



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 4BIC / pdb_00004bic
Title : Crystal Structures of Ask1-inhibitor Complexes
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Deposited on : 2013-04-10
Resolution : 2.62 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

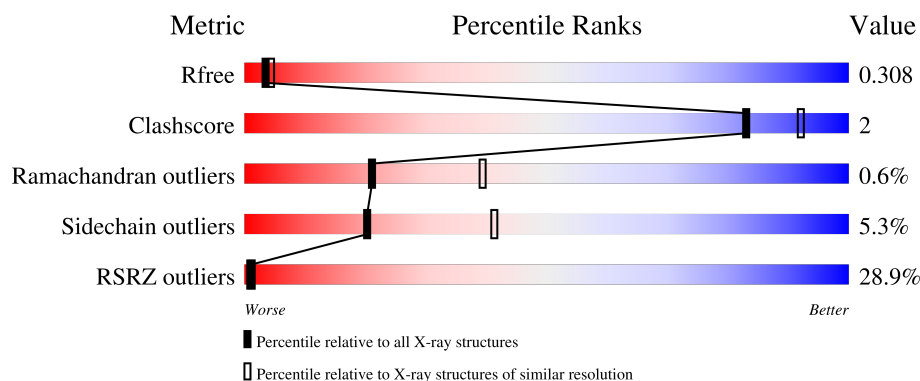
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.62 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	4951 (2.64-2.60)
Clashscore	190562	5303 (2.64-2.60)
Ramachandran outliers	187476	5217 (2.64-2.60)
Sidechain outliers	187428	5217 (2.64-2.60)
RSRZ outliers	180081	4950 (2.64-2.60)

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	318	12% 72% 9% 18%
1	B	318	35% 72% 9% 19%

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	GOL	A	1941	-	X	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4162 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.

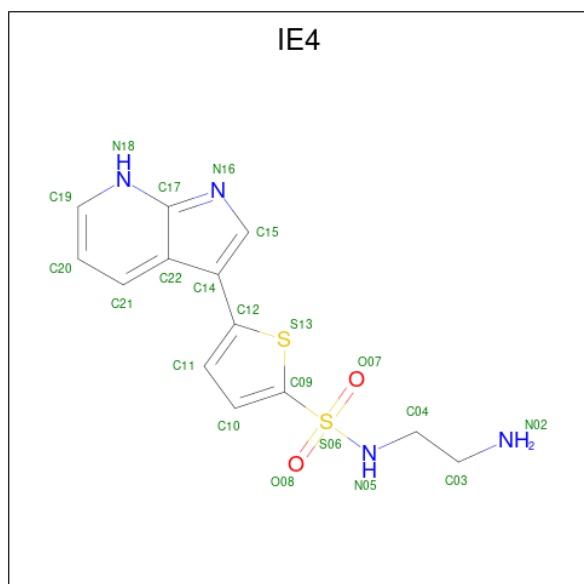
- Molecule 1 is a protein called MITOGEN-ACTIVATED PROTEIN KINASE KINASE KINASE 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	260	Total	C	N	O	S	0	0	0
			2021	1302	330	380	9			
1	B	259	Total	C	N	O	S	0	0	0
			1979	1271	323	376	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	838	GLU	THR	engineered mutation	UNP Q99683
B	838	GLU	THR	engineered mutation	UNP Q99683

- Molecule 2 is N-(2-aminoethyl)-5-{1H-pyrrolo[2,3-b]pyridin-3-yl}thiophene-2-sulfonamide (CCD ID: IE4) (formula: C₁₃H₁₄N₄O₂S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			21	13	4	2	2		
2	A	1	Total	C	N	O	S	0	0
			21	13	4	2	2		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

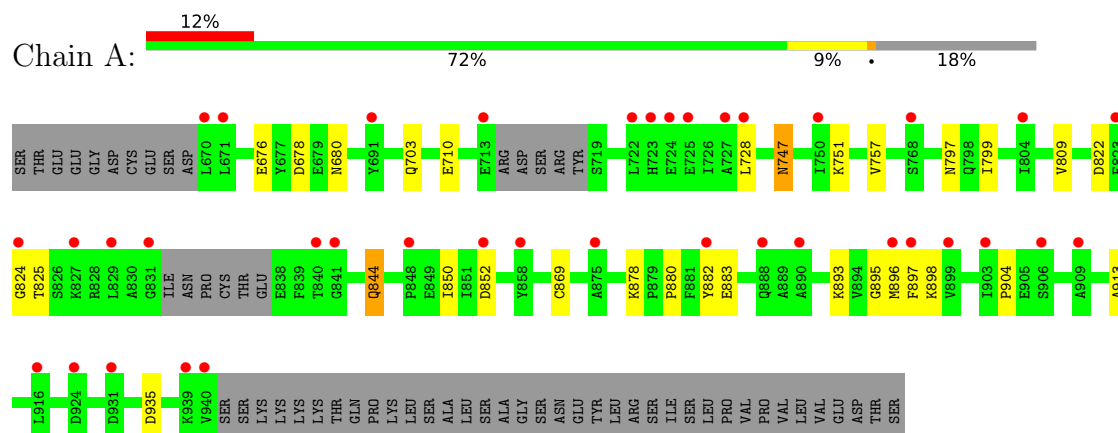
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	74	Total	O	0	0
			74	74		
4	B	34	Total	O	0	0
			34	34		

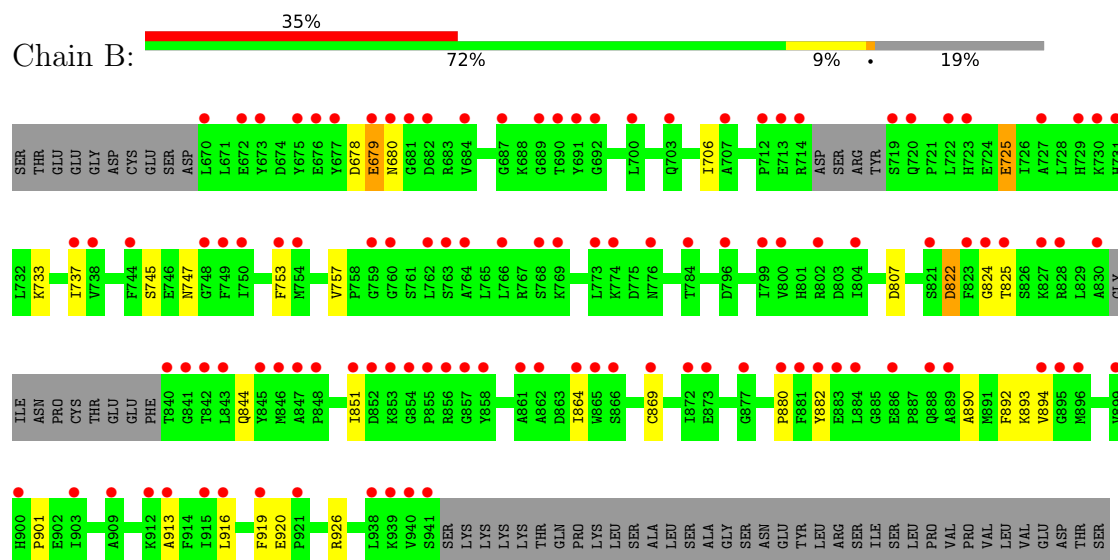
3 Residue-property plots [i](#)

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: MITOGEN-ACTIVATED PROTEIN KINASE KINASE KINASE 5



• Molecule 1: MITOGEN-ACTIVATED PROTEIN KINASE KINASE KINASE 5



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	77.95Å 77.95Å 418.43Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.84 – 2.62 29.84 – 2.62	Depositor EDS
% Data completeness (in resolution range)	99.3 (29.84-2.62) 99.1 (29.84-2.62)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.50 (at 2.61Å)	Xtriage
Refinement program	BUSTER 2.11.1	Depositor
R, R_{free}	0.207 , 0.265 (Not available) , 0.308	Depositor DCC
R_{free} test set	1006 reflections (4.15%)	wwPDB-VP
Wilson B-factor (Å ²)	55.0	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	4162	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.12% 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: GOL, IE4

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	0/2064	1.34	13/2786 (0.5%)
1	B	0.82	0/2020	1.36	9/2732 (0.3%)
All	All	0.86	0/4084	1.35	22/5518 (0.4%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	850	ILE	N-CA-C	-8.08	104.86	111.81
1	A	935	ASP	CA-C-N	6.29	128.99	120.38
1	A	935	ASP	C-N-CA	6.29	128.99	120.38
1	B	807	ASP	CA-CB-CG	6.17	118.77	112.60
1	B	678	ASP	CA-C-N	6.04	128.98	120.28
1	B	678	ASP	C-N-CA	6.04	128.98	120.28
1	A	882	TYR	CA-C-N	5.95	128.25	120.28
1	A	882	TYR	C-N-CA	5.95	128.25	120.28
1	A	904	PRO	CA-C-N	5.94	128.52	120.38
1	A	904	PRO	C-N-CA	5.94	128.52	120.38
1	A	678	ASP	CA-CB-CG	5.84	118.44	112.60
1	B	892	PHE	N-CA-C	-5.80	104.87	111.14
1	B	882	TYR	CA-C-N	5.80	128.63	120.28
1	B	882	TYR	C-N-CA	5.80	128.63	120.28
1	B	678	ASP	CA-CB-CG	5.74	118.34	112.60
1	A	897	PHE	N-CA-C	5.17	119.59	113.28
1	A	824	GLY	CA-C-N	5.11	127.55	120.29
1	A	824	GLY	C-N-CA	5.11	127.55	120.29
1	A	913	ALA	CA-C-N	5.06	127.02	120.44
1	A	913	ALA	C-N-CA	5.06	127.02	120.44
1	B	913	ALA	CA-C-N	5.01	126.95	120.44
1	B	913	ALA	C-N-CA	5.01	126.95	120.44

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	2021	0	1980	9	3
1	B	1979	0	1915	9	3
2	A	42	0	28	3	0
3	A	6	0	8	1	0
3	B	6	0	8	0	0
4	A	74	0	0	2	1
4	B	34	0	0	0	1
All	All	4162	0	3939	19	4

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 2.

All (19) 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:2:IE4:H18	1:B:757:VAL:H	1.31	0.79
1:A:844:GLN:HG2	4:A:2058:HOH:O	1.84	0.78
1:A:757:VAL:H	2:A:1:IE4:H18	1.33	0.73
1:A:895:GLY:O	1:A:898:LYS:NZ	2.27	0.67
1:B:869:CYS:HB3	1:B:880:PRO:HG3	1.80	0.63
1:B:706:ILE:HD12	1:B:753:PHE:HD2	1.70	0.56
1:B:890:ALA:O	1:B:894:VAL:HG23	2.09	0.53
1:A:869:CYS:HB3	1:A:880:PRO:HG3	1.91	0.52
1:B:725:GLU:HG2	1:B:824:GLY:HA2	1.93	0.49
1:A:797:ASN:HB2	1:A:799:ILE:HD12	1.96	0.48
1:B:679:GLU:CD	1:B:679:GLU:H	2.22	0.47
1:B:901:PRO:HD2	1:B:919:PHE:CZ	2.50	0.45
1:B:920:GLU:HB3	1:B:926:ARG:HG3	1.98	0.45
1:A:710:GLU:HG2	1:A:751:LYS:HG2	1.99	0.44
1:A:728:LEU:HD11	1:A:799:ILE:HG12	2.00	0.43
2:A:2:IE4:S13	2:A:2:IE4:H21	2.59	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:703:GLN:O	3:A:1941:GOL:O1	2.25	0.42
1:A:676:GLU:HG2	4:A:2023:HOH:O	2.20	0.42
1:B:844:GLN:HB3	1:B:880:PRO:HG2	2.03	0.41

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:852:ASP:O	1:B:893:LYS:NZ[8_665]	2.08	0.12
1:A:747:ASN:O	4:B:2004:HOH:O[1_655]	2.13	0.07
1:A:680:ASN:OD1	1:B:747:ASN:ND2[1_655]	2.17	0.03
1:B:680:ASN:O	4:A:2003:HOH:O[1_455]	2.19	0.01

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	254/318 (80%)	240 (94%)	12 (5%)	2 (1%)	16	31
1	B	253/318 (80%)	242 (96%)	10 (4%)	1 (0%)	30	49
All	All	507/636 (80%)	482 (95%)	22 (4%)	3 (1%)	21	39

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	822	ASP
1	A	747	ASN
1	B	822	ASP

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	120/274 (44%)	113 (94%)	7 (6%)	18	37
1	B	199/274 (73%)	189 (95%)	10 (5%)	22	44
All	All	319/548 (58%)	302 (95%)	17 (5%)	20	41

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	809	VAL
1	A	825	THR
1	A	844	GLN
1	A	878	LYS
1	A	883	GLU
1	A	893	LYS
1	A	896	MET
1	B	679	GLU
1	B	725	GLU
1	B	733	LYS
1	B	737	ILE
1	B	745	SER
1	B	822	ASP
1	B	825	THR
1	B	851	ILE
1	B	864	ILE
1	B	916	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	798	GLN
1	B	756	GLN

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](#)

4 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	IE4	A	1	-	23,23,23	2.22	4 (17%)	28,33,33	2.96	11 (39%)
3	GOL	B	1942	-	5,5,5	0.82	0	5,5,5	0.99	0
3	GOL	A	1941	-	5,5,5	0.94	0	5,5,5	2.42	4 (80%)
2	IE4	A	2	-	23,23,23	2.15	4 (17%)	28,33,33	3.09	11 (39%)

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	IE4	A	1	-	-	3/15/15/15	0/3/3/3
3	GOL	B	1942	-	-	2/4/4/4	-
3	GOL	A	1941	-	-	4/4/4/4	-
2	IE4	A	2	-	-	6/15/15/15	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	IE4	S06-N05	8.55	1.74	1.61
2	A	2	IE4	S06-N05	8.23	1.73	1.61
2	A	2	IE4	C22-C17	-3.82	1.38	1.45
2	A	1	IE4	C22-C17	-3.45	1.39	1.45
2	A	1	IE4	C17-N18	-2.84	1.33	1.38
2	A	2	IE4	C17-N18	-2.37	1.33	1.38
2	A	1	IE4	O08-S06	2.24	1.46	1.43
2	A	2	IE4	C19-C20	2.05	1.39	1.35

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	IE4	O07-S06-C09	-7.82	97.79	107.21
2	A	2	IE4	O07-S06-C09	-7.30	98.42	107.21
2	A	1	IE4	C20-C19-N18	6.82	127.62	120.22
2	A	2	IE4	S13-C09-S06	-6.34	112.17	120.43
2	A	1	IE4	O07-S06-N05	6.31	113.31	107.41
2	A	2	IE4	C20-C19-N18	5.90	126.62	120.22
2	A	1	IE4	C19-N18-C17	-5.44	111.24	122.89
2	A	2	IE4	O07-S06-N05	5.18	112.25	107.41
2	A	2	IE4	C19-N18-C17	-4.90	112.38	122.89
2	A	2	IE4	C15-N16-C17	4.24	112.17	106.25
2	A	1	IE4	S13-C09-S06	-4.22	114.93	120.43
2	A	2	IE4	C10-C09-S06	3.89	135.11	126.65
2	A	2	IE4	C14-C15-N16	-3.87	104.87	112.05
2	A	2	IE4	C22-C17-N16	-3.32	107.79	115.07
3	A	1941	GOL	C3-C2-C1	-3.26	99.84	111.80
2	A	2	IE4	C15-C14-C22	3.03	110.15	106.98
2	A	1	IE4	C20-C21-C22	-2.94	114.79	120.83
3	A	1941	GOL	O2-C2-C3	2.93	121.30	109.18
2	A	1	IE4	C22-C17-N16	-2.71	109.13	115.07
2	A	2	IE4	O08-S06-O07	2.70	124.76	119.82
2	A	1	IE4	C14-C15-N16	-2.64	107.14	112.05
2	A	1	IE4	C10-C09-S06	2.58	132.26	126.65
2	A	1	IE4	C15-N16-C17	2.37	109.56	106.25
3	A	1941	GOL	O2-C2-C1	-2.20	100.07	109.18
3	A	1941	GOL	O1-C1-C2	-2.15	100.72	110.38
2	A	1	IE4	O08-S06-O07	2.12	123.70	119.82

There are no chirality outliers.

All (15) torsion outliers are listed below:

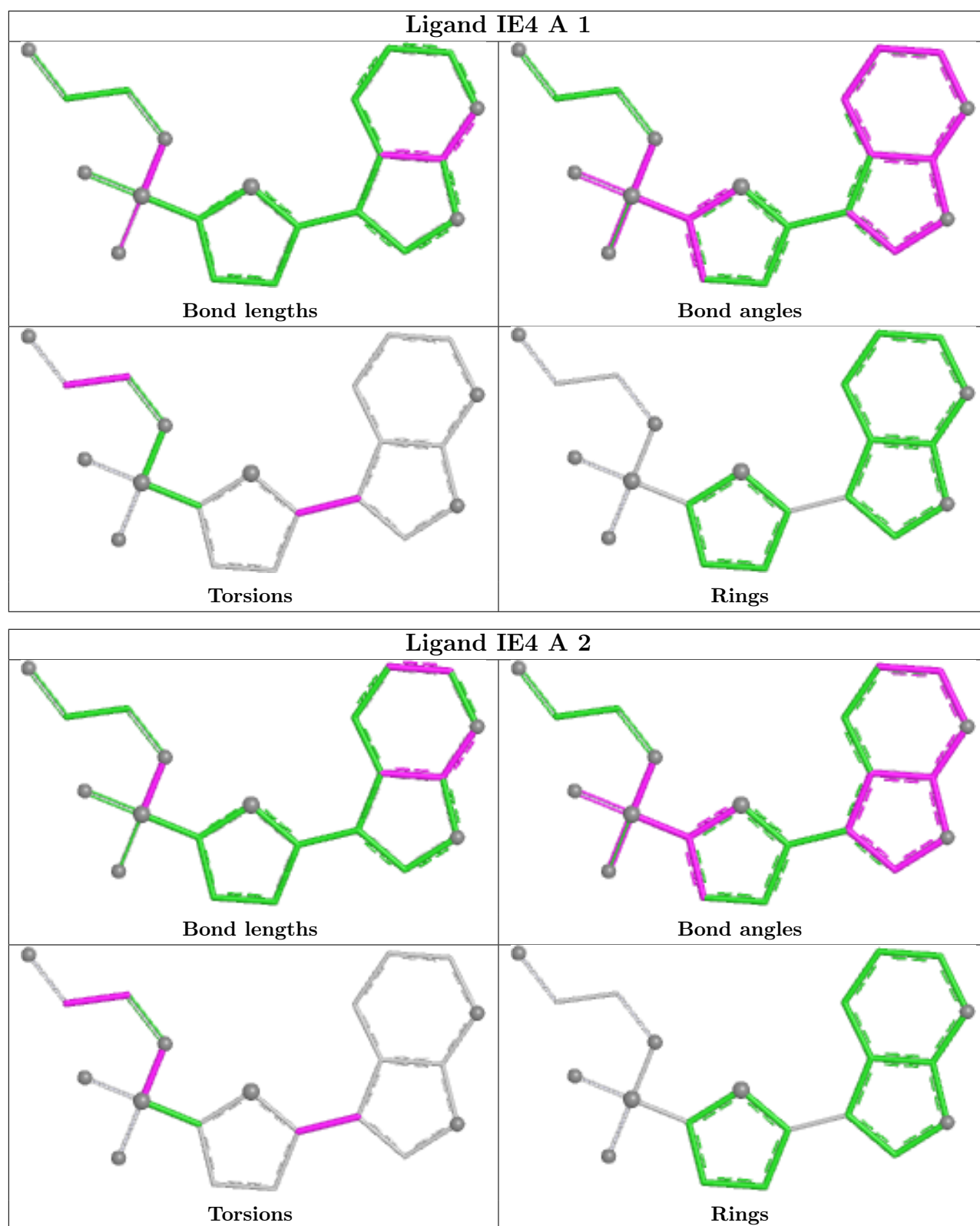
Mol	Chain	Res	Type	Atoms
2	A	1	IE4	C11-C12-C14-C22
2	A	1	IE4	S13-C12-C14-C22
2	A	1	IE4	N02-C03-C04-N05
2	A	2	IE4	C11-C12-C14-C22
2	A	2	IE4	N02-C03-C04-N05
3	B	1942	GOL	O1-C1-C2-O2
2	A	2	IE4	C04-N05-S06-O08
3	A	1941	GOL	C1-C2-C3-O3
3	B	1942	GOL	O1-C1-C2-C3
3	A	1941	GOL	O2-C2-C3-O3
2	A	2	IE4	C04-N05-S06-O07
2	A	2	IE4	C04-N05-S06-C09
2	A	2	IE4	S13-C12-C14-C22
3	A	1941	GOL	O1-C1-C2-C3
3	A	1941	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	IE4	1	0
3	A	1941	GOL	1	0
2	A	2	IE4	2	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 ⓘ

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	260/318 (81%)	1.11	38 (14%) 6 4	32, 52, 95, 130	0
1	B	259/318 (81%)	1.94	112 (43%) 0 0	40, 69, 109, 132	0
All	All	519/636 (81%)	1.52	150 (28%) 1 1	32, 59, 104, 132	0

All (150) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	940	VAL	5.2
1	B	858	TYR	5.2
1	A	670	LEU	5.1
1	B	670	LEU	4.8
1	B	940	VAL	4.6
1	A	829	LEU	4.6
1	A	897	PHE	4.6
1	A	723	HIS	4.6
1	B	855	PRO	4.3
1	B	896	MET	4.3
1	B	881	PHE	4.2
1	B	723	HIS	4.2
1	B	884	LEU	3.9
1	B	691	TYR	3.8
1	B	921	PRO	3.8
1	B	841	GLY	3.8
1	B	888	GLN	3.8
1	A	939	LYS	3.7
1	B	768	SER	3.7
1	B	900	HIS	3.4
1	B	714	ARG	3.4
1	B	853	LYS	3.4
1	B	824	GLY	3.4
1	B	889	ALA	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	899	VAL	3.3
1	B	880	PRO	3.3
1	B	680	ASN	3.3
1	B	753	PHE	3.3
1	A	827	LYS	3.3
1	B	712	PRO	3.3
1	B	769	LYS	3.3
1	B	690	THR	3.2
1	B	744	PHE	3.2
1	A	713	GLU	3.2
1	B	729	HIS	3.2
1	B	941	SER	3.2
1	B	722	LEU	3.1
1	B	866	SER	3.1
1	A	848	PRO	3.1
1	A	823	PHE	3.1
1	B	894	VAL	3.1
1	B	919	PHE	3.1
1	B	939	LYS	3.0
1	B	869	CYS	3.0
1	B	763	SER	3.0
1	B	821	SER	3.0
1	B	916	LEU	3.0
1	B	675	TYR	3.0
1	B	830	ALA	3.0
1	B	851	ILE	2.9
1	B	681	GLY	2.9
1	A	840	THR	2.9
1	B	825	THR	2.9
1	A	724	GLU	2.9
1	B	737	ILE	2.9
1	B	848	PRO	2.9
1	B	845	TYR	2.9
1	A	906	SER	2.8
1	A	931	ASP	2.8
1	B	800	VAL	2.8
1	A	824	GLY	2.8
1	A	831	GLY	2.8
1	B	748	GLY	2.8
1	B	749	PHE	2.8
1	B	843	LEU	2.7
1	B	854	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	882	TYR	2.7
1	B	750	ILE	2.7
1	B	799	ILE	2.7
1	A	896	MET	2.7
1	B	877	GLY	2.7
1	B	759	GLY	2.7
1	B	823	PHE	2.7
1	B	852	ASP	2.7
1	B	912	LYS	2.6
1	A	899	VAL	2.6
1	B	847	ALA	2.6
1	B	784	THR	2.6
1	B	802	ARG	2.6
1	A	909	ALA	2.6
1	B	703	GLN	2.5
1	B	764	ALA	2.5
1	A	722	LEU	2.5
1	B	774	LYS	2.5
1	B	840	THR	2.5
1	A	852	ASP	2.5
1	B	865	TRP	2.5
1	B	895	GLY	2.5
1	A	924	ASP	2.4
1	B	846	MET	2.4
1	B	684	VAL	2.4
1	B	676	GLU	2.4
1	B	719	SER	2.4
1	B	687	GLY	2.4
1	B	828	ARG	2.4
1	A	888	GLN	2.3
1	B	720	GLN	2.3
1	A	727	ALA	2.3
1	B	909	ALA	2.3
1	B	861	ALA	2.3
1	A	804	ILE	2.3
1	A	903	ILE	2.3
1	A	691	TYR	2.3
1	A	725	GLU	2.3
1	B	856	ARG	2.3
1	B	689	GLY	2.3
1	B	857	GLY	2.3
1	B	804	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	872	ILE	2.3
1	B	913	ALA	2.3
1	B	766	LEU	2.3
1	B	864	ILE	2.2
1	B	754	MET	2.2
1	B	886	GLU	2.2
1	B	700	LEU	2.2
1	B	727	ALA	2.2
1	B	682	ASP	2.2
1	B	938	LEU	2.2
1	B	730	LYS	2.2
1	A	890	ALA	2.2
1	B	776	ASN	2.2
1	A	728	LEU	2.2
1	A	916	LEU	2.2
1	B	692	GLY	2.2
1	B	738	VAL	2.2
1	A	750	ILE	2.2
1	B	731	HIS	2.2
1	B	862	ALA	2.2
1	A	875	ALA	2.1
1	B	672	GLU	2.1
1	B	679	GLU	2.1
1	B	873	GLU	2.1
1	B	827	LYS	2.1
1	B	903	ILE	2.1
1	B	883	GLU	2.1
1	A	858	TYR	2.1
1	B	677	TYR	2.1
1	B	762	LEU	2.1
1	A	841	GLY	2.1
1	B	760	GLY	2.1
1	B	915	ILE	2.1
1	B	796	ASP	2.1
1	A	671	LEU	2.1
1	B	713	GLU	2.0
1	B	707	ALA	2.0
1	A	768	SER	2.0
1	B	842	THR	2.0
1	B	773	LEU	2.0
1	A	882	TYR	2.0
1	B	673	TYR	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](#)

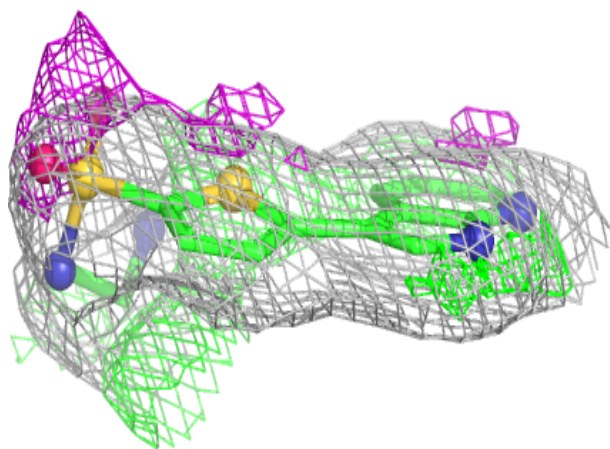
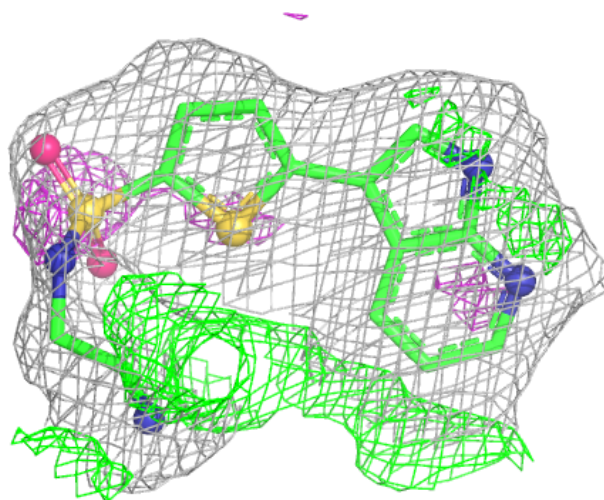
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

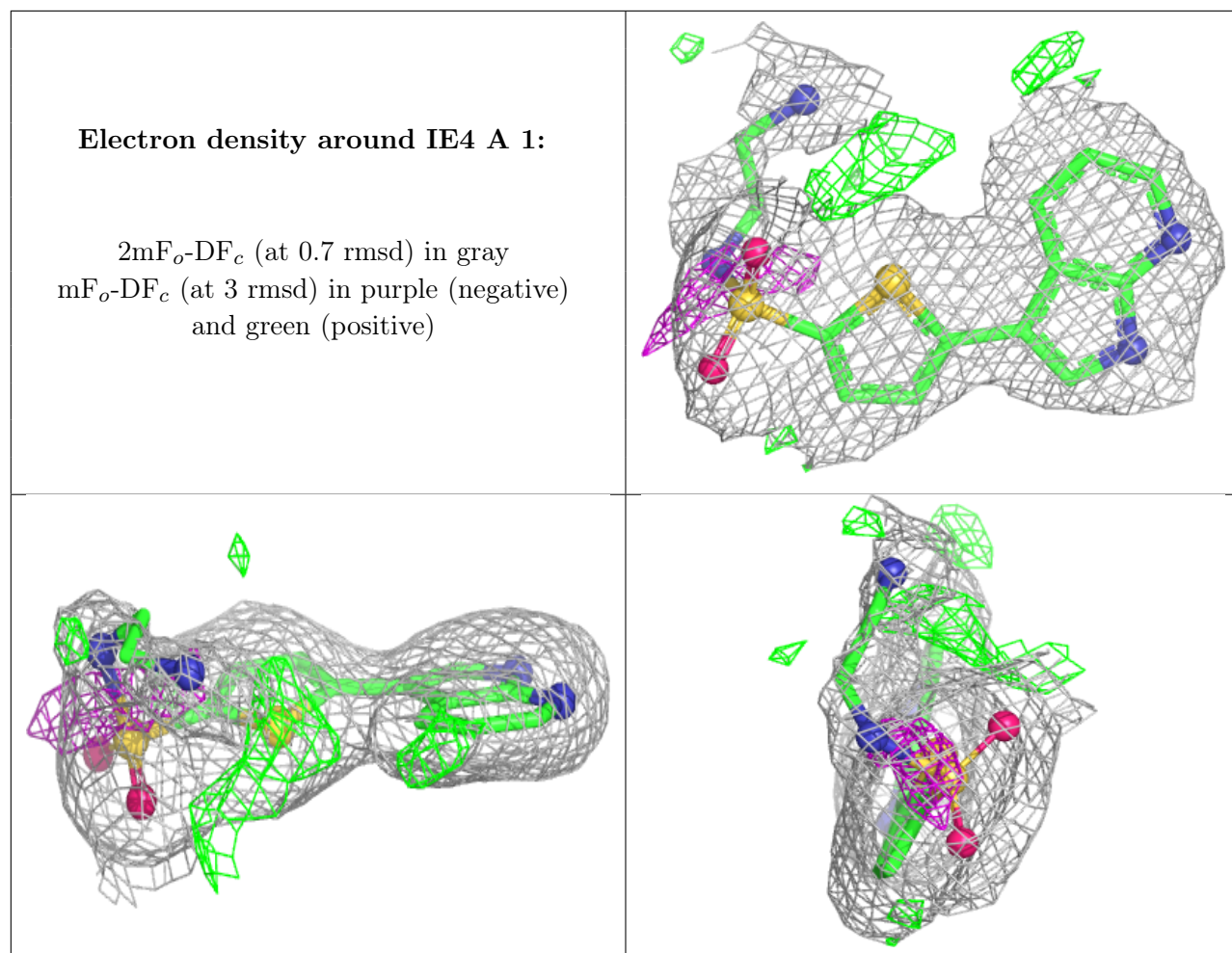
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	IE4	A	2	21/21	0.71	0.21	41,54,79,79	0
3	GOL	A	1941	6/6	0.87	0.14	53,53,54,55	0
3	GOL	B	1942	6/6	0.87	0.14	54,57,58,58	0
2	IE4	A	1	21/21	0.89	0.13	31,40,68,71	0

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 IE4 A 2:

$2mF_o-DF_c$ (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.