



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:50 AM UTC

PDB ID : 4ZJJ / pdb_00004zjj
Title : PAK1 in complex with (S)-N-(tert-butyl)-3-((2-chloro-5-ethyl-8-fluoro-dibenzodiazepin-11-yl)amino)pyrrolidine-1-carboxamide
Authors : Gutmann, S.; Rummel, G.
Deposited on : 2015-04-29
Resolution : 2.20 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

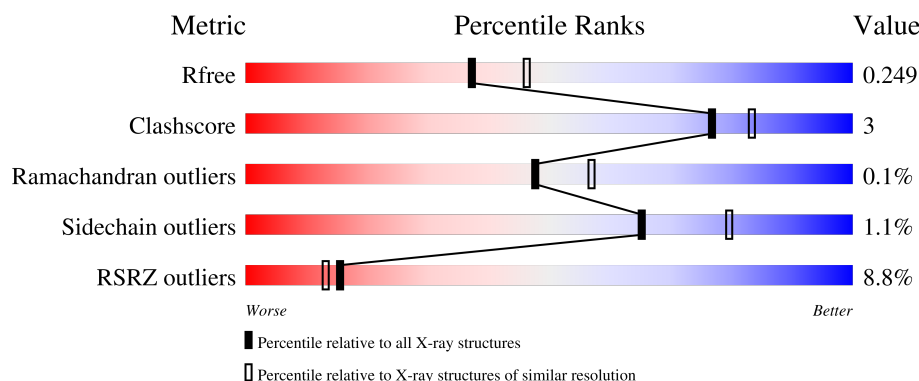
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	299	4% 87% 6% • 6%
1	B	299	7% 85% 6% • 7%
1	C	299	9% 82% 9% • 8%
1	D	299	11% 84% 7% 8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9072 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.

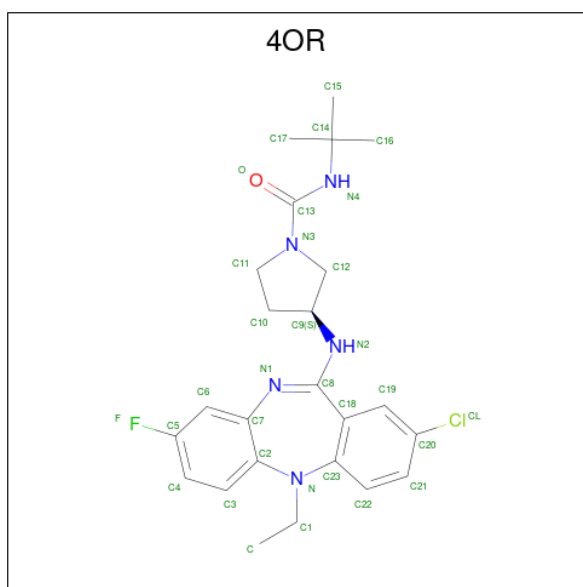
- Molecule 1 is a protein called Serine/threonine-protein kinase PAK 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	280	Total	C	N	O	S	0	1	0
			2126	1351	352	406	17			
1	B	277	Total	C	N	O	S	0	1	0
			2049	1306	335	391	17			
1	C	276	Total	C	N	O	S	0	0	0
			2012	1276	339	384	13			
1	D	274	Total	C	N	O	S	0	1	0
			1990	1261	338	376	15			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	247	GLY	-	expression tag	UNP Q13153
A	248	PRO	-	expression tag	UNP Q13153
B	247	GLY	-	expression tag	UNP Q13153
B	248	PRO	-	expression tag	UNP Q13153
C	247	GLY	-	expression tag	UNP Q13153
C	248	PRO	-	expression tag	UNP Q13153
D	247	GLY	-	expression tag	UNP Q13153
D	248	PRO	-	expression tag	UNP Q13153

- Molecule 2 is (S)-N-(tert-butyl)-3-((2-chloro-5-ethyl-8-fluoro-dibenzodiazepin-11-yl)amino)p yrrolidine-1-carboxamide (CCD ID: 4OR) (formula: C₂₄H₂₉ClFN₅O).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 32	C 24	Cl 1	F 1	N 5	O 1	0	0
2	B	1	Total 32	C 24	Cl 1	F 1	N 5	O 1	0	0
2	C	1	Total 32	C 24	Cl 1	F 1	N 5	O 1	0	0
2	D	1	Total 32	C 24	Cl 1	F 1	N 5	O 1	0	0

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	227	Total	O	0	0
			227	227		
4	B	145	Total	O	0	0
			145	145		
4	C	212	Total	O	0	0
			212	212		

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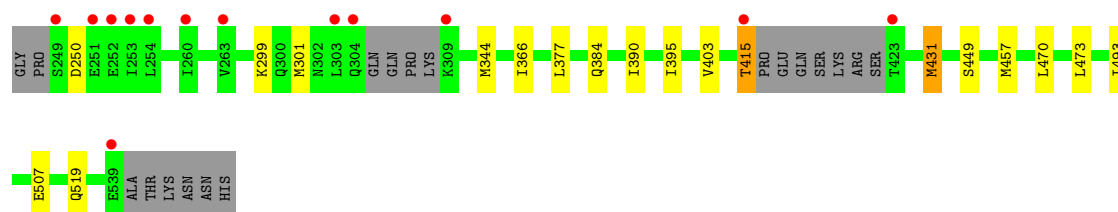
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	181	Total 181	O 181	0	0

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.

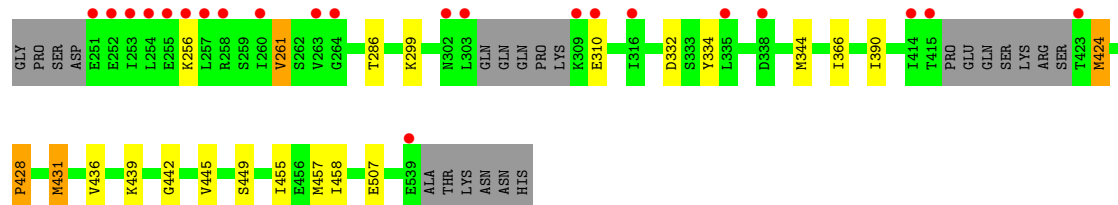
- Molecule 1: Serine/threonine-protein kinase PAK 1

Chain A: 



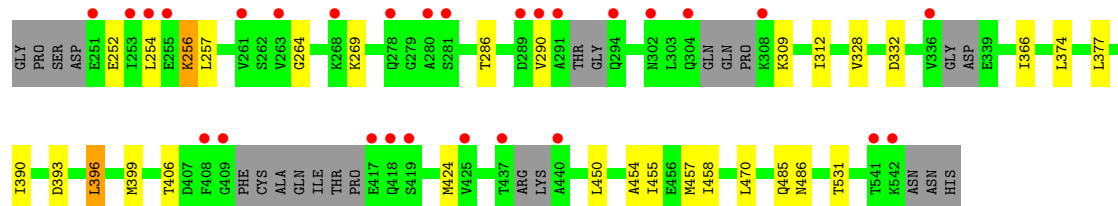
- Molecule 1: Serine/threonine-protein kinase PAK 1

Chain B: 



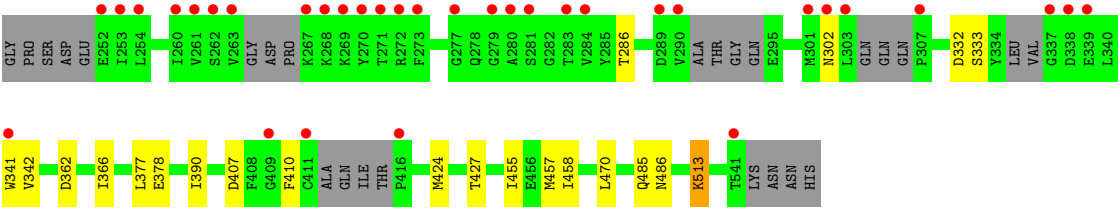
- Molecule 1: Serine/threonine-protein kinase PAK 1

Chain C: 



- Molecule 1: Serine/threonine-protein kinase PAK 1

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	110.33Å 113.79Å 116.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	116.36 – 2.20 116.36 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.9 (116.36-2.20) 99.9 (116.36-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.55 (at 2.20Å)	Xtriage
Refinement program	BUSTER 2.11.4	Depositor
R, R_{free}	0.201 , 0.238 0.208 , 0.249	Depositor DCC
R_{free} test set	3751 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	25.5	Xtriage
Anisotropy	0.689	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 69.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.019 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9072	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 38.73 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.5308e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: MG, 4OR

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	0.86	1/2160 (0.0%)	1.14	0/2929
1	B	0.85	2/2083 (0.1%)	1.21	3/2831 (0.1%)
1	C	0.85	0/2042	1.24	9/2774 (0.3%)
1	D	0.85	0/2021	1.22	4/2745 (0.1%)
All	All	0.85	3/8306 (0.0%)	1.20	16/11279 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	431	MET	SD-CE	-6.71	1.62	1.79
1	B	424	MET	SD-CE	-5.83	1.65	1.79
1	A	431	MET	SD-CE	-5.43	1.66	1.79

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	256	LYS	N-CA-C	-10.04	100.44	112.89
1	C	256	LYS	N-CA-C	-8.99	102.31	113.28
1	C	393	ASP	CA-CB-CG	5.86	118.46	112.60
1	D	362	ASP	CA-CB-CG	5.68	118.28	112.60
1	B	286	THR	N-CA-C	-5.44	101.41	110.17
1	D	302	ASN	CA-C-N	5.40	131.41	121.70
1	D	302	ASN	C-N-CA	5.40	131.41	121.70
1	C	286	THR	N-CA-C	-5.36	101.54	110.17
1	B	310	GLU	N-CA-C	-5.36	106.59	113.23
1	D	286	THR	N-CA-C	-5.16	101.86	110.17
1	C	256	LYS	CA-C-N	5.11	128.08	120.31
1	C	256	LYS	C-N-CA	5.11	128.08	120.31
1	C	252	GLU	CA-C-N	5.02	127.33	120.46
1	C	252	GLU	C-N-CA	5.02	127.33	120.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	257	LEU	CA-C-N	5.01	127.50	120.28
1	C	257	LEU	C-N-CA	5.01	127.50	120.28

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	2126	0	2083	13	0
1	B	2049	0	1946	12	0
1	C	2012	0	1881	16	0
1	D	1990	0	1844	13	0
2	A	32	0	29	1	0
2	B	32	0	29	0	0
2	C	32	0	29	1	0
2	D	32	0	29	1	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	227	0	0	0	0
4	B	145	0	0	0	0
4	C	212	0	0	1	0
4	D	181	0	0	0	0
All	All	9072	0	7870	47	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 3.

All (47) 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:473:LEU:HD11	1:D:424:MET:HE1	1.73	0.71
1:B:261:VAL:HG13	1:B:334:TYR:HA	1.84	0.59
1:A:377:LEU:HD21	1:A:390:ILE:HD11	1.86	0.58
1:A:431:MET:HE1	1:D:424:MET:HE2	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:377:LEU:HD21	1:D:390:ILE:HD11	1.88	0.55
1:C:396:LEU:HD22	1:C:406:THR:HG21	1.88	0.55
1:A:301:MET:HE1	2:A:601:4OR:C17	2.36	0.54
1:A:470:LEU:HD11	1:D:427:THR:HG21	1.90	0.53
1:C:377:LEU:HD21	1:C:390:ILE:HD11	1.89	0.53
1:C:399:MET:HE3	4:C:720:HOH:O	2.08	0.52
1:B:424:MET:HE3	1:B:436:VAL:HA	1.93	0.51
1:D:407:ASP:OD2	1:D:410:PHE:HB2	2.11	0.50
1:D:485:GLN:HE21	1:D:486:ASN:HD21	1.58	0.49
1:C:254:LEU:C	1:C:256:LYS:H	2.21	0.49
1:B:366:ILE:HG12	1:B:457:MET:HB3	1.94	0.48
1:A:384:GLN:HB3	1:A:415:THR:HG22	1.96	0.48
1:A:299:LYS:HD2	1:A:344:MET:HE1	1.94	0.48
1:D:378:GLU:OE2	1:D:513:LYS:HG2	2.14	0.48
1:C:309:LYS:HA	1:C:312:ILE:HD12	1.96	0.47
1:A:390:ILE:HB	1:A:449:SER:HB2	1.96	0.46
1:A:366:ILE:HG12	1:A:457:MET:HB3	1.97	0.46
1:A:493:ILE:HD11	1:A:519:GLN:HB2	1.97	0.46
1:C:485:GLN:HE21	1:C:486:ASN:HD21	1.63	0.45
1:A:507:GLU:HB2	1:B:507:GLU:HB2	1.98	0.45
1:B:299:LYS:HD2	1:B:344:MET:HE1	1.98	0.45
1:C:455:ILE:HA	1:C:458:ILE:HD12	1.99	0.44
1:D:455:ILE:HA	1:D:458:ILE:HD12	2.00	0.44
1:C:470:LEU:HD22	1:D:470:LEU:HD22	1.99	0.43
1:B:424:MET:HE2	1:B:431:MET:SD	2.58	0.43
1:B:442:GLY:O	1:B:445:VAL:HG22	2.18	0.43
1:B:390:ILE:HB	1:B:449:SER:HB2	2.00	0.43
1:B:428:PRO:HD3	1:C:424:MET:HB3	2.00	0.43
1:A:377:LEU:HD21	1:A:390:ILE:CD1	2.48	0.43
1:C:374:LEU:HD21	1:C:450:LEU:HD22	2.01	0.42
1:B:431:MET:HE1	1:C:424:MET:SD	2.59	0.42
1:C:264:GLY:HA3	1:C:269:LYS:CE	2.49	0.42
1:C:454:ALA:HA	1:C:457:MET:HE3	2.01	0.42
1:A:395:ILE:HG23	1:A:403:VAL:HG13	2.01	0.42
1:D:333:SER:HA	1:D:341:TRP:O	2.19	0.42
1:D:342:VAL:HG21	2:D:601:4OR:H16	2.01	0.41
1:C:328:VAL:O	2:C:601:4OR:H3	2.19	0.41
1:B:424:MET:HB2	1:B:439:LYS:HE2	2.02	0.41
1:D:377:LEU:HD21	1:D:390:ILE:CD1	2.48	0.41
1:C:366:ILE:HG12	1:C:457:MET:HB3	2.01	0.41
1:C:399:MET:O	1:C:531:THR:HG22	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:366:ILE:HG12	1:D:457:MET:HB3	2.02	0.41
1:B:455:ILE:HA	1:B:458:ILE:HD12	2.02	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	275/299 (92%)	272 (99%)	3 (1%)	0	100	100
1	B	272/299 (91%)	266 (98%)	6 (2%)	0	100	100
1	C	264/299 (88%)	258 (98%)	5 (2%)	1 (0%)	30	34
1	D	263/299 (88%)	259 (98%)	4 (2%)	0	100	100
All	All	1074/1196 (90%)	1055 (98%)	18 (2%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	290	VAL

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	223/259 (86%)	221 (99%)	2 (1%)	70	84

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	204/259 (79%)	201 (98%)	3 (2%)	57	73
1	C	194/259 (75%)	192 (99%)	2 (1%)	68	81
1	D	190/259 (73%)	188 (99%)	2 (1%)	65	79
All	All	811/1036 (78%)	802 (99%)	9 (1%)	65	79

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	250	ASP
1	A	415	THR
1	B	261	VAL
1	B	332	ASP
1	B	428	PRO
1	C	332	ASP
1	C	396	LEU
1	D	332	ASP
1	D	513	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	314	ASN
1	A	326	ASN
1	A	375	GLN
1	A	485	GLN
1	A	519	GLN
1	B	294	GLN
1	B	326	ASN
1	B	375	GLN
1	C	326	ASN
1	C	375	GLN
1	C	479	ASN
1	C	485	GLN
1	D	326	ASN
1	D	375	GLN
1	D	485	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

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	4OR	A	601	-	34,35,35	0.62	1 (2%)	44,52,52	0.79	1 (2%)
2	4OR	D	601	-	34,35,35	0.87	2 (5%)	44,52,52	0.92	1 (2%)
2	4OR	C	601	-	34,35,35	0.83	2 (5%)	44,52,52	0.64	1 (2%)
2	4OR	B	601	-	34,35,35	0.72	1 (2%)	44,52,52	0.87	2 (4%)

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	4OR	A	601	-	-	1/13/40/40	0/4/4/4
2	4OR	D	601	-	-	2/13/40/40	0/3/4/4
2	4OR	C	601	-	-	1/13/40/40	0/4/4/4
2	4OR	B	601	-	-	1/13/40/40	0/4/4/4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	601	4OR	C13-N4	4.15	1.44	1.37
2	B	601	4OR	C13-N4	3.94	1.44	1.37
2	C	601	4OR	C13-N4	3.48	1.43	1.37
2	A	601	4OR	C13-N4	3.16	1.43	1.37
2	C	601	4OR	C12-C9	2.98	1.58	1.53
2	D	601	4OR	C12-C9	2.44	1.57	1.53

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	601	4OR	C9-N2-C8	4.81	127.66	122.29
2	B	601	4OR	C9-N2-C8	4.42	127.22	122.29
2	A	601	4OR	C9-N2-C8	4.39	127.19	122.29
2	C	601	4OR	C14-N4-C13	2.74	129.47	125.15
2	B	601	4OR	C14-N4-C13	2.68	129.38	125.15

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	601	4OR	C-C1-N-C23
2	D	601	4OR	C-C1-N-C23
2	B	601	4OR	C12-C9-N2-C8
2	D	601	4OR	C12-C9-N2-C8
2	A	601	4OR	C-C1-N-C23

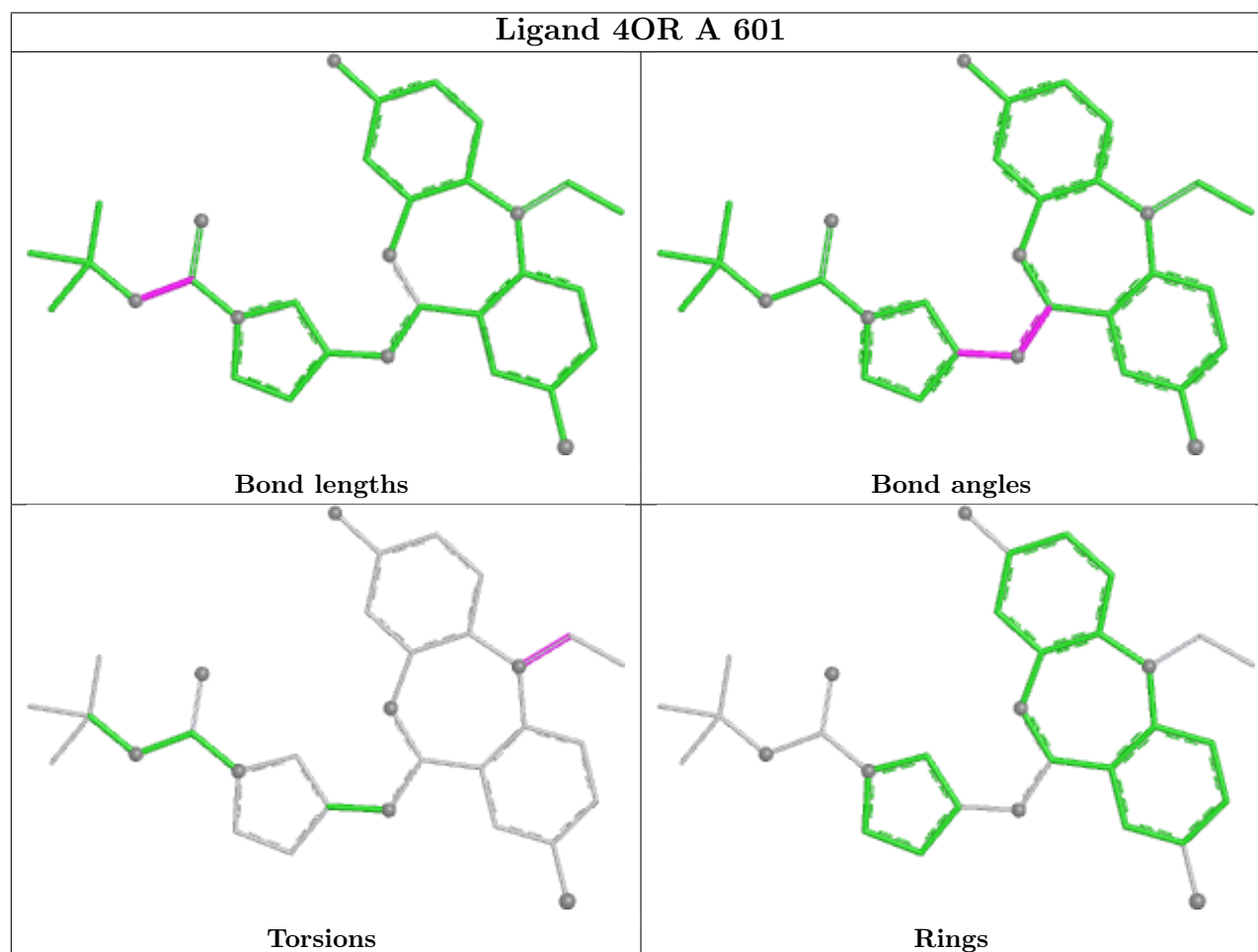
There are no ring outliers.

3 monomers are involved in 3 short contacts:

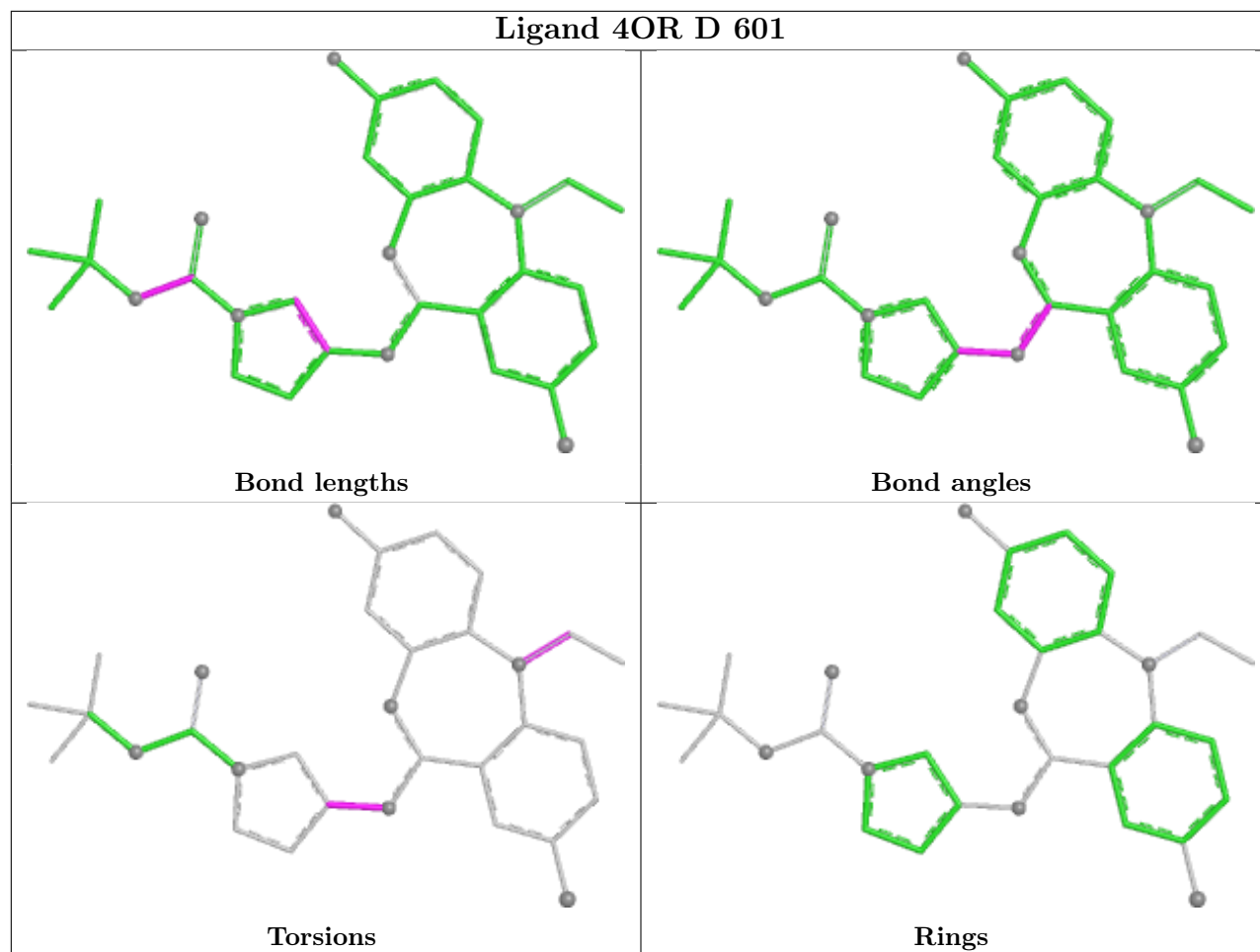
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	4OR	1	0
2	D	601	4OR	1	0
2	C	601	4OR	1	0

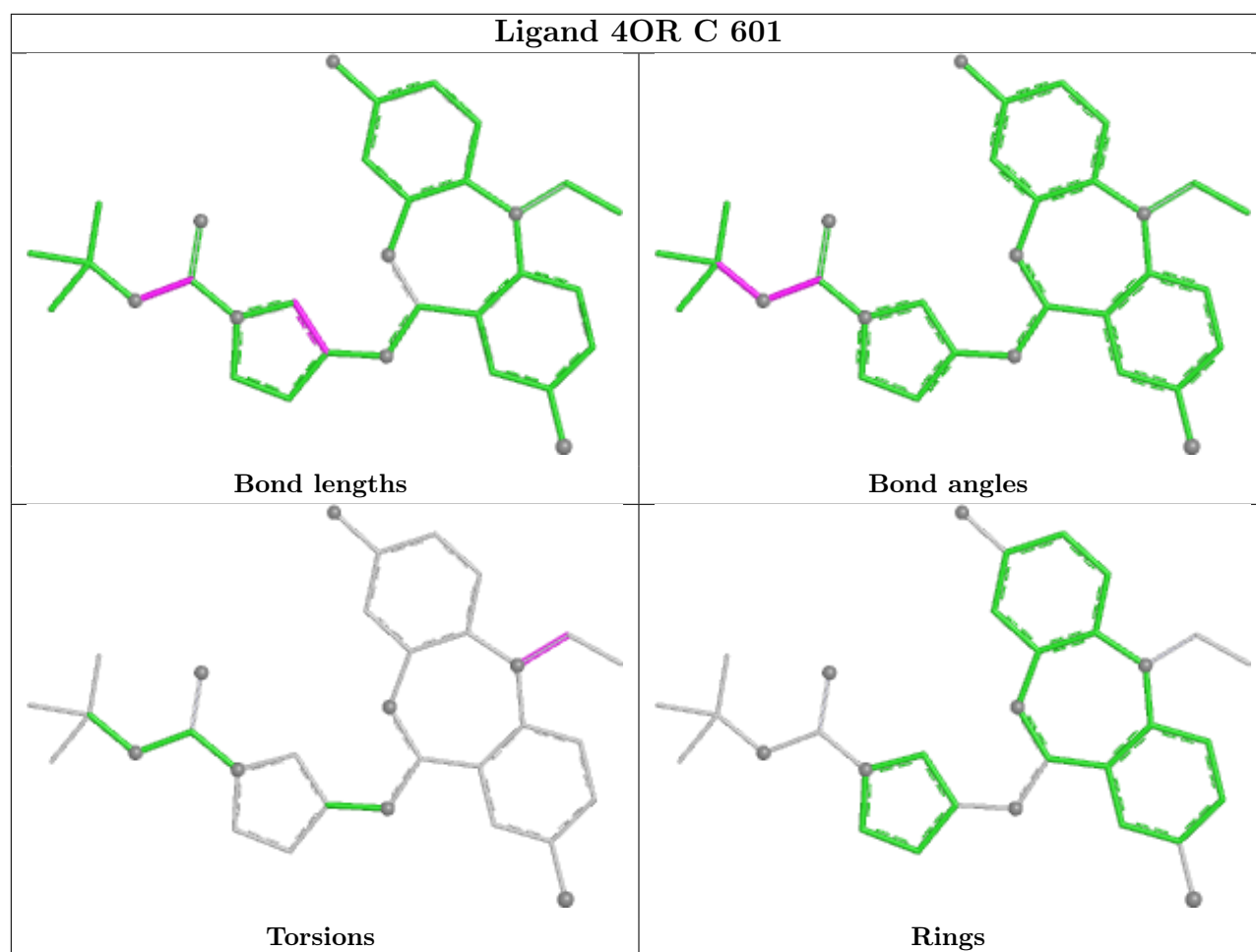
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

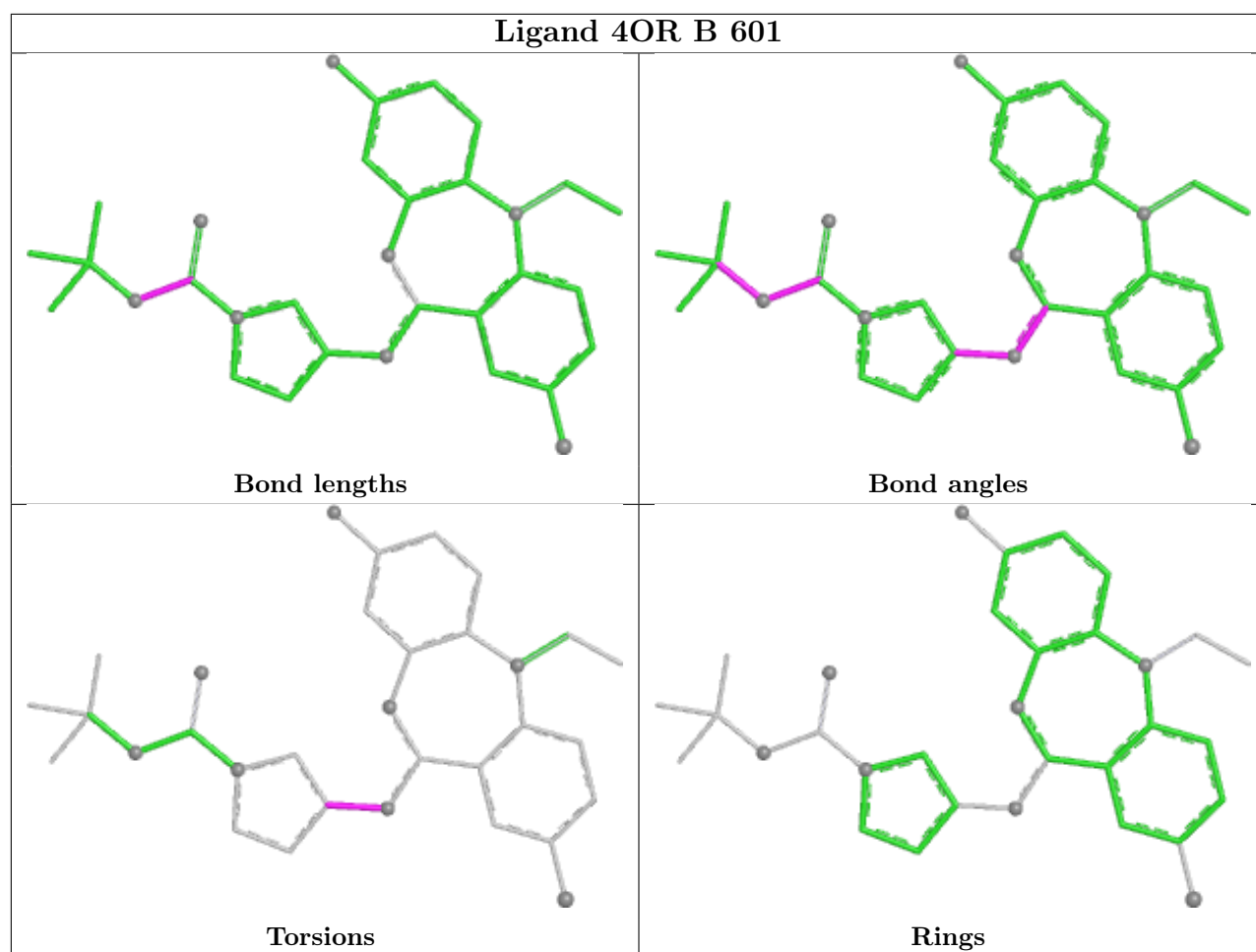
highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified 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 4OR D 601







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	280/299 (93%)	-0.09	13 (4%)	37 34	10, 23, 51, 86	1 (0%)
1	B	277/299 (92%)	0.34	22 (7%)	18 16	13, 35, 70, 103	1 (0%)
1	C	276/299 (92%)	0.32	28 (10%)	12 10	15, 30, 68, 98	0
1	D	274/299 (91%)	0.45	34 (12%)	8 6	12, 30, 64, 90	1 (0%)
All	All	1107/1196 (92%)	0.25	97 (8%)	15 13	10, 30, 65, 103	3 (0%)

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	290	VAL	8.6
1	C	291	ALA	8.0
1	C	336	VAL	7.5
1	D	303	LEU	7.4
1	B	415	THR	6.9
1	B	303	LEU	6.5
1	A	415	THR	6.1
1	C	409	GLY	6.1
1	D	263	VAL	5.8
1	D	252	GLU	5.6
1	B	253	ILE	5.5
1	C	290	VAL	5.4
1	B	254	LEU	5.3
1	D	341	TRP	5.1
1	D	277	GLY	4.9
1	A	539	GLU	4.8
1	D	411	CYS	4.8
1	C	304	GLN	4.7
1	D	339	GLU	4.4
1	A	303	LEU	4.3
1	C	417	GLU	4.2

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Mol	Chain	Res	Type	RSRZ
1	C	418	GLN	4.2
1	B	251	GLU	4.2
1	D	253	ILE	4.1
1	D	270	TYR	4.0
1	D	302	ASN	4.0
1	C	294	GLN	3.8
1	C	253	ILE	3.6
1	B	539	GLU	3.6
1	A	249	SER	3.5
1	B	309	LYS	3.5
1	A	423	THR	3.5
1	C	254	LEU	3.5
1	D	271	THR	3.4
1	D	337	GLY	3.4
1	B	423	THR	3.4
1	B	338	ASP	3.2
1	D	283	THR	3.2
1	D	281	SER	3.2
1	D	269	LYS	3.1
1	B	263	VAL	3.1
1	A	304	GLN	3.1
1	C	255	GLU	3.0
1	D	338	ASP	3.0
1	A	253	ILE	3.0
1	D	416	PRO	2.9
1	D	301	MET	2.9
1	C	302	ASN	2.9
1	D	267	LYS	2.8
1	A	260	ILE	2.8
1	C	419	SER	2.8
1	C	278	GLN	2.8
1	A	254	LEU	2.8
1	D	289	ASP	2.8
1	D	284	VAL	2.8
1	B	255	GLU	2.7
1	D	268	LYS	2.7
1	D	307	PRO	2.7
1	C	251	GLU	2.7
1	C	280	ALA	2.7
1	B	252	GLU	2.7
1	C	437	THR	2.6
1	B	260	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	262	SER	2.6
1	B	257	LEU	2.5
1	D	279	GLY	2.5
1	D	261	VAL	2.5
1	B	414	ILE	2.5
1	B	264	GLY	2.5
1	C	408	PHE	2.4
1	C	425	VAL	2.4
1	B	310	GLU	2.4
1	C	440	ALA	2.4
1	D	272	ARG	2.4
1	C	268	LYS	2.3
1	C	308	LYS	2.3
1	C	263	VAL	2.3
1	D	260	ILE	2.3
1	A	309	LYS	2.3
1	D	254	LEU	2.3
1	D	541	THR	2.3
1	B	258	ARG	2.3
1	C	289	ASP	2.3
1	A	263	VAL	2.2
1	B	302	ASN	2.2
1	D	273	PHE	2.2
1	C	281	SER	2.2
1	C	261	VAL	2.1
1	C	542	LYS	2.1
1	B	316	ILE	2.1
1	A	252	GLU	2.1
1	D	409	GLY	2.1
1	C	541	THR	2.1
1	A	251	GLU	2.1
1	D	280	ALA	2.1
1	B	335	LEU	2.1
1	B	256	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

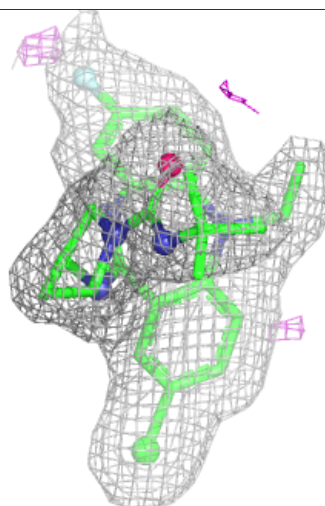
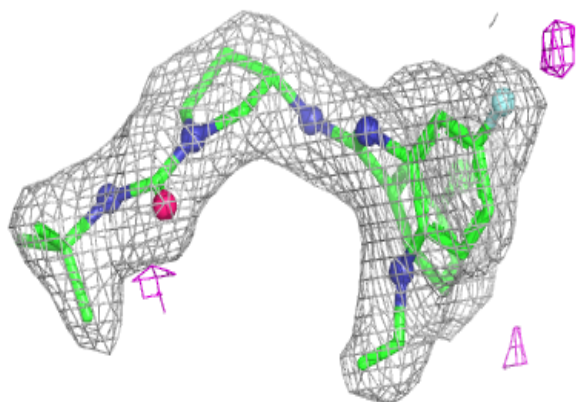
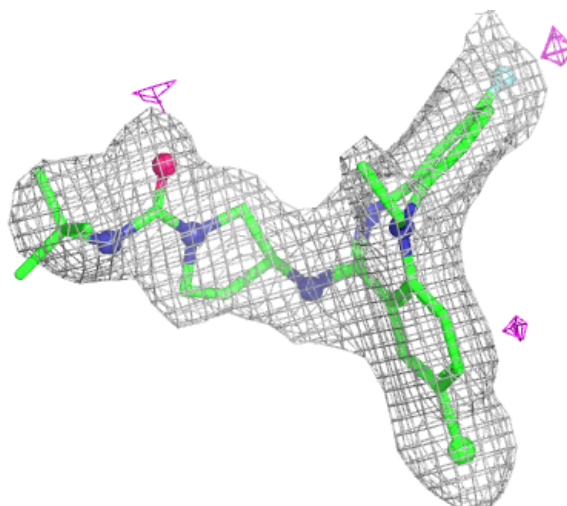
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	MG	D	602	1/1	0.92	0.08	42,42,42,42	0
2	4OR	C	601	32/32	0.94	0.09	20,33,46,48	0
2	4OR	B	601	32/32	0.95	0.08	28,34,39,42	0
2	4OR	D	601	32/32	0.97	0.06	19,23,39,41	0
3	MG	C	602	1/1	0.97	0.04	36,36,36,36	0
2	4OR	A	601	32/32	0.97	0.07	15,21,28,30	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.

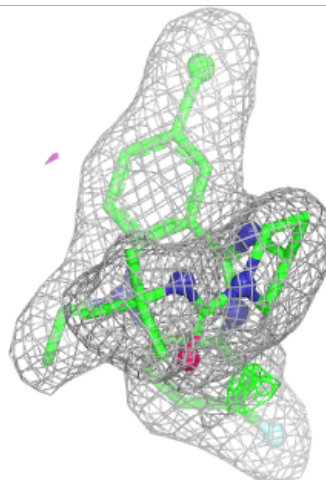
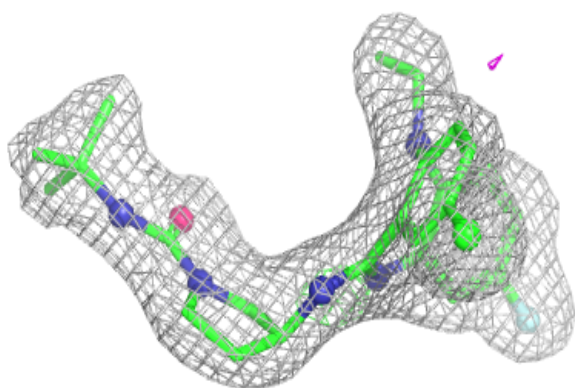
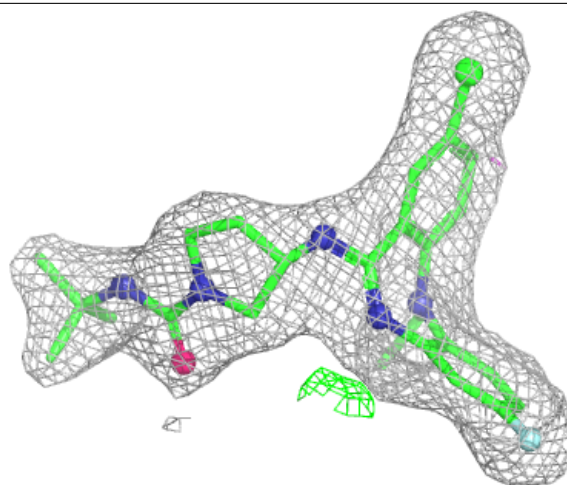
Electron density around 4OR C 601:

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



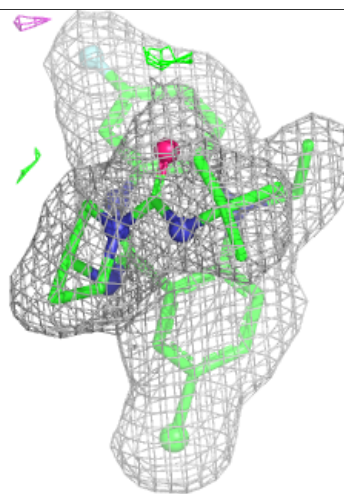
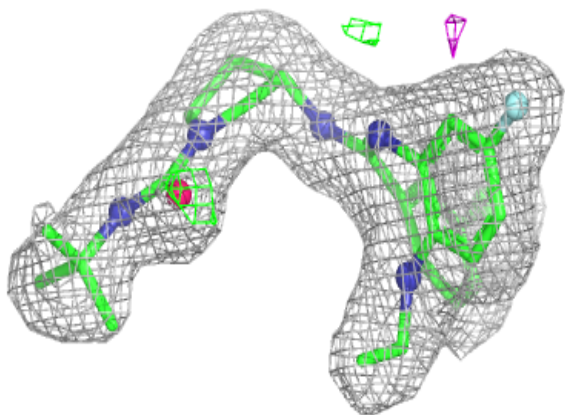
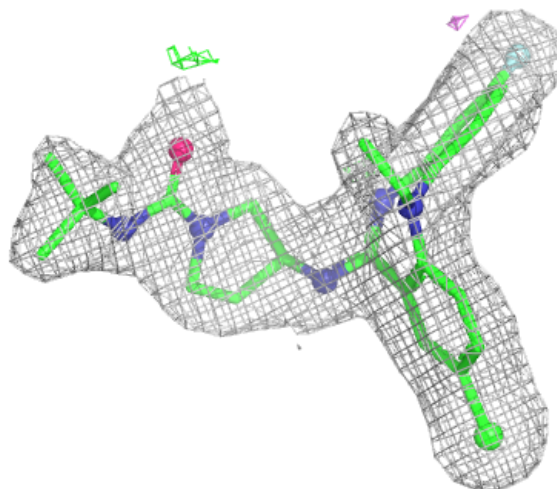
Electron density around 4OR B 601:

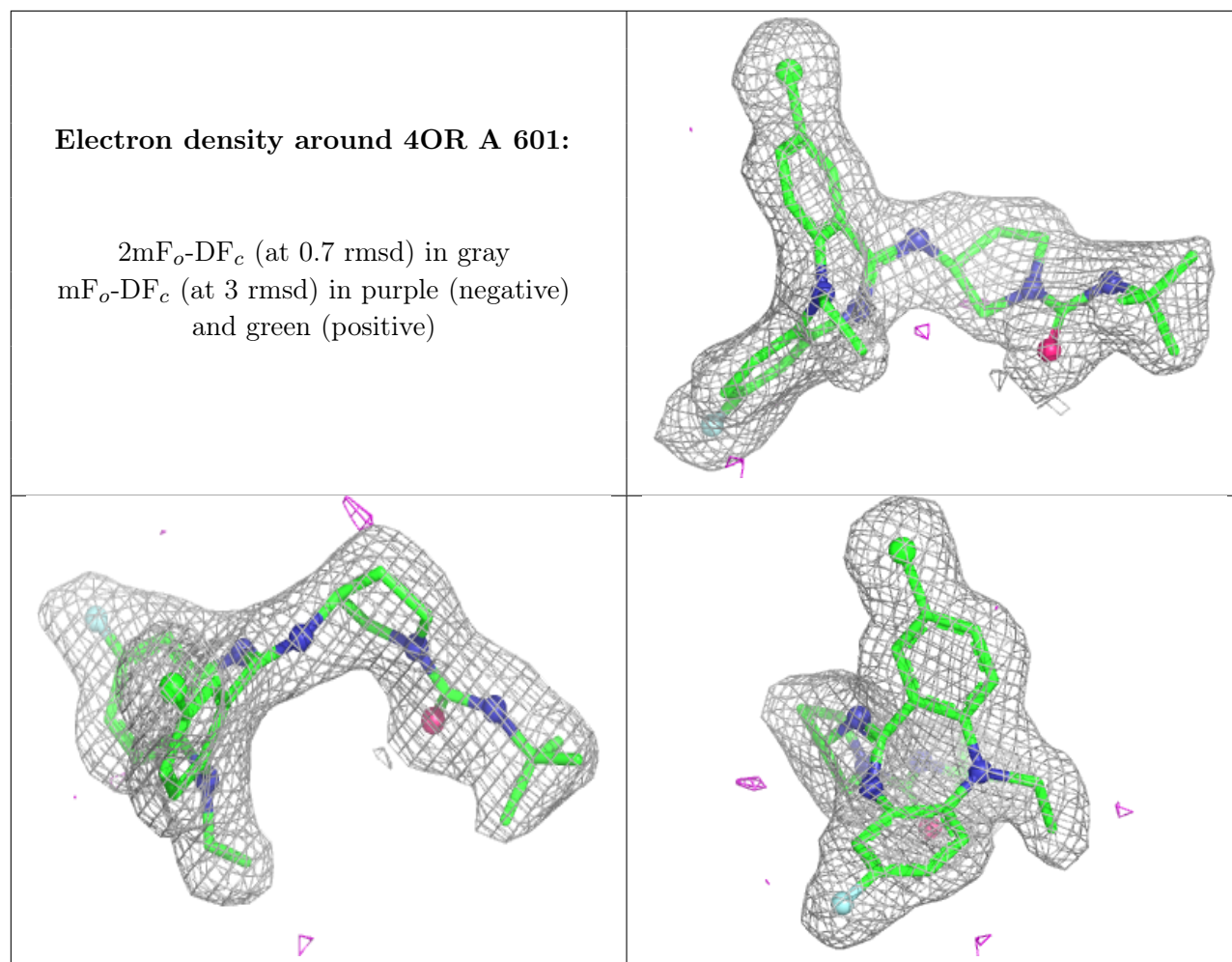
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 4OR D 601:

$2mF_o-DF_c$ (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.