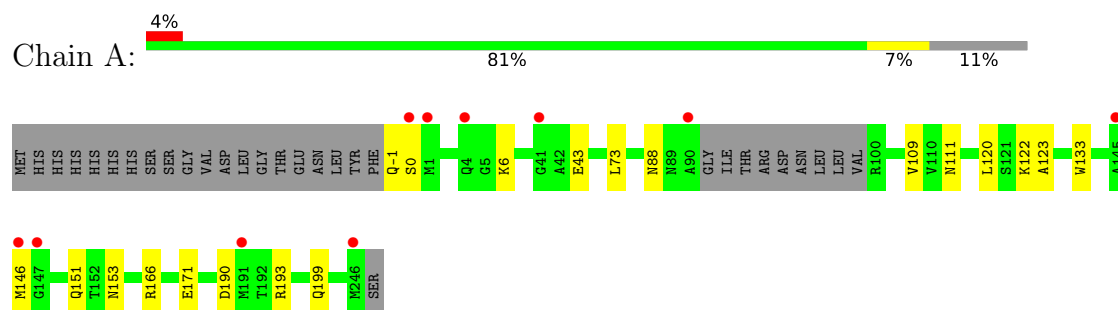
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	17	Total	O	0	0
			17	17		
3	B	31	Total	O	0	0
			31	31		
3	C	21	Total	O	0	0
			21	21		
3	D	20	Total	O	0	0
			20	20		

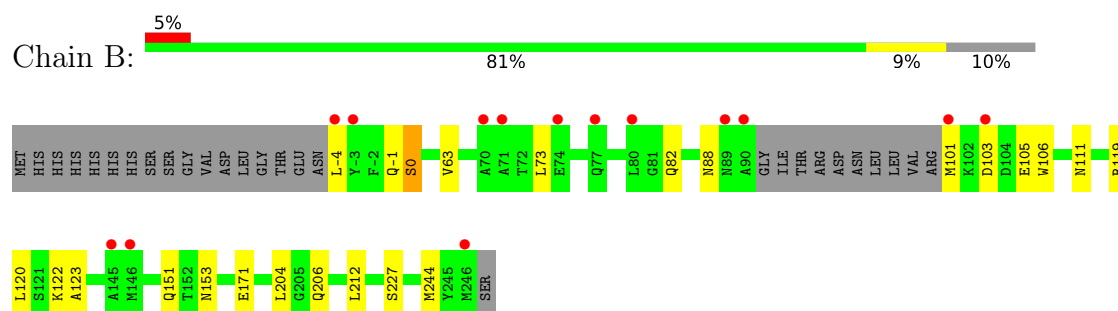
3 Residue-property plots

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.

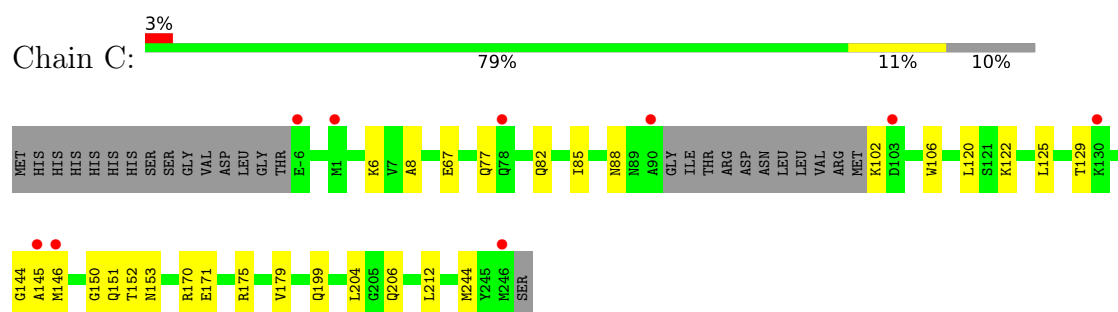
• Molecule 1: 3-OXOACYL-[ACYL-CARRIER-PROTEIN] REDUCTASE FABG



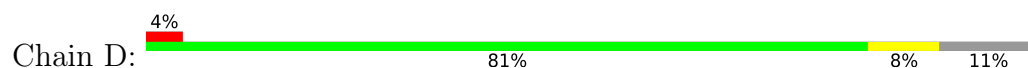
• Molecule 1: 3-OXOACYL-[ACYL-CARRIER-PROTEIN] REDUCTASE FABG



• 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	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	109.43Å 140.92Å 143.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.27 – 2.30 50.27 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.1 (50.27-2.30) 99.1 (50.27-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.213 , 0.255 0.215 , 0.254	Depositor DCC
R_{free} test set	2482 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	41.6	Xtriage
Anisotropy	0.468	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 28.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.011 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7192	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% 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: O74

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	0.90	0/1784	0.94	1/2413 (0.0%)
1	B	0.97	0/1799	0.98	1/2434 (0.0%)
1	C	1.03	1/1799 (0.1%)	1.01	1/2435 (0.0%)
1	D	0.95	1/1779 (0.1%)	0.96	0/2403
All	All	0.96	2/7161 (0.0%)	0.98	3/9685 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	179	VAL	C-O	7.44	1.32	1.24
1	D	90	ALA	CA-C	5.85	1.60	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	103	ASP	N-CA-CB	-6.07	100.23	110.49
1	A	166	ARG	CB-CA-C	-5.19	102.18	110.79
1	C	125	LEU	N-CA-C	5.01	117.63	111.82

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	1755	0	1800	14	0
1	B	1771	0	1812	14	0
1	C	1774	0	1809	15	0
1	D	1753	0	1803	12	0
2	A	25	0	14	9	0
2	C	25	0	14	3	0
3	A	17	0	0	0	0
3	B	31	0	0	0	0
3	C	21	0	0	2	0
3	D	20	0	0	1	0
All	All	7192	0	7252	59	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 (59) 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:111:ASN:HD22	2:A:1247:O74:HAH	1.28	0.97
1:A:111:ASN:HD22	2:A:1247:O74:CAH	1.90	0.84
2:A:1247:O74:N1	2:A:1247:O74:HAF	1.92	0.83
1:A:190:ASP:OD1	1:A:193:ARG:NH1	2.26	0.68
2:A:1247:O74:N1	2:A:1247:O74:CAF	2.50	0.68
1:D:187:ILE:HG23	1:D:213:GLY:O	1.98	0.64
1:B:122:LYS:NZ	1:B:171:GLU:OE2	2.31	0.64
1:C:152:THR:OG1	1:D:170:ARG:NH1	2.34	0.60
1:C:145:ALA:HB1	1:C:150:GLY:HA3	1.83	0.59
2:A:1247:O74:OAU	2:A:1247:O74:N3	2.22	0.59
2:C:1247:O74:HAF	2:C:1247:O74:N1	2.18	0.58
1:C:204:LEU:HD21	1:C:212:LEU:HD12	1.84	0.58
1:A:-1:GLN:HG3	1:A:0:SER:H	1.68	0.57
2:A:1247:O74:HAK	1:B:111:ASN:OD1	2.05	0.57
1:B:204:LEU:HD21	1:B:212:LEU:HD12	1.88	0.55
1:C:146:MET:SD	1:C:199:GLN:OE1	2.64	0.55
1:D:186:PHE:CZ	1:D:203:LEU:HD11	2.41	0.55
1:A:-1:GLN:HG3	1:A:0:SER:N	2.22	0.55
1:C:122:LYS:NZ	1:C:171:GLU:OE2	2.41	0.54
1:A:109:VAL:HG11	1:A:153:ASN:HD21	1.74	0.53
1:B:101:MET:HB3	1:B:105:GLU:OE1	2.09	0.52
2:A:1247:O74:HAF	2:A:1247:O74:HAT	1.92	0.52
1:A:122:LYS:NZ	1:A:171:GLU:OE2	2.42	0.51
1:A:111:ASN:HD22	2:A:1247:O74:CAK	2.25	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:0:SER:O	1:B:227:SER:HB2	2.12	0.49
1:A:111:ASN:ND2	2:A:1247:O74:HAH	2.12	0.48
2:C:1247:O74:HAK	1:D:111:ASN:OD1	2.14	0.48
1:B:206[A]:GLN:HB2	1:B:244:MET:HE3	1.96	0.47
1:C:144:GLY:O	1:C:145:ALA:C	2.57	0.47
1:C:170:ARG:HD3	3:C:2020:HOH:O	2.13	0.47
2:C:1247:O74:N1	2:C:1247:O74:CAF	2.76	0.47
1:D:170:ARG:HD2	3:D:2017:HOH:O	2.15	0.46
1:D:122[B]:LYS:NZ	1:D:171:GLU:OE2	2.44	0.46
1:C:129:THR:HG23	1:C:175:ARG:HD3	1.97	0.45
1:C:77:GLN:HG2	1:C:82:GLN:OE1	2.16	0.45
1:B:-4:LEU:HD23	1:B:-1:GLN:HE22	1.82	0.45
1:D:122[B]:LYS:HA	1:D:125:LEU:HD12	1.99	0.45
1:A:88:ASN:OD1	1:A:120:LEU:HD23	2.18	0.44
1:B:206[A]:GLN:NE2	1:D:148:ASN:OD1	2.51	0.44
1:A:6:LYS:HE3	1:A:133:TRP:CH2	2.53	0.43
1:B:106:TRP:HB3	1:B:153:ASN:ND2	2.33	0.43
1:B:106:TRP:HB3	1:B:153:ASN:HD22	1.83	0.43
1:C:102:LYS:HA	3:C:2008:HOH:O	2.18	0.43
1:C:206[A]:GLN:HB2	1:C:244:MET:HE3	2.00	0.43
1:D:129:THR:HG23	1:D:175:ARG:HD3	2.01	0.42
1:C:106:TRP:HB3	1:C:153:ASN:ND2	2.34	0.42
1:B:88:ASN:OD1	1:B:120:LEU:HD23	2.19	0.42
1:C:88:ASN:OD1	1:C:120:LEU:HD23	2.20	0.42
1:D:73[A]:LEU:CD1	1:D:123:ALA:HB1	2.50	0.41
1:C:8:ALA:HA	1:C:85:ILE:O	2.21	0.41
1:A:73[A]:LEU:CD1	1:A:123:ALA:HB1	2.50	0.41
1:A:109:VAL:CG1	1:A:153:ASN:HD21	2.33	0.41
1:A:146:MET:SD	1:A:199:GLN:OE1	2.79	0.41
1:D:88:ASN:OD1	1:D:120:LEU:HD23	2.21	0.41
1:B:73[A]:LEU:CD1	1:B:123:ALA:HB1	2.51	0.41
1:B:206[B]:GLN:HB2	1:B:244:MET:HE3	2.03	0.41
1:C:206[B]:GLN:HB2	1:C:244:MET:HE3	2.01	0.41
1:D:63:VAL:O	1:D:119:ARG:HD2	2.21	0.40
1:B:63:VAL:O	1:B:119:ARG:HD2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	239/269 (89%)	236 (99%)	3 (1%)	0	100	100
1	B	240/269 (89%)	236 (98%)	4 (2%)	0	100	100
1	C	240/269 (89%)	234 (98%)	6 (2%)	0	100	100
1	D	239/269 (89%)	236 (99%)	3 (1%)	0	100	100
All	All	958/1076 (89%)	942 (98%)	16 (2%)	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	178/201 (89%)	176 (99%)	2 (1%)	65	81
1	B	179/201 (89%)	176 (98%)	3 (2%)	53	72
1	C	179/201 (89%)	176 (98%)	3 (2%)	53	72
1	D	177/201 (88%)	173 (98%)	4 (2%)	44	63
All	All	713/804 (89%)	701 (98%)	12 (2%)	53	72

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	GLU
1	A	151	GLN

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Mol	Chain	Res	Type
1	B	0	SER
1	B	82	GLN
1	B	151	GLN
1	C	6	LYS
1	C	67	GLU
1	C	151	GLN
1	D	43	GLU
1	D	132	ARG
1	D	151	GLN
1	D	203	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-1	GLN
1	A	19	GLN
1	A	79	HIS
1	A	111	ASN
1	A	153	ASN
1	A	199	GLN
1	B	-1	GLN
1	B	19	GLN
1	B	79	HIS
1	B	151	GLN
1	B	153	ASN
1	C	4	GLN
1	C	79	HIS
1	C	151	GLN
1	C	153	ASN
1	D	4	GLN
1	D	77	GLN
1	D	79	HIS
1	D	82	GLN
1	D	89	ASN
1	D	151	GLN
1	D	153	ASN
1	D	199	GLN

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](#)

2 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	O74	A	1247	-	28,28,28	2.30	8 (28%)	32,39,39	2.47	8 (25%)
2	O74	C	1247	-	28,28,28	2.33	9 (32%)	32,39,39	2.25	8 (25%)

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	O74	A	1247	-	-	0/4/19/19	0/4/4/4
2	O74	C	1247	-	-	0/4/19/19	0/4/4/4

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1247	O74	C5-C6	-7.82	1.37	1.46
2	A	1247	O74	C5-C6	-6.41	1.39	1.46
2	C	1247	O74	CAE-NAJ	-5.03	1.31	1.41
2	C	1247	O74	OAU-CAV	4.67	1.37	1.24
2	A	1247	O74	C6-NAJ	-4.27	1.28	1.36
2	A	1247	O74	CAE-NAJ	-4.24	1.33	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1247	O74	CAX-CAV	-3.81	1.35	1.45
2	A	1247	O74	OAU-CAV	3.78	1.34	1.24
2	A	1247	O74	CAB-CL	-3.27	1.65	1.73
2	C	1247	O74	CAS-CAV	-3.04	1.38	1.47
2	A	1247	O74	C4-N3	-2.91	1.34	1.39
2	A	1247	O74	CAS-CAV	-2.79	1.38	1.47
2	C	1247	O74	C4-N3	-2.66	1.35	1.39
2	C	1247	O74	C6-NAJ	-2.34	1.31	1.36
2	C	1247	O74	CAB-CL	-2.12	1.68	1.73
2	C	1247	O74	C2-N3	-2.10	1.33	1.36
2	C	1247	O74	CAX-CAV	-2.03	1.40	1.45

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1247	O74	CAE-CAB-CL	-8.03	110.82	119.52
2	C	1247	O74	CAB-CAE-NAJ	-7.02	108.43	119.25
2	A	1247	O74	CAB-CAE-NAJ	-6.94	108.56	119.25
2	C	1247	O74	CAE-CAB-CL	-4.65	114.49	119.52
2	C	1247	O74	CAT-CAS-C2	-4.62	117.28	121.12
2	A	1247	O74	OAU-CAV-CAS	-4.21	115.48	121.53
2	A	1247	O74	CAF-CAE-NAJ	3.75	130.51	121.82
2	C	1247	O74	CAF-CAE-NAJ	3.50	129.93	121.82
2	A	1247	O74	CAC-CAB-CL	3.44	125.20	118.42
2	C	1247	O74	CAI-C5-C4	3.25	122.47	118.81
2	C	1247	O74	CAW-CAT-CAS	2.80	122.18	117.59
2	C	1247	O74	CAF-CAE-CAB	2.79	121.62	117.89
2	C	1247	O74	CAC-CAB-CL	2.47	123.28	118.42
2	A	1247	O74	CAD-CAC-CAB	-2.30	115.96	119.42
2	A	1247	O74	CAF-CAE-CAB	2.26	120.91	117.89
2	A	1247	O74	CAX-CAV-CAS	2.15	121.21	117.45

There are no chirality outliers.

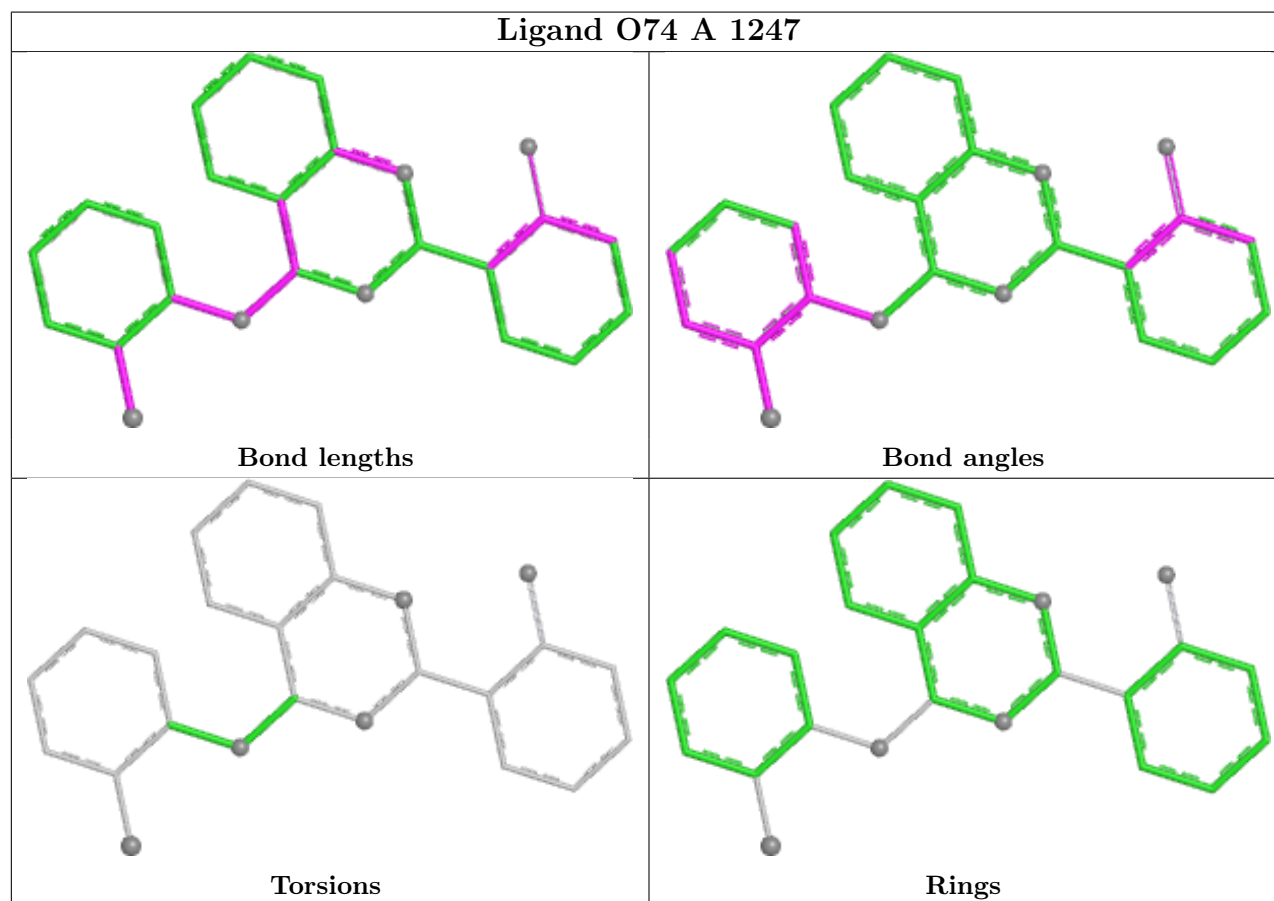
There are no torsion outliers.

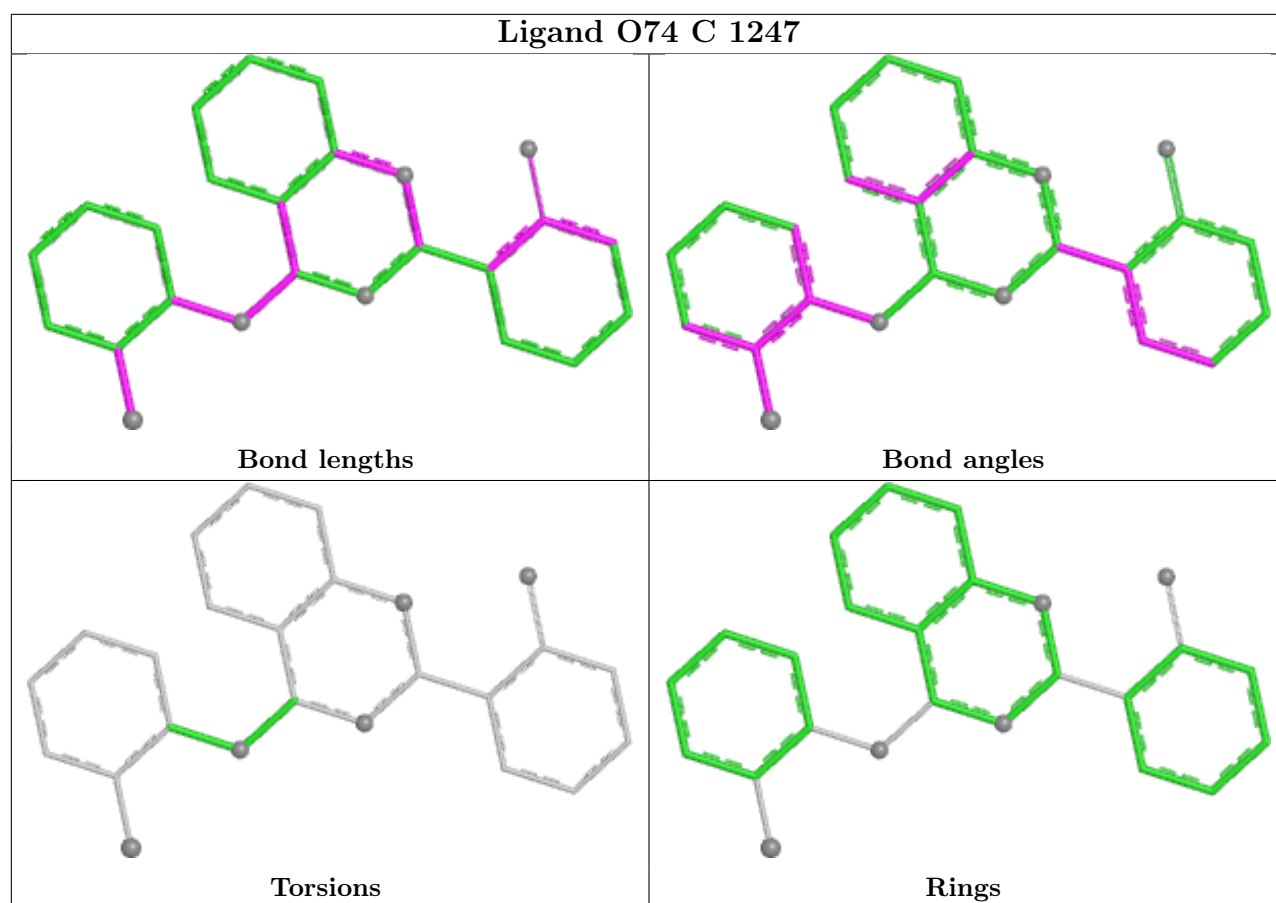
There are no ring outliers.

2 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1247	O74	9	0
2	C	1247	O74	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 [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	239/269 (88%)	0.34	10 (4%)	40	42	28, 50, 88, 111	4 (1%)
1	B	241/269 (89%)	0.34	14 (5%)	29	31	21, 46, 91, 127	3 (1%)
1	C	242/269 (89%)	0.19	9 (3%)	45	48	23, 45, 84, 111	2 (0%)
1	D	240/269 (89%)	0.30	10 (4%)	40	42	25, 47, 82, 109	3 (1%)
All	All	962/1076 (89%)	0.29	43 (4%)	38	40	21, 48, 85, 127	12 (1%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	90	ALA	7.3
1	B	246	MET	5.3
1	B	-4	LEU	5.1
1	A	1	MET	5.0
1	A	191	MET	4.4
1	D	1	MET	4.3
1	A	246	MET	4.2
1	D	247	SER	4.0
1	D	91	GLY	3.8
1	D	90	ALA	3.6
1	B	101	MET	3.5
1	B	146	MET	3.4
1	B	145	ALA	3.3
1	B	90	ALA	3.2
1	D	100	ARG	3.1
1	C	246	MET	3.0
1	B	-3	TYR	3.0
1	C	90	ALA	2.9
1	D	0	SER	2.9
1	A	146	MET	2.8
1	B	74	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	73[A]	LEU	2.6
1	C	1	MET	2.6
1	C	146	MET	2.6
1	C	-6	GLU	2.5
1	D	122[A]	LYS	2.4
1	B	77	GLN	2.4
1	B	71	ALA	2.4
1	B	80	LEU	2.4
1	A	147	GLY	2.3
1	C	145	ALA	2.3
1	B	89	ASN	2.3
1	A	41	GLY	2.2
1	B	103	ASP	2.2
1	D	101	MET	2.2
1	C	78	GLN	2.2
1	A	0	SER	2.2
1	D	80	LEU	2.1
1	A	4	GLN	2.1
1	C	103	ASP	2.1
1	C	130	LYS	2.1
1	A	145	ALA	2.0
1	B	70	ALA	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	O74	A	1247	25/25	0.89	0.12	41,48,80,83	0

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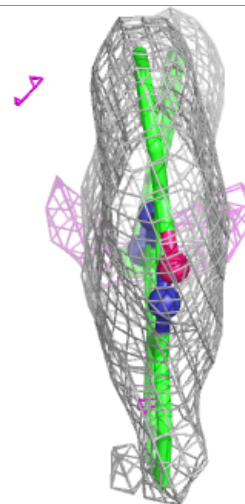
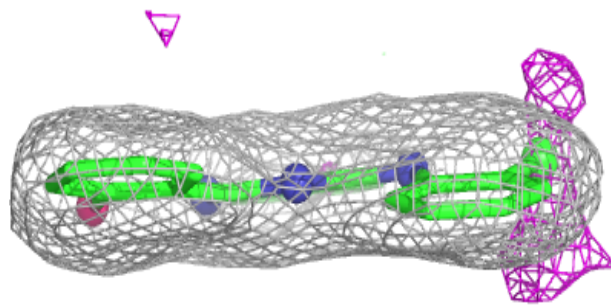
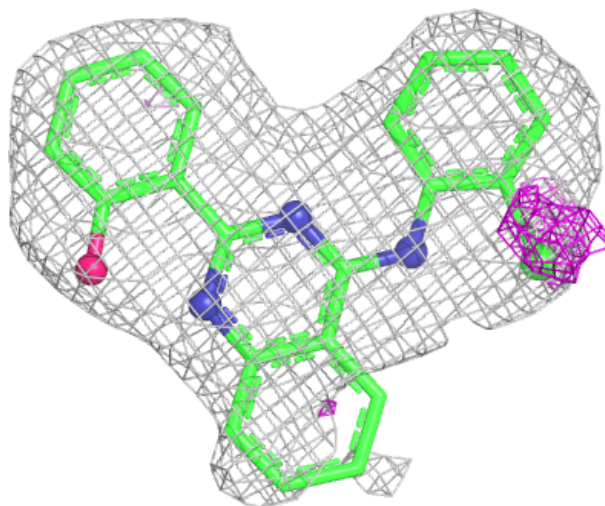
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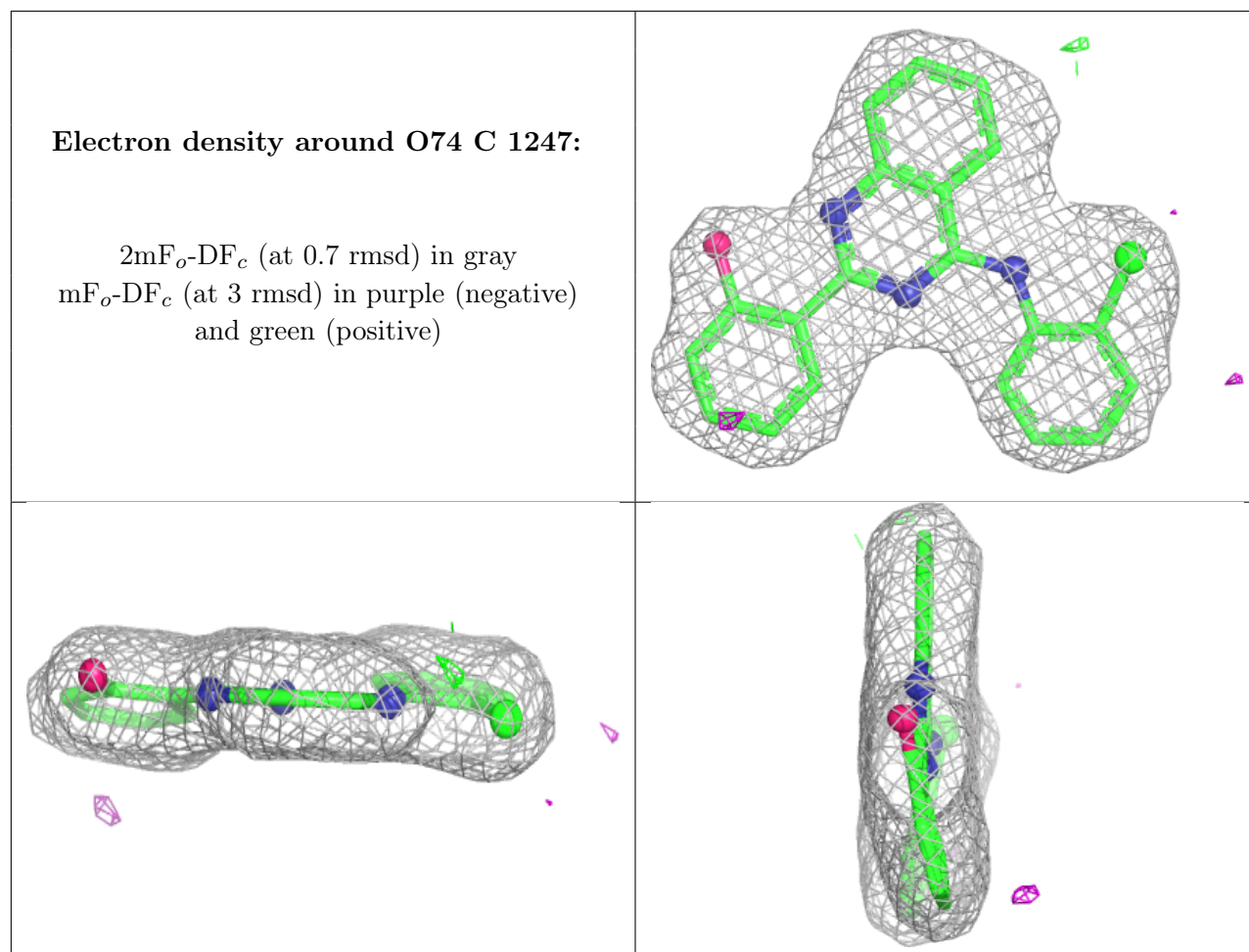
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	O74	C	1247	25/25	0.94	0.08	39,47,54,72	0

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 O74 A 1247:

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 ⓘ

There are no such residues in this entry.