



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:14 PM UTC

PDB ID : 8XLO / pdb_00008xlo
Title : FGFR1 kinase domain with a dual-warhead covalent inhibitor CXF-007
Authors : Chen, X.J.; Chen, Y.H.
Deposited on : 2023-12-26
Resolution : 2.36 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

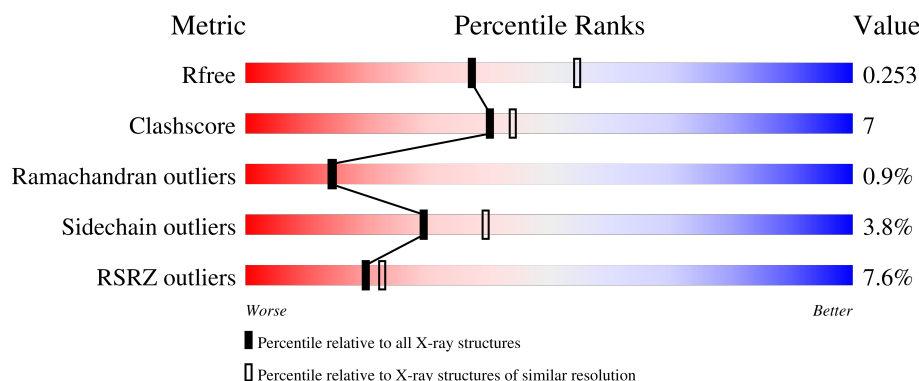
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.36 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	310	
1	B	310	

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
3	SO4	A	804	-	-	X	-
3	SO4	B	804	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4664 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.

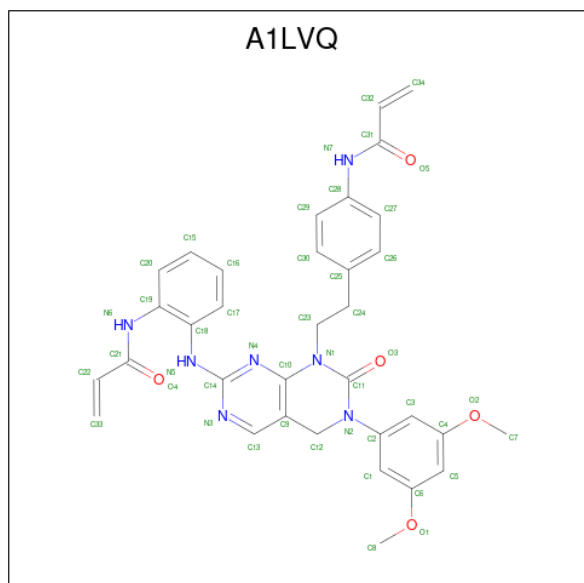
- Molecule 1 is a protein called Fibroblast growth factor receptor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	293	Total	C	N	O	S	0	0	0
			2278	1446	387	427	18			
1	B	289	Total	C	N	O	S	0	1	0
			2227	1415	378	416	18			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	456	GLY	-	expression tag	UNP P11362
A	457	PRO	-	expression tag	UNP P11362
A	584	SER	CYS	engineered mutation	UNP P11362
B	456	GLY	-	expression tag	UNP P11362
B	457	PRO	-	expression tag	UNP P11362
B	584	SER	CYS	engineered mutation	UNP P11362

- Molecule 2 is CXF007 (CCD ID: A1LVQ) (formula: $C_{34}H_{33}N_7O_5$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			46	34	7	5		
2	B	1	Total	C	N	O	0	0
			46	34	7	5		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

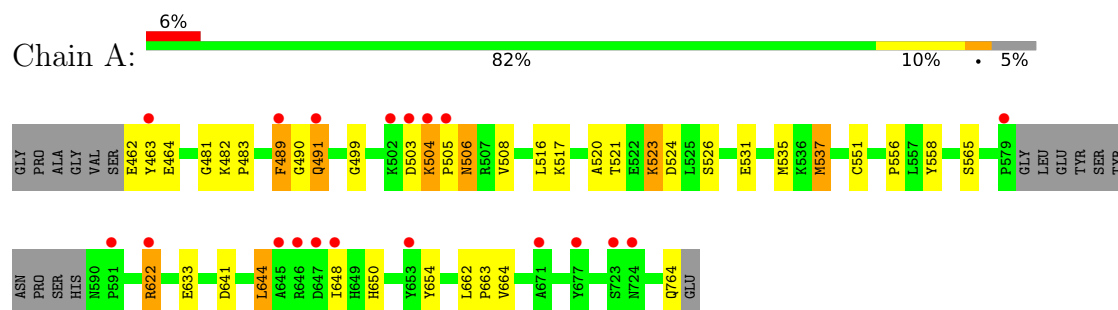
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	25	Total	O	0	0
			25	25		
4	B	12	Total	O	0	0
			12	12		

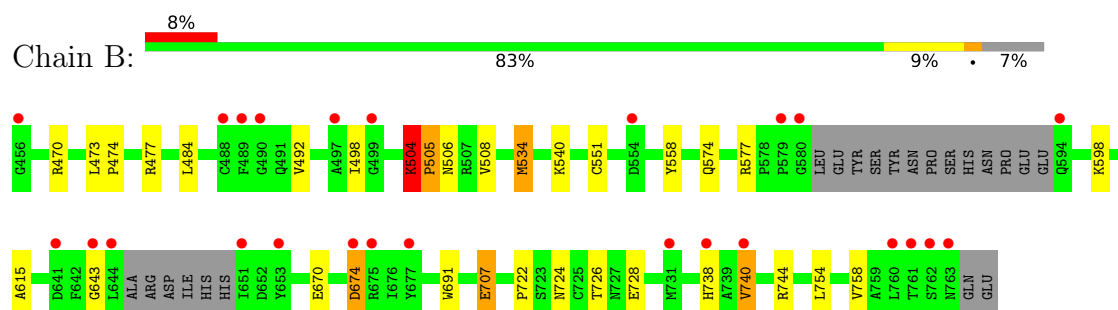
3 Residue-property plots [i](#)

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: Fibroblast growth factor receptor 1



- Molecule 1: Fibroblast growth factor receptor 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	209.21Å 57.48Å 66.41Å 90.00° 107.86° 90.00°	Depositor
Resolution (Å)	46.68 – 2.36 46.68 – 2.36	Depositor EDS
% Data completeness (in resolution range)	99.6 (46.68-2.36) 99.6 (46.68-2.36)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 2.37Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.208 , 0.255 0.212 , 0.253	Depositor DCC
R_{free} test set	1423 reflections (4.56%)	wwPDB-VP
Wilson B-factor (Å ²)	58.0	Xtriage
Anisotropy	0.545	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 53.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.018 for -h-2*k,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4664	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.58% 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: A1LVQ, SO4

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	0.46	0/2326	0.79	2/3155 (0.1%)
1	B	0.30	0/2277	0.79	5/3092 (0.2%)
All	All	0.39	0/4603	0.79	7/6247 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	504	LYS	CA-C-N	8.76	130.79	119.84
1	B	504	LYS	C-N-CA	8.76	130.79	119.84
1	A	506	ASN	N-CA-C	-7.10	102.92	112.94
1	B	674	ASP	N-CA-C	-5.37	108.50	114.62
1	B	740	VAL	CA-C-N	5.19	124.99	119.28
1	B	740	VAL	C-N-CA	5.19	124.99	119.28
1	A	491	GLN	N-CA-C	-5.09	106.58	112.89

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	2278	0	2220	34	0
1	B	2227	0	2164	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	46	0	0	4	0
2	B	46	0	0	3	0
3	A	15	0	0	4	0
3	B	15	0	0	3	0
4	A	25	0	0	0	0
4	B	12	0	0	0	0
All	All	4664	0	4384	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 7.

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:489:PHE:CZ	1:A:491:GLN:NE2	2.26	1.02
1:B:504:LYS:H	1:B:505:PRO:CD	1.83	0.91
2:A:801:A1LVQ:C33	3:A:804:SO4:S	2.61	0.89
1:A:489:PHE:CE2	1:A:491:GLN:NE2	2.42	0.88
1:B:504:LYS:N	1:B:505:PRO:CD	2.35	0.88
1:B:504:LYS:N	1:B:505:PRO:HD2	1.87	0.88
2:B:801:A1LVQ:C33	3:B:804:SO4:O2	2.27	0.82
1:A:622:ARG:HB3	1:A:622:ARG:CZ	2.09	0.80
2:B:801:A1LVQ:C33	3:B:804:SO4:S	2.69	0.80
1:A:622:ARG:H	1:A:622:ARG:NH1	1.89	0.70
1:B:477:ARG:HG2	1:B:498:ILE:HB	1.77	0.65
2:A:801:A1LVQ:C33	3:A:804:SO4:O1	2.45	0.64
1:B:504:LYS:H	1:B:505:PRO:HD3	1.64	0.61
1:B:707:GLU:OE1	1:B:707:GLU:N	2.26	0.61
1:A:551:CYS:HB2	1:A:558:TYR:HB2	1.82	0.61
1:A:463:TYR:CZ	1:A:556:PRO:HA	2.35	0.61
1:B:574:GLN:HA	1:B:577:ARG:HD2	1.83	0.60
2:A:801:A1LVQ:C33	3:A:804:SO4:O4	2.50	0.60
1:A:462:GLU:HG2	1:A:462:GLU:O	2.05	0.55
1:B:504:LYS:O	1:B:506:ASN:N	2.41	0.53
1:A:622:ARG:CG	1:A:622:ARG:HH11	2.22	0.53
1:A:622:ARG:NH1	1:A:622:ARG:CB	2.73	0.52
1:A:504:LYS:O	1:A:504:LYS:HG2	2.10	0.51
1:A:463:TYR:CE2	1:A:556:PRO:HA	2.46	0.51
1:B:484:LEU:HD12	1:B:492:VAL:HG12	1.92	0.51
1:A:483:PRO:HB3	1:A:490:GLY:HA2	1.94	0.50
1:A:641:ASP:O	1:A:644:LEU:HD12	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:670:GLU:OE2	1:B:744:ARG:NH1	2.44	0.49
1:A:622:ARG:HE	1:A:644:LEU:HB2	1.76	0.49
1:A:523:LYS:HD3	1:A:526:SER:OG	2.14	0.48
1:B:551:CYS:HB2	1:B:558:TYR:HB2	1.95	0.48
1:A:622:ARG:CZ	1:A:622:ARG:CB	2.84	0.47
1:A:537:MET:HE3	1:A:537:MET:HB2	1.80	0.47
1:B:473:LEU:HD12	1:B:474:PRO:HD2	1.97	0.47
1:A:654:TYR:CD1	1:A:654:TYR:N	2.83	0.46
1:A:622:ARG:NH1	1:A:622:ARG:N	2.60	0.46
1:B:498:ILE:HG22	1:B:506:ASN:HA	1.98	0.45
1:A:641:ASP:HB3	1:A:644:LEU:CD1	2.47	0.45
2:A:801:A1LVQ:C33	3:A:804:SO4:O2	2.64	0.45
1:A:641:ASP:HB3	1:A:644:LEU:HD11	1.99	0.45
1:B:534:MET:SD	1:B:643:GLY:HA2	2.57	0.45
1:A:622:ARG:NH1	1:A:622:ARG:CG	2.81	0.44
1:B:726:THR:OG1	1:B:728:GLU:OE1	2.31	0.43
1:A:622:ARG:HH11	1:A:622:ARG:HG2	1.83	0.43
1:A:481:GLY:C	1:A:482:LYS:HD3	2.44	0.43
1:A:521:THR:HG23	1:A:524:ASP:OD2	2.19	0.42
1:B:670:GLU:O	1:B:674:ASP:HB3	2.19	0.42
1:A:565:SER:OG	1:A:633:GLU:HG2	2.20	0.41
1:A:662:LEU:HA	1:A:663:PRO:HD2	1.89	0.41
1:B:691:TRP:CD1	1:B:722:PRO:HG3	2.56	0.41
1:B:754:LEU:O	1:B:758:VAL:HG23	2.21	0.41
1:A:516:LEU:HD23	1:A:520:ALA:HB1	2.02	0.41
1:A:499:GLY:HA2	1:A:505:PRO:HA	2.03	0.41
1:A:504:LYS:HE2	1:A:506:ASN:HB2	2.03	0.41
1:A:662:LEU:HB3	1:A:664:VAL:HG12	2.03	0.41
1:A:764:GLN:HE22	1:B:615:ALA:HB1	1.86	0.40
2:B:801:A1LVQ:C33	3:B:804:SO4:O1	2.69	0.40
1:A:531:GLU:O	1:A:535:MET:HG3	2.22	0.40
1:A:622:ARG:HB3	1:A:622:ARG:NH1	2.31	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	289/310 (93%)	277 (96%)	9 (3%)	3 (1%)	12	12
1	B	284/310 (92%)	275 (97%)	7 (2%)	2 (1%)	18	20
All	All	573/620 (92%)	552 (96%)	16 (3%)	5 (1%)	14	14

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	648	ILE
1	A	650	HIS
1	B	504	LYS
1	B	505	PRO
1	A	504	LYS

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	242/271 (89%)	233 (96%)	9 (4%)	30	40
1	B	235/271 (87%)	226 (96%)	9 (4%)	29	39
All	All	477/542 (88%)	459 (96%)	18 (4%)	29	39

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	464	GLU
1	A	489	PHE
1	A	503	ASP
1	A	508	VAL
1	A	517	LYS
1	A	523	LYS
1	A	537	MET

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Mol	Chain	Res	Type
1	A	622	ARG
1	A	644	LEU
1	B	470	ARG
1	B	508	VAL
1	B	534	MET
1	B	540	LYS
1	B	598	LYS
1	B	707	GLU
1	B	724	ASN
1	B	738	HIS
1	B	740	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	635	ASN
1	A	743	GLN
1	B	543	ASN
1	B	763	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](#)

8 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	A1LVQ	B	801	1	50,50,50	0.95	2 (4%)	62,69,69	1.32	4 (6%)
2	A1LVQ	A	801	1	50,50,50	0.95	2 (4%)	62,69,69	1.32	3 (4%)
3	SO4	B	802	-	4,4,4	0.24	0	6,6,6	0.08	0
3	SO4	B	804	-	4,4,4	0.24	0	6,6,6	0.08	0
3	SO4	A	802	-	4,4,4	0.24	0	6,6,6	0.07	0
3	SO4	B	803	-	4,4,4	0.24	0	6,6,6	0.08	0
3	SO4	A	804	-	4,4,4	0.22	0	6,6,6	0.08	0
3	SO4	A	803	-	4,4,4	0.24	0	6,6,6	0.08	0

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	A1LVQ	B	801	1	-	9/29/45/45	0/5/5/5
2	A1LVQ	A	801	1	-	9/29/45/45	0/5/5/5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	A1LVQ	C34-C32	4.14	1.50	1.30
2	B	801	A1LVQ	C34-C32	4.12	1.50	1.30
2	B	801	A1LVQ	C13-C9	3.99	1.45	1.37
2	A	801	A1LVQ	C13-C9	3.95	1.45	1.37

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	A1LVQ	C34-C32-C31	-8.21	113.20	122.22
2	B	801	A1LVQ	C34-C32-C31	-7.78	113.67	122.22
2	B	801	A1LVQ	C10-N1-C11	-2.69	119.81	122.23
2	B	801	A1LVQ	C9-C13-N3	-2.27	120.14	123.83
2	A	801	A1LVQ	C9-C13-N3	-2.22	120.23	123.83
2	B	801	A1LVQ	C13-C9-C10	2.05	118.01	115.51
2	A	801	A1LVQ	O3-C11-N1	-2.01	119.55	122.15

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	A1LVQ	O5-C31-N7-C28
2	A	801	A1LVQ	C32-C31-N7-C28
2	A	801	A1LVQ	O4-C21-C22-C33
2	A	801	A1LVQ	N6-C21-C22-C33
2	A	801	A1LVQ	O5-C31-C32-C34
2	A	801	A1LVQ	N7-C31-C32-C34
2	B	801	A1LVQ	O4-C21-N6-C19
2	B	801	A1LVQ	C22-C21-N6-C19
2	B	801	A1LVQ	O4-C21-C22-C33
2	B	801	A1LVQ	N6-C21-C22-C33
2	B	801	A1LVQ	O5-C31-C32-C34
2	B	801	A1LVQ	N7-C31-C32-C34
2	A	801	A1LVQ	C23-C24-C25-C30
2	A	801	A1LVQ	C23-C24-C25-C26
2	A	801	A1LVQ	C3-C2-N2-C11
2	B	801	A1LVQ	C1-C2-N2-C11
2	B	801	A1LVQ	C3-C2-N2-C11
2	B	801	A1LVQ	N4-C14-N5-C18

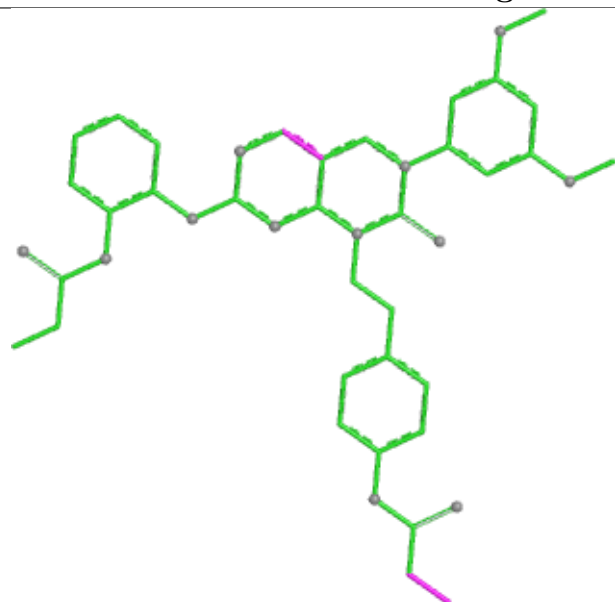
There are no ring outliers.

4 monomers are involved in 7 short contacts:

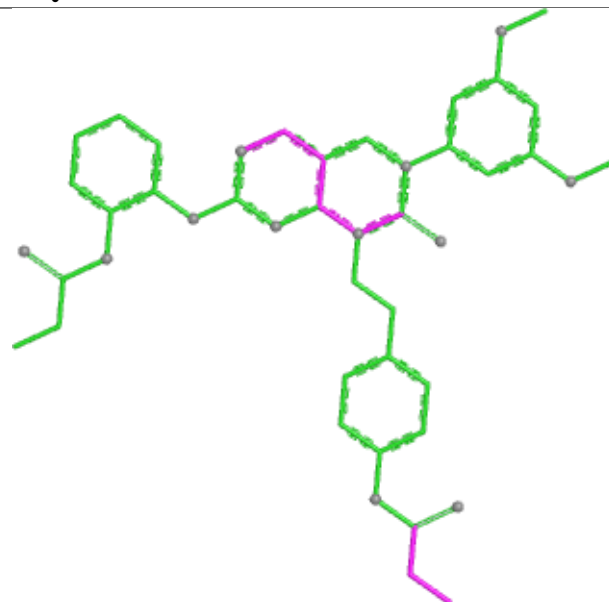
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	801	A1LVQ	3	0
2	A	801	A1LVQ	4	0
3	B	804	SO4	3	0
3	A	804	SO4	4	0

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.

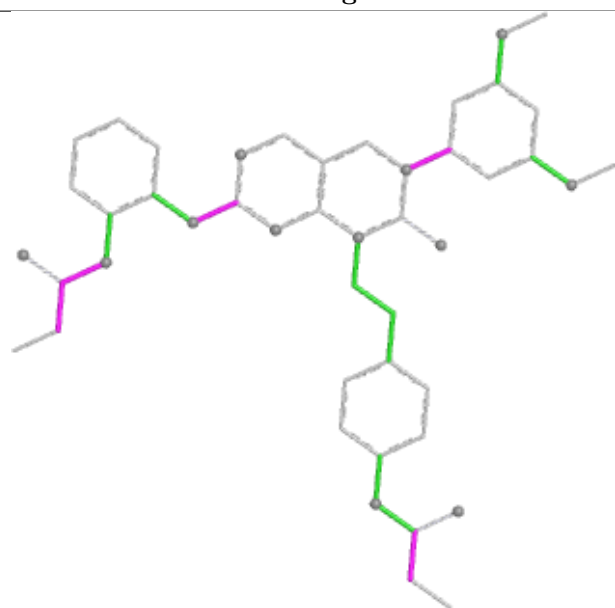
Ligand A1LVQ B 801



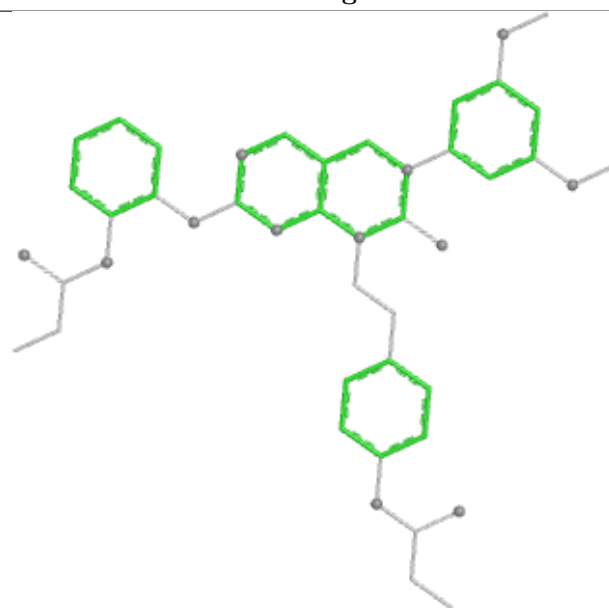
Bond lengths



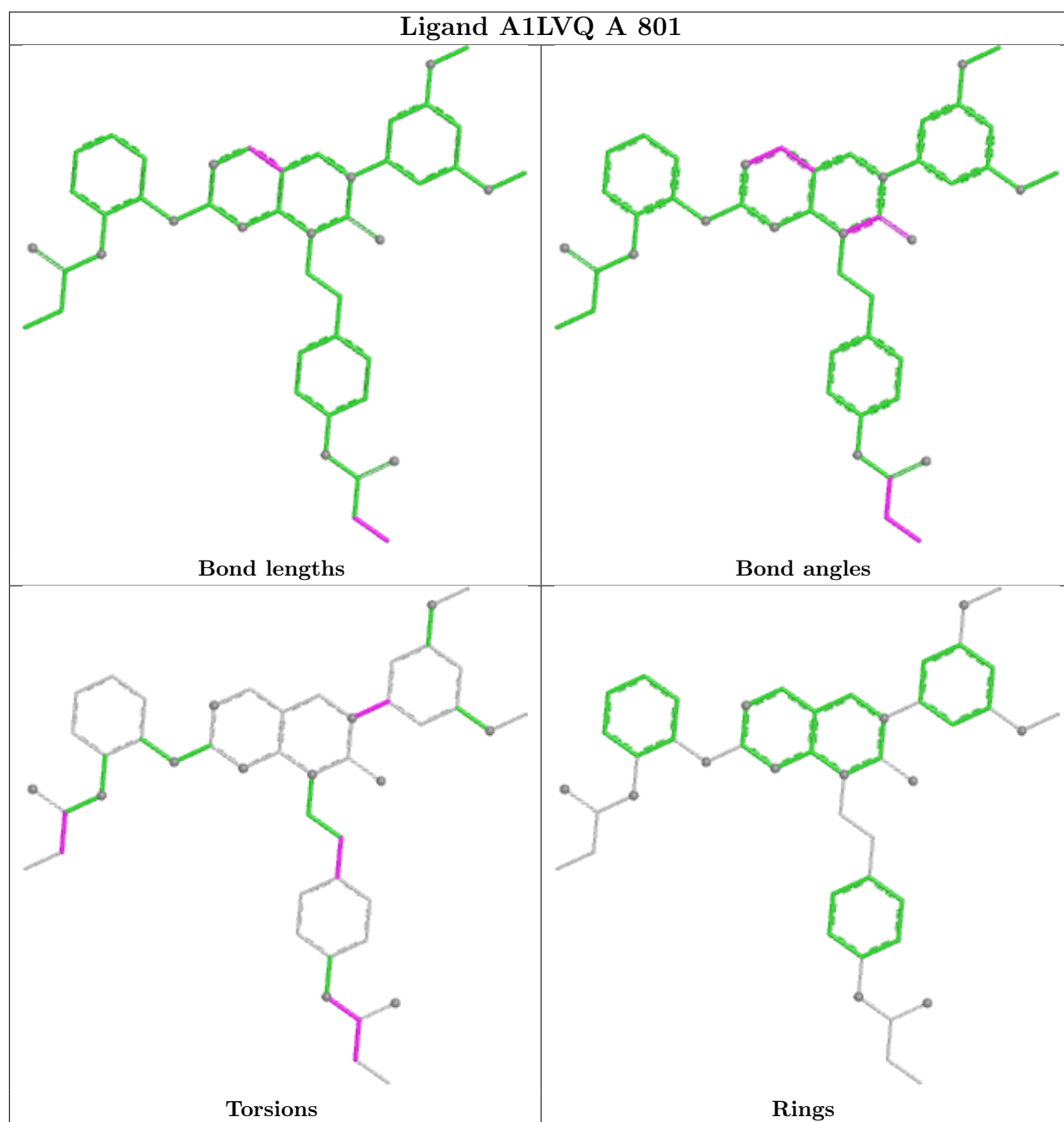
Bond angles



Torsions



Rings



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	293/310 (94%)	0.57	19 (6%)	25 28	48, 65, 113, 143	0
1	B	289/310 (93%)	0.79	25 (8%)	16 18	48, 71, 104, 126	1 (0%)
All	All	582/620 (93%)	0.68	44 (7%)	20 22	48, 68, 108, 143	1 (0%)

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	489	PHE	4.6
1	B	651	ILE	4.5
1	B	644	LEU	4.4
1	B	643	GLY	4.2
1	B	489	PHE	4.1
1	A	463	TYR	3.9
1	A	677	TYR	3.6
1	A	622	ARG	3.6
1	B	674	ASP	3.5
1	A	645	ALA	3.4
1	B	760	LEU	3.3
1	A	591	PRO	3.3
1	A	504	LYS	3.2
1	B	580	GLY	3.0
1	B	675	ARG	3.0
1	A	646	ARG	2.9
1	B	731	MET	2.9
1	B	677	TYR	2.8
1	B	763	ASN	2.7
1	B	456	GLY	2.7
1	B	499	GLY	2.6
1	B	738	HIS	2.6
1	A	671	ALA	2.6
1	A	647	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	502	LYS	2.4
1	A	503	ASP	2.4
1	B	490	GLY	2.4
1	A	505	PRO	2.3
1	B	761	THR	2.3
1	B	594	GLN	2.3
1	A	724	ASN	2.3
1	B	579	PRO	2.3
1	B	762	SER	2.3
1	B	653	TYR	2.3
1	B	554	ASP	2.2
1	A	579	PRO	2.2
1	B	740	VAL	2.2
1	B	488	CYS	2.2
1	A	653	TYR	2.1
1	A	648	ILE	2.1
1	A	491	GLN	2.1
1	B	497	ALA	2.1
1	A	723	SER	2.0
1	B	641	ASP	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
3	SO4	A	803	5/5	0.64	0.14	88,89,108,125	0
3	SO4	B	803	5/5	0.68	0.11	87,89,110,121	0
3	SO4	A	804	5/5	0.72	0.14	86,98,113,136	0

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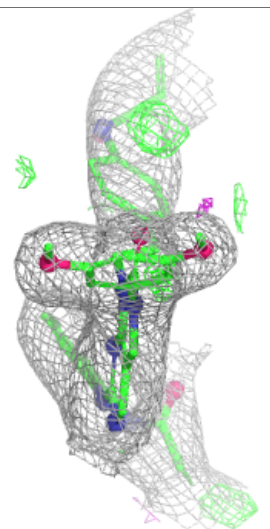
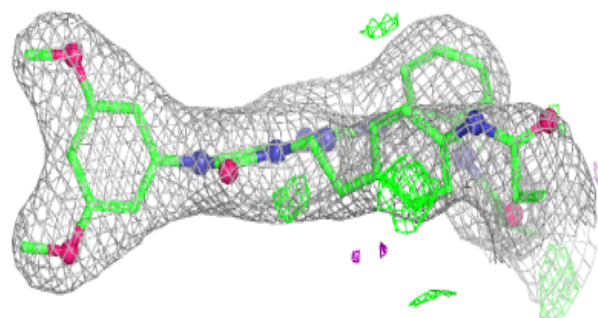
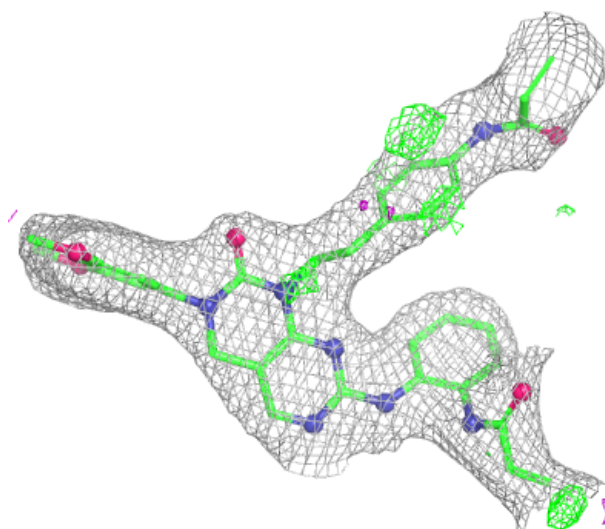
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	B	804	5/5	0.74	0.14	81,120,138,146	0
2	A1LVQ	A	801	46/46	0.93	0.12	50,65,99,105	0
2	A1LVQ	B	801	46/46	0.94	0.11	52,67,97,99	0
3	SO4	A	802	5/5	0.96	0.09	66,70,73,75	0
3	SO4	B	802	5/5	0.97	0.07	64,69,75,83	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.

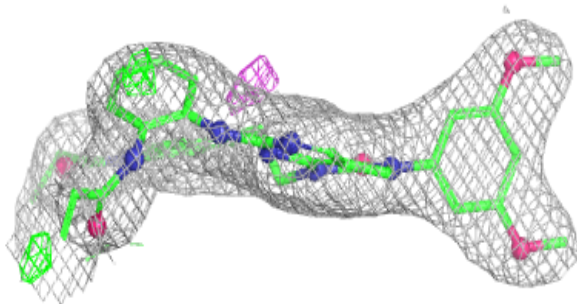
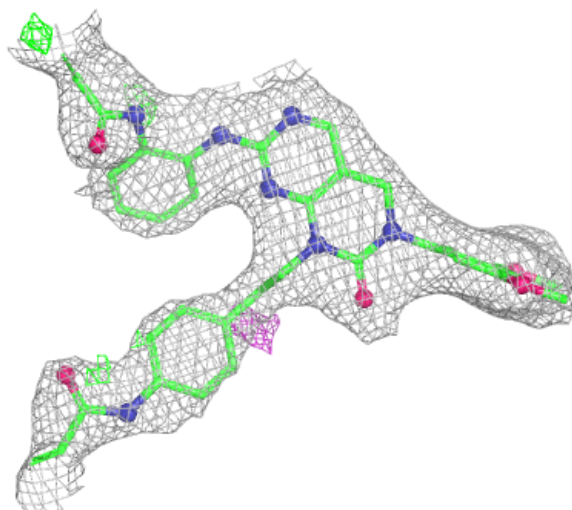
Electron density around A1LVQ A 801:

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 A1LVQ B 801:

$2mF_o-DF_c$ (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.