



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 01:11 am BST

PDB ID : 5JYO
Title : Allosteric inhibition of Kidney Isoform of Glutaminase
Authors : Sivaraman, J.; Jayaraman, S.
Deposited on : 2016-05-15
Resolution : 2.10 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

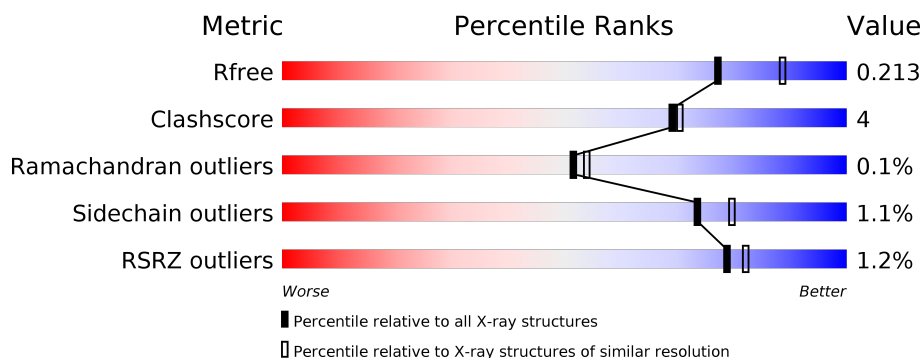
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 on 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	333	<div> <div>2%</div> <div> <div></div> <div>86%</div> <div>8%</div> <div>6%</div> </div> </div>
1	B	333	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>8%</div> <div>7%</div> </div> </div>
1	C	333	<div> <div>%</div> <div> <div></div> <div>83%</div> <div>11%</div> <div>7%</div> </div> </div>
1	D	333	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>7%</div> <div>7%</div> </div> </div>
1	E	333	<div> <div>2%</div> <div> <div></div> <div>86%</div> <div>7%</div> <div>7%</div> </div> </div>
1	F	333	<div> <div>%</div> <div> <div></div> <div>85%</div> <div>8%</div> <div>7%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	333	<div><div>%</div><div><div></div><div>84%</div><div>10%</div><div>7%</div></div></div>
1	H	333	<div><div>2%</div><div><div></div><div>89%</div><div>5%</div><div>7%</div></div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 20485 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 Glutaminase kidney isoform, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	311	Total	C	N	O	S	0	0	0
			2385	1517	394	450	24			
1	A	312	Total	C	N	O	S	0	0	0
			2393	1523	395	451	24			
1	B	311	Total	C	N	O	S	0	0	0
			2385	1517	394	450	24			
1	C	311	Total	C	N	O	S	0	0	0
			2385	1519	393	449	24			
1	E	311	Total	C	N	O	S	0	0	0
			2385	1517	394	450	24			
1	F	311	Total	C	N	O	S	0	0	0
			2385	1517	394	450	24			
1	G	311	Total	C	N	O	S	0	0	0
			2385	1517	394	450	24			
1	H	311	Total	C	N	O	S	0	0	0
			2385	1517	394	450	24			

There are 160 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	201	MET	-	expression tag	UNP O94925
D	202	GLY	-	expression tag	UNP O94925
D	203	SER	-	expression tag	UNP O94925
D	204	SER	-	expression tag	UNP O94925
D	205	HIS	-	expression tag	UNP O94925
D	206	HIS	-	expression tag	UNP O94925
D	207	HIS	-	expression tag	UNP O94925
D	208	HIS	-	expression tag	UNP O94925
D	209	HIS	-	expression tag	UNP O94925
D	210	HIS	-	expression tag	UNP O94925
D	211	SER	-	expression tag	UNP O94925
D	212	SER	-	expression tag	UNP O94925
D	213	GLY	-	expression tag	UNP O94925

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Chain	Residue	Modelled	Actual	Comment	Reference
D	214	LEU	-	expression tag	UNP O94925
D	215	VAL	-	expression tag	UNP O94925
D	216	PRO	-	expression tag	UNP O94925
D	217	ARG	-	expression tag	UNP O94925
D	218	GLY	-	expression tag	UNP O94925
D	219	SER	-	expression tag	UNP O94925
D	220	MET	-	expression tag	UNP O94925
A	201	MET	-	expression tag	UNP O94925
A	202	GLY	-	expression tag	UNP O94925
A	203	SER	-	expression tag	UNP O94925
A	204	SER	-	expression tag	UNP O94925
A	205	HIS	-	expression tag	UNP O94925
A	206	HIS	-	expression tag	UNP O94925
A	207	HIS	-	expression tag	UNP O94925
A	208	HIS	-	expression tag	UNP O94925
A	209	HIS	-	expression tag	UNP O94925
A	210	HIS	-	expression tag	UNP O94925
A	211	SER	-	expression tag	UNP O94925
A	212	SER	-	expression tag	UNP O94925
A	213	GLY	-	expression tag	UNP O94925
A	214	LEU	-	expression tag	UNP O94925
A	215	VAL	-	expression tag	UNP O94925
A	216	PRO	-	expression tag	UNP O94925
A	217	ARG	-	expression tag	UNP O94925
A	218	GLY	-	expression tag	UNP O94925
A	219	SER	-	expression tag	UNP O94925
A	220	MET	-	expression tag	UNP O94925
B	201	MET	-	expression tag	UNP O94925
B	202	GLY	-	expression tag	UNP O94925
B	203	SER	-	expression tag	UNP O94925
B	204	SER	-	expression tag	UNP O94925
B	205	HIS	-	expression tag	UNP O94925
B	206	HIS	-	expression tag	UNP O94925
B	207	HIS	-	expression tag	UNP O94925
B	208	HIS	-	expression tag	UNP O94925
B	209	HIS	-	expression tag	UNP O94925
B	210	HIS	-	expression tag	UNP O94925
B	211	SER	-	expression tag	UNP O94925
B	212	SER	-	expression tag	UNP O94925
B	213	GLY	-	expression tag	UNP O94925
B	214	LEU	-	expression tag	UNP O94925
B	215	VAL	-	expression tag	UNP O94925

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Chain	Residue	Modelled	Actual	Comment	Reference
B	216	PRO	-	expression tag	UNP O94925
B	217	ARG	-	expression tag	UNP O94925
B	218	GLY	-	expression tag	UNP O94925
B	219	SER	-	expression tag	UNP O94925
B	220	MET	-	expression tag	UNP O94925
C	201	MET	-	expression tag	UNP O94925
C	202	GLY	-	expression tag	UNP O94925
C	203	SER	-	expression tag	UNP O94925
C	204	SER	-	expression tag	UNP O94925
C	205	HIS	-	expression tag	UNP O94925
C	206	HIS	-	expression tag	UNP O94925
C	207	HIS	-	expression tag	UNP O94925
C	208	HIS	-	expression tag	UNP O94925
C	209	HIS	-	expression tag	UNP O94925
C	210	HIS	-	expression tag	UNP O94925
C	211	SER	-	expression tag	UNP O94925
C	212	SER	-	expression tag	UNP O94925
C	213	GLY	-	expression tag	UNP O94925
C	214	LEU	-	expression tag	UNP O94925
C	215	VAL	-	expression tag	UNP O94925
C	216	PRO	-	expression tag	UNP O94925
C	217	ARG	-	expression tag	UNP O94925
C	218	GLY	-	expression tag	UNP O94925
C	219	SER	-	expression tag	UNP O94925
C	220	MET	-	expression tag	UNP O94925
E	201	MET	-	expression tag	UNP O94925
E	202	GLY	-	expression tag	UNP O94925
E	203	SER	-	expression tag	UNP O94925
E	204	SER	-	expression tag	UNP O94925
E	205	HIS	-	expression tag	UNP O94925
E	206	HIS	-	expression tag	UNP O94925
E	207	HIS	-	expression tag	UNP O94925
E	208	HIS	-	expression tag	UNP O94925
E	209	HIS	-	expression tag	UNP O94925
E	210	HIS	-	expression tag	UNP O94925
E	211	SER	-	expression tag	UNP O94925
E	212	SER	-	expression tag	UNP O94925
E	213	GLY	-	expression tag	UNP O94925
E	214	LEU	-	expression tag	UNP O94925
E	215	VAL	-	expression tag	UNP O94925
E	216	PRO	-	expression tag	UNP O94925
E	217	ARG	-	expression tag	UNP O94925

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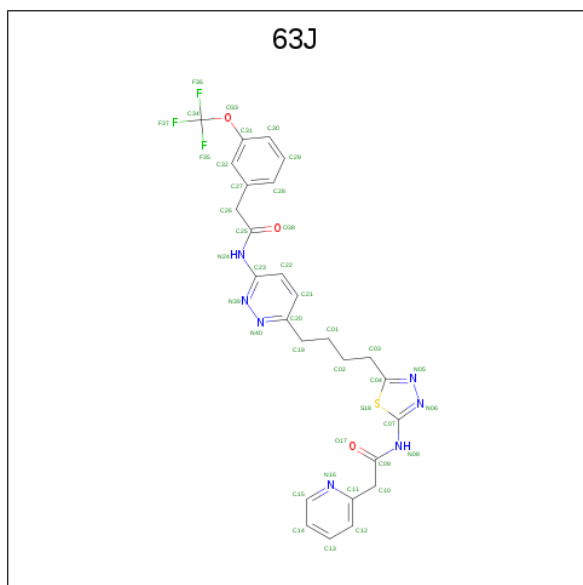
Chain	Residue	Modelled	Actual	Comment	Reference
E	218	GLY	-	expression tag	UNP O94925
E	219	SER	-	expression tag	UNP O94925
E	220	MET	-	expression tag	UNP O94925
F	201	MET	-	expression tag	UNP O94925
F	202	GLY	-	expression tag	UNP O94925
F	203	SER	-	expression tag	UNP O94925
F	204	SER	-	expression tag	UNP O94925
F	205	HIS	-	expression tag	UNP O94925
F	206	HIS	-	expression tag	UNP O94925
F	207	HIS	-	expression tag	UNP O94925
F	208	HIS	-	expression tag	UNP O94925
F	209	HIS	-	expression tag	UNP O94925
F	210	HIS	-	expression tag	UNP O94925
F	211	SER	-	expression tag	UNP O94925
F	212	SER	-	expression tag	UNP O94925
F	213	GLY	-	expression tag	UNP O94925
F	214	LEU	-	expression tag	UNP O94925
F	215	VAL	-	expression tag	UNP O94925
F	216	PRO	-	expression tag	UNP O94925
F	217	ARG	-	expression tag	UNP O94925
F	218	GLY	-	expression tag	UNP O94925
F	219	SER	-	expression tag	UNP O94925
F	220	MET	-	expression tag	UNP O94925
G	201	MET	-	expression tag	UNP O94925
G	202	GLY	-	expression tag	UNP O94925
G	203	SER	-	expression tag	UNP O94925
G	204	SER	-	expression tag	UNP O94925
G	205	HIS	-	expression tag	UNP O94925
G	206	HIS	-	expression tag	UNP O94925
G	207	HIS	-	expression tag	UNP O94925
G	208	HIS	-	expression tag	UNP O94925
G	209	HIS	-	expression tag	UNP O94925
G	210	HIS	-	expression tag	UNP O94925
G	211	SER	-	expression tag	UNP O94925
G	212	SER	-	expression tag	UNP O94925
G	213	GLY	-	expression tag	UNP O94925
G	214	LEU	-	expression tag	UNP O94925
G	215	VAL	-	expression tag	UNP O94925
G	216	PRO	-	expression tag	UNP O94925
G	217	ARG	-	expression tag	UNP O94925
G	218	GLY	-	expression tag	UNP O94925
G	219	SER	-	expression tag	UNP O94925

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Chain	Residue	Modelled	Actual	Comment	Reference
G	220	MET	-	expression tag	UNP O94925
H	201	MET	-	expression tag	UNP O94925
H	202	GLY	-	expression tag	UNP O94925
H	203	SER	-	expression tag	UNP O94925
H	204	SER	-	expression tag	UNP O94925
H	205	HIS	-	expression tag	UNP O94925
H	206	HIS	-	expression tag	UNP O94925
H	207	HIS	-	expression tag	UNP O94925
H	208	HIS	-	expression tag	UNP O94925
H	209	HIS	-	expression tag	UNP O94925
H	210	HIS	-	expression tag	UNP O94925
H	211	SER	-	expression tag	UNP O94925
H	212	SER	-	expression tag	UNP O94925
H	213	GLY	-	expression tag	UNP O94925
H	214	LEU	-	expression tag	UNP O94925
H	215	VAL	-	expression tag	UNP O94925
H	216	PRO	-	expression tag	UNP O94925
H	217	ARG	-	expression tag	UNP O94925
H	218	GLY	-	expression tag	UNP O94925
H	219	SER	-	expression tag	UNP O94925
H	220	MET	-	expression tag	UNP O94925

- Molecule 2 is 2-(pyridin-2-yl)-N-(5-{4-[6-({[3-(trifluoromethoxy)phenyl]acetyl}amino)pyridazin-3-yl]butyl}-1,3,4-thiadiazol-2-yl)acetamide (three-letter code: 63J) (formula: C₂₆H₂₄F₃N₇O₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	D	1	Total	C	F	N	O	S	1	0
			40	26	3	7	3	1		
2	C	1	Total	C	F	N	O	S	1	0
			40	26	3	7	3	1		
2	F	1	Total	C	F	N	O	S	0	0
			40	26	3	7	3	1		
2	F	1	Total	C	F	N	O	S	1	0
			40	26	3	7	3	1		

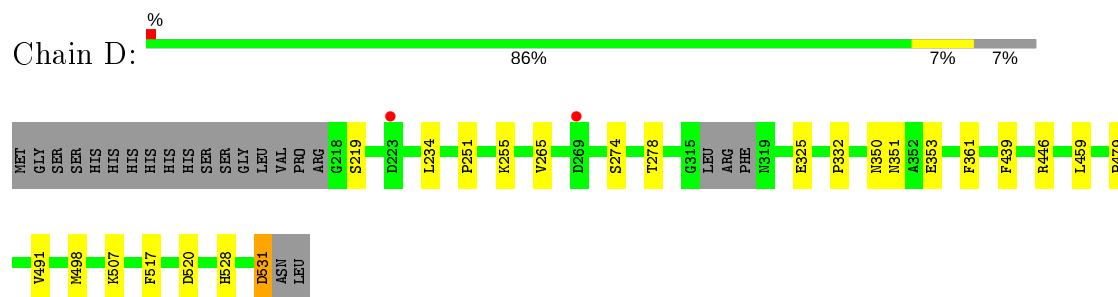
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	160	Total	O	0	0
			160	160		
3	A	153	Total	O	0	0
			153	153		
3	B	156	Total	O	0	0
			156	156		
3	C	148	Total	O	0	0
			148	148		
3	E	151	Total	O	0	0
			151	151		
3	F	163	Total	O	0	0
			163	163		
3	G	151	Total	O	0	0
			151	151		
3	H	155	Total	O	0	0
			155	155		

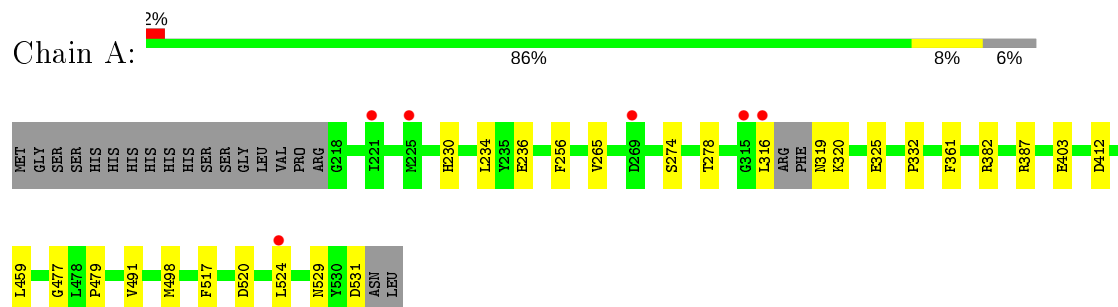
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

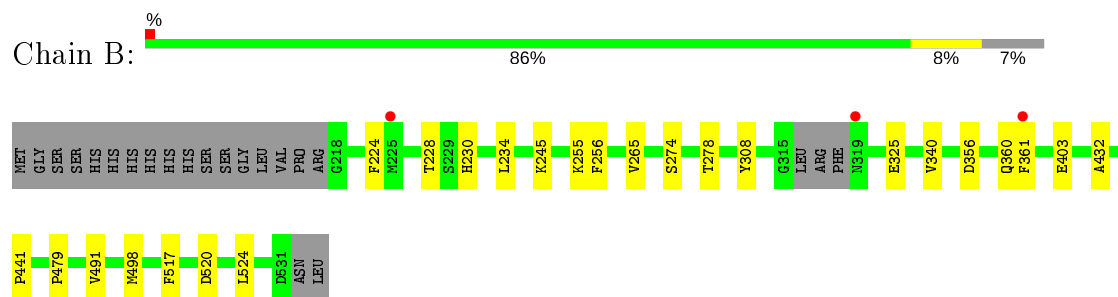
- Molecule 1: Glutaminase kidney isoform, mitochondrial



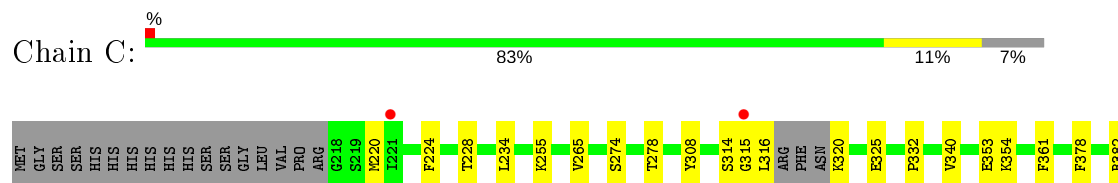
- Molecule 1: Glutaminase kidney isoform, mitochondrial



- Molecule 1: Glutaminase kidney isoform, mitochondrial

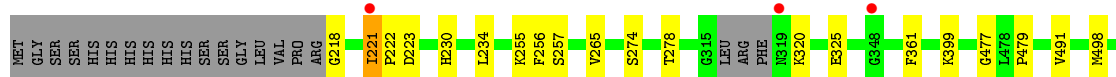
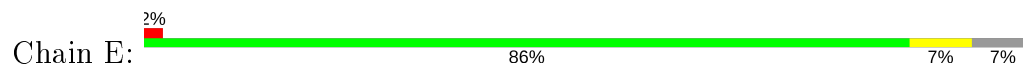


- Molecule 1: Glutaminase kidney isoform, mitochondrial

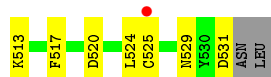
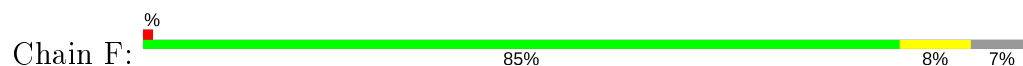




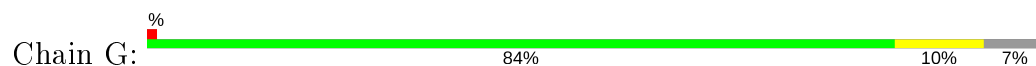
- Molecule 1: Glutaminase kidney isoform, mitochondrial



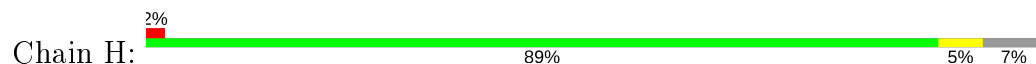
- Molecule 1: Glutaminase kidney isoform, mitochondrial



- Molecule 1: Glutaminase kidney isoform, mitochondrial



- Molecule 1: Glutaminase kidney isoform, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	126.40Å 126.63Å 126.27Å 112.88° 102.81° 112.74°	Depositor
Resolution (Å)	21.36 – 2.10 21.36 – 2.10	Depositor EDS
% Data completeness (in resolution range)	96.5 (21.36-2.10) 91.4 (21.36-2.10)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.09Å)	Xtriage
Refinement program	PHENIX (dev_2420)	Depositor
R, R_{free}	0.190 , 0.213 0.190 , 0.213	Depositor DCC
R_{free} test set	1981 reflections (0.59%)	wwPDB-VP
Wilson B-factor (Å ²)	38.9	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 9.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.419 for h+k+l,-h,-k 0.419 for -k,-l,h+k+l 0.459 for l,-h-k-l,h 0.447 for -l,-k,-h 0.437 for -h-k-l,l,k 0.437 for k,h,-h-k-l 0.447 for -h,h+k+l,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	20485	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 63J

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	0.44	0/2450	0.53	0/3312
1	B	0.44	0/2442	0.53	0/3301
1	C	0.43	0/2442	0.54	0/3301
1	D	0.41	0/2442	0.52	0/3301
1	E	0.45	0/2442	0.53	0/3301
1	F	0.42	0/2442	0.55	0/3301
1	G	0.46	0/2442	0.55	0/3301
1	H	0.42	0/2442	0.53	0/3301
All	All	0.43	0/19544	0.54	0/26419

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2393	0	2324	16	0
1	B	2385	0	2313	18	0
1	C	2385	0	2318	25	0
1	D	2385	0	2313	16	0
1	E	2385	0	2313	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	2385	0	2313	21	0
1	G	2385	0	2313	23	0
1	H	2385	0	2313	11	0
2	C	40	0	0	0	0
2	D	40	0	0	0	0
2	F	80	0	0	0	0
3	A	153	0	0	2	0
3	B	156	0	0	1	0
3	C	148	0	0	3	0
3	D	160	0	0	1	0
3	E	151	0	0	2	0
3	F	163	0	0	5	0
3	G	151	0	0	5	1
3	H	155	0	0	1	1
All	All	20485	0	18520	138	1

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 (138) 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:B:245:LYS:NZ	3:B:601:HOH:O	1.93	0.99
1:D:498:MET:HE1	1:D:517:PHE:HE2	1.45	0.80
1:G:498:MET:HE1	1:G:517:PHE:HE2	1.45	0.79
1:E:498:MET:HE1	1:E:517:PHE:HE2	1.48	0.79
1:H:498:MET:HE1	1:H:517:PHE:HE2	1.46	0.78
1:B:498:MET:HE1	1:B:517:PHE:HE2	1.48	0.77
1:C:498:MET:HE1	1:C:517:PHE:HE2	1.53	0.73
1:D:498:MET:HE1	1:D:517:PHE:CE2	2.24	0.72
1:F:498:MET:HE1	1:F:517:PHE:HE2	1.54	0.72
1:C:325:GLU:H	1:C:325:GLU:CD	1.93	0.71
1:G:498:MET:HE1	1:G:517:PHE:CE2	2.26	0.70
1:B:325:GLU:H	1:B:325:GLU:CD	1.94	0.69
1:A:498:MET:HE1	1:A:517:PHE:HE2	1.57	0.69
1:F:498:MET:HE1	1:F:517:PHE:CE2	2.28	0.68
1:F:325:GLU:CD	1:F:325:GLU:H	1.97	0.68
1:A:325:GLU:CD	1:A:325:GLU:H	1.97	0.68
1:E:498:MET:HE1	1:E:517:PHE:CE2	2.28	0.68
1:H:498:MET:HE1	1:H:517:PHE:CE2	2.27	0.67
1:A:498:MET:HE1	1:A:517:PHE:CE2	2.29	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:478:LEU:HD13	1:F:490:LEU:HD21	1.76	0.67
1:E:221:ILE:O	1:E:223:ASP:N	2.28	0.66
1:B:498:MET:HE1	1:B:517:PHE:CE2	2.29	0.66
1:E:218:GLY:O	3:E:601:HOH:O	2.13	0.66
1:A:403:GLU:CD	1:A:403:GLU:H	1.99	0.65
1:F:513:LYS:NZ	3:F:703:HOH:O	2.29	0.64
1:E:399:LYS:NZ	3:E:602:HOH:O	2.20	0.64
1:C:498:MET:HE1	1:C:517:PHE:CE2	2.33	0.63
1:C:507:LYS:HE2	1:C:507:LYS:H	1.64	0.62
1:C:353:GLU:OE1	3:C:701:HOH:O	2.16	0.62
1:B:403:GLU:H	1:B:403:GLU:CD	2.02	0.62
1:E:325:GLU:CD	1:E:325:GLU:H	2.03	0.62
1:D:234:LEU:HD22	1:D:520:ASP:HB3	1.82	0.61
1:F:234:LEU:HD22	1:F:520:ASP:HB3	1.83	0.60
1:G:425:GLU:OE2	3:G:601:HOH:O	2.15	0.60
1:A:234:LEU:HD22	1:A:520:ASP:HB3	1.82	0.60
1:C:234:LEU:HD22	1:C:520:ASP:HB3	1.85	0.59
1:C:265:VAL:HG22	1:C:498:MET:HG2	1.84	0.58
1:G:274:SER:HB3	1:G:278:THR:HG21	1.86	0.57
1:G:403:GLU:H	1:G:403:GLU:CD	2.09	0.56
1:F:265:VAL:HG22	1:F:498:MET:HG2	1.86	0.56
1:A:274:SER:HB3	1:A:278:THR:HG21	1.88	0.56
1:G:360:GLN:NE2	3:G:604:HOH:O	2.39	0.55
1:B:265:VAL:HG22	1:B:498:MET:HG2	1.88	0.55
1:H:265:VAL:HG22	1:H:498:MET:HG2	1.90	0.54
1:E:256:PHE:CE2	1:G:255:LYS:HD2	2.44	0.53
1:C:316:LEU:N	3:C:703:HOH:O	2.40	0.53
1:H:274:SER:HB3	1:H:278:THR:HG21	1.89	0.53
1:G:507:LYS:HE2	1:G:507:LYS:H	1.73	0.53
1:E:234:LEU:HD22	1:E:520:ASP:HB3	1.91	0.52
1:F:219:SER:HA	3:F:702:HOH:O	2.10	0.51
1:G:251:PRO:O	1:G:255:LYS:HG3	2.10	0.51
1:B:255:LYS:HE2	1:H:256:PHE:CZ	2.46	0.51
1:D:350:ASN:OD1	1:D:353:GLU:HG3	2.11	0.51
1:H:354:LYS:NZ	3:H:604:HOH:O	2.42	0.51
1:H:251:PRO:O	1:H:255:LYS:HG3	2.10	0.51
1:A:316:LEU:C	1:A:319:ASN:HB2	2.31	0.51
1:D:265:VAL:HG22	1:D:498:MET:HG2	1.93	0.50
1:B:230:HIS:HB3	1:B:524:LEU:HD21	1.93	0.50
1:F:387:ARG:NH1	3:F:709:HOH:O	2.45	0.50
1:A:230:HIS:HB3	1:A:524:LEU:HD21	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:234:LEU:HD22	1:G:520:ASP:HB3	1.95	0.49
1:B:356:ASP:O	1:B:360:GLN:HG2	2.12	0.49
1:C:507:LYS:H	1:C:507:LYS:CE	2.25	0.49
1:F:524:LEU:HD23	1:F:525:CYS:SG	2.52	0.49
1:A:477:GLY:O	1:A:529:ASN:HB2	2.12	0.49
1:G:477:GLY:O	1:G:529:ASN:HB2	2.12	0.49
1:D:251:PRO:O	1:D:255:LYS:HG3	2.13	0.49
1:F:319:ASN:ND2	3:F:701:HOH:O	2.22	0.49
1:A:387:ARG:NH1	3:A:603:HOH:O	2.44	0.48
1:F:332:PRO:HD2	1:F:459:LEU:HD13	1.94	0.48
1:C:274:SER:HB3	1:C:278:THR:HG21	1.95	0.48
1:F:274:SER:HB3	1:F:278:THR:HG21	1.95	0.48
1:D:325:GLU:CD	1:D:325:GLU:H	2.17	0.48
1:B:234:LEU:HD22	1:B:520:ASP:HB3	1.95	0.48
1:A:256:PHE:CZ	1:C:255:LYS:HE3	2.49	0.48
1:E:477:GLY:O	1:E:529:ASN:HB2	2.14	0.48
1:H:234:LEU:HD22	1:H:520:ASP:HB3	1.96	0.48
1:F:314:SER:HB3	1:F:315:GLY:HA2	1.96	0.47
1:D:498:MET:HB3	1:D:498:MET:HE3	1.74	0.47
1:G:399:LYS:NZ	3:G:608:HOH:O	2.47	0.47
1:G:230:HIS:HB3	1:G:524:LEU:HD21	1.96	0.47
1:D:531:ASP:OD1	1:C:454:ARG:HD3	2.15	0.47
1:E:265:VAL:HG22	1:E:498:MET:HG2	1.97	0.47
1:E:479:PRO:HD2	1:E:491:VAL:O	2.15	0.47
1:G:351:ASN:OD1	3:G:602:HOH:O	2.20	0.47
1:G:314:SER:OG	1:G:315:GLY:HA2	2.15	0.47
1:A:382:ARG:HD3	1:A:412:ASP:OD1	2.16	0.46
1:G:507:LYS:CE	1:G:507:LYS:H	2.27	0.46
1:B:479:PRO:HD2	1:B:491:VAL:O	2.16	0.46
1:C:435:ALA:HB2	1:C:491:VAL:HG13	1.98	0.45
1:C:477:GLY:O	1:C:529:ASN:HB2	2.15	0.45
1:C:354:LYS:NZ	3:C:707:HOH:O	2.49	0.45
1:G:265:VAL:HG12	1:G:498:MET:HG2	1.98	0.45
1:C:314:SER:OG	1:C:315:GLY:HA2	2.16	0.45
1:F:490:LEU:HD23	1:F:491:VAL:N	2.32	0.45
1:A:265:VAL:HG22	1:A:498:MET:HG2	1.99	0.45
1:E:498:MET:HE3	1:E:498:MET:HB3	1.75	0.45
1:F:507:LYS:HB2	1:F:507:LYS:HE2	1.83	0.44
1:G:332:PRO:HD2	1:G:459:LEU:HD13	2.00	0.44
1:F:507:LYS:CE	1:F:507:LYS:H	2.31	0.44
1:B:224:PHE:O	1:B:228:THR:HG23	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:439:PHE:CE2	1:D:446:ARG:HB2	2.53	0.43
1:D:274:SER:HB3	1:D:278:THR:HG21	1.99	0.43
1:C:382:ARG:NH1	1:C:412:ASP:OD1	2.51	0.43
1:E:230:HIS:HB3	1:E:524:LEU:HD21	1.99	0.43
1:H:332:PRO:HD2	1:H:459:LEU:HD13	2.00	0.43
1:D:332:PRO:HD2	1:D:459:LEU:HD13	2.00	0.43
1:C:332:PRO:HD2	1:C:459:LEU:HD13	2.00	0.43
1:C:308:TYR:HB3	1:C:340:VAL:HG11	2.01	0.42
1:B:274:SER:HB3	1:B:278:THR:HG21	1.99	0.42
1:F:477:GLY:O	1:F:529:ASN:HB2	2.20	0.42
1:A:236:GLU:HG3	3:A:656:HOH:O	2.19	0.42
1:F:507:LYS:HE2	1:F:507:LYS:H	1.84	0.42
1:F:507:LYS:NZ	3:F:716:HOH:O	2.52	0.42
1:F:435:ALA:HB2	1:F:491:VAL:HG13	2.01	0.42
1:A:479:PRO:HD2	1:A:491:VAL:O	2.20	0.42
1:G:479:PRO:HD2	1:G:491:VAL:O	2.20	0.42
1:H:498:MET:HB3	1:H:498:MET:HE3	1.76	0.42
1:E:257:SER:HB2	1:G:257:SER:HA	2.01	0.42
1:A:332:PRO:HD2	1:A:459:LEU:HD13	2.00	0.42
1:C:432:ALA:HB1	1:C:441:PRO:HG2	2.02	0.42
1:B:498:MET:HB3	1:B:498:MET:HE3	1.89	0.41
1:C:498:MET:HE3	1:C:498:MET:HB3	1.87	0.41
1:B:325:GLU:N	1:B:325:GLU:CD	2.70	0.41
1:B:432:ALA:HB1	1:B:441:PRO:HG2	2.02	0.41
1:D:479:PRO:HD2	1:D:491:VAL:O	2.20	0.41
1:D:507:LYS:H	1:D:507:LYS:CE	2.34	0.40
1:G:387:ARG:NH1	3:G:609:HOH:O	2.54	0.40
1:G:378:PHE:CE1	1:G:416:GLN:HG3	2.57	0.40
1:B:256:PHE:CE2	1:H:255:LYS:HD2	2.56	0.40
1:C:378:PHE:CE1	1:C:416:GLN:HG3	2.57	0.40
1:D:528:HIS:CG	1:C:454:ARG:HD2	2.56	0.40
1:B:308:TYR:HB3	1:B:340:VAL:HG11	2.03	0.40
1:D:351:ASN:OD1	3:D:701:HOH:O	2.22	0.40
1:E:256:PHE:CZ	1:G:255:LYS:HD2	2.57	0.40
1:C:224:PHE:O	1:C:228:THR:HG23	2.21	0.40
1:C:479:PRO:HD2	1:C:491:VAL:O	2.21	0.40
1:E:274:SER:HB3	1:E:278:THR:HG21	2.03	0.40

All (1) 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)
3:G:730:HOH:O	3:H:682:HOH:O[1_544]	2.16	0.04

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	308/333 (92%)	303 (98%)	5 (2%)	0	100	100
1	B	307/333 (92%)	302 (98%)	5 (2%)	0	100	100
1	C	307/333 (92%)	297 (97%)	10 (3%)	0	100	100
1	D	307/333 (92%)	304 (99%)	3 (1%)	0	100	100
1	E	307/333 (92%)	301 (98%)	4 (1%)	2 (1%)	22	18
1	F	307/333 (92%)	300 (98%)	7 (2%)	0	100	100
1	G	307/333 (92%)	299 (97%)	8 (3%)	0	100	100
1	H	307/333 (92%)	304 (99%)	3 (1%)	0	100	100
All	All	2457/2664 (92%)	2410 (98%)	45 (2%)	2 (0%)	51	54

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	222	PRO
1	E	221	ILE

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	264/283 (93%)	261 (99%)	3 (1%)	73	79
1	B	263/283 (93%)	262 (100%)	1 (0%)	91	94
1	C	263/283 (93%)	260 (99%)	3 (1%)	73	79
1	D	263/283 (93%)	260 (99%)	3 (1%)	73	79
1	E	263/283 (93%)	260 (99%)	3 (1%)	73	79
1	F	263/283 (93%)	259 (98%)	4 (2%)	65	71
1	G	263/283 (93%)	260 (99%)	3 (1%)	73	79
1	H	263/283 (93%)	260 (99%)	3 (1%)	73	79
All	All	2105/2264 (93%)	2082 (99%)	23 (1%)	73	79

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	219	SER
1	D	361	PHE
1	D	531	ASP
1	A	320	LYS
1	A	361	PHE
1	A	531	ASP
1	B	361	PHE
1	C	220	MET
1	C	320	LYS
1	C	361	PHE
1	E	255	LYS
1	E	320	LYS
1	E	361	PHE
1	F	320	LYS
1	F	361	PHE
1	F	507	LYS
1	F	531	ASP
1	G	220	MET
1	G	361	PHE
1	G	531	ASP
1	H	319	ASN
1	H	361	PHE
1	H	531	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	330	HIS
1	B	319	ASN
1	B	360	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 carbohydrates 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	63J	F	601	-	40,43,43	1.49	4 (10%)	45,58,58	1.56	10 (22%)
2	63J	D	601	-	40,43,43	1.43	5 (12%)	45,58,58	1.36	8 (17%)
2	63J	F	602	-	40,43,43	1.47	4 (10%)	45,58,58	1.44	9 (20%)
2	63J	C	601	-	40,43,43	1.52	5 (12%)	45,58,58	1.42	8 (17%)

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	63J	F	601	-	-	9/25/28/28	0/4/4/4
2	63J	D	601	-	-	2/25/28/28	0/4/4/4
2	63J	F	602	-	-	11/25/28/28	0/4/4/4
2	63J	C	601	-	-	6/25/28/28	0/4/4/4

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	601	63J	C03-C04	5.33	1.52	1.49
2	F	601	63J	C03-C04	5.07	1.52	1.49
2	F	602	63J	C03-C04	5.01	1.52	1.49
2	D	601	63J	C03-C04	4.99	1.52	1.49
2	F	601	63J	C07-N08	4.52	1.44	1.36
2	F	602	63J	C07-N08	4.21	1.44	1.36
2	C	601	63J	C07-N08	4.02	1.43	1.36
2	F	601	63J	C25-N24	3.90	1.44	1.35
2	C	601	63J	C25-N24	3.78	1.44	1.35
2	D	601	63J	C25-N24	3.70	1.43	1.35
2	D	601	63J	C07-N08	3.69	1.43	1.36
2	F	602	63J	C25-N24	3.52	1.43	1.35
2	D	601	63J	C09-N08	3.40	1.43	1.35
2	C	601	63J	C09-N08	3.35	1.43	1.35
2	F	602	63J	C09-N08	3.18	1.42	1.35
2	F	601	63J	C09-N08	3.09	1.42	1.35
2	C	601	63J	C20-N40	2.41	1.35	1.33
2	D	601	63J	C20-N40	2.11	1.34	1.33

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	601	63J	C15-N16-C11	3.93	122.80	117.42
2	F	601	63J	C10-C09-N08	3.84	122.71	114.77
2	D	601	63J	C15-N16-C11	3.61	122.37	117.42
2	F	601	63J	C15-N16-C11	3.58	122.32	117.42
2	F	602	63J	C10-C11-N16	3.21	120.81	116.41
2	F	601	63J	C10-C11-N16	3.12	120.69	116.41
2	F	601	63J	C26-C25-N24	3.09	121.15	114.77
2	F	602	63J	C26-C25-N24	3.07	121.11	114.77
2	F	602	63J	C15-N16-C11	2.97	121.48	117.42
2	C	601	63J	C20-N40-N39	2.87	121.20	119.77
2	C	601	63J	C10-C09-N08	2.80	120.55	114.77
2	D	601	63J	C10-C11-N16	2.76	120.19	116.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	602	63J	C12-C11-N16	-2.75	118.21	122.17
2	F	601	63J	C07-N08-C09	-2.68	122.29	129.54
2	C	601	63J	C10-C11-N16	2.68	120.08	116.41
2	F	601	63J	O17-C09-N08	-2.67	118.75	123.63
2	D	601	63J	C26-C25-N24	2.66	120.26	114.77
2	C	601	63J	C12-C11-N16	-2.61	118.41	122.17
2	F	602	63J	C13-C12-C11	2.55	122.11	118.93
2	F	602	63J	C20-N40-N39	2.51	121.02	119.77
2	F	601	63J	C12-C11-N16	-2.49	118.59	122.17
2	C	601	63J	C26-C25-N24	2.47	119.88	114.77
2	D	601	63J	C20-N40-N39	2.43	120.98	119.77
2	F	602	63J	C23-N39-N40	2.40	121.07	119.37
2	F	601	63J	C20-N40-N39	2.34	120.94	119.77
2	D	601	63J	C10-C09-N08	2.26	119.44	114.77
2	F	602	63J	C10-C09-N08	2.22	119.36	114.77
2	D	601	63J	C12-C11-N16	-2.20	119.00	122.17
2	F	602	63J	C27-C26-C25	-2.17	106.16	112.57
2	F	601	63J	O38-C25-C26	-2.10	117.25	122.03
2	C	601	63J	O17-C09-N08	-2.07	119.86	123.63
2	F	601	63J	C23-N39-N40	2.07	120.83	119.37
2	D	601	63J	C23-N39-N40	2.06	120.83	119.37
2	D	601	63J	C14-C15-N16	-2.06	120.07	123.43
2	C	601	63J	C31-C32-C27	2.05	121.99	119.73

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	601	63J	C01-C02-C03-C04
2	F	601	63J	C02-C03-C04-N05
2	F	602	63J	C02-C03-C04-N05
2	F	602	63J	F37-C34-O33-C31
2	F	602	63J	F35-C34-O33-C31
2	C	601	63J	C25-C26-C27-C28
2	F	601	63J	C02-C01-C19-C20
2	F	602	63J	C02-C01-C19-C20
2	C	601	63J	C25-C26-C27-C32
2	F	602	63J	F36-C34-O33-C31
2	F	602	63J	O17-C09-C10-C11
2	F	602	63J	C19-C01-C02-C03
2	D	601	63J	C19-C01-C02-C03
2	F	601	63J	C19-C01-C02-C03

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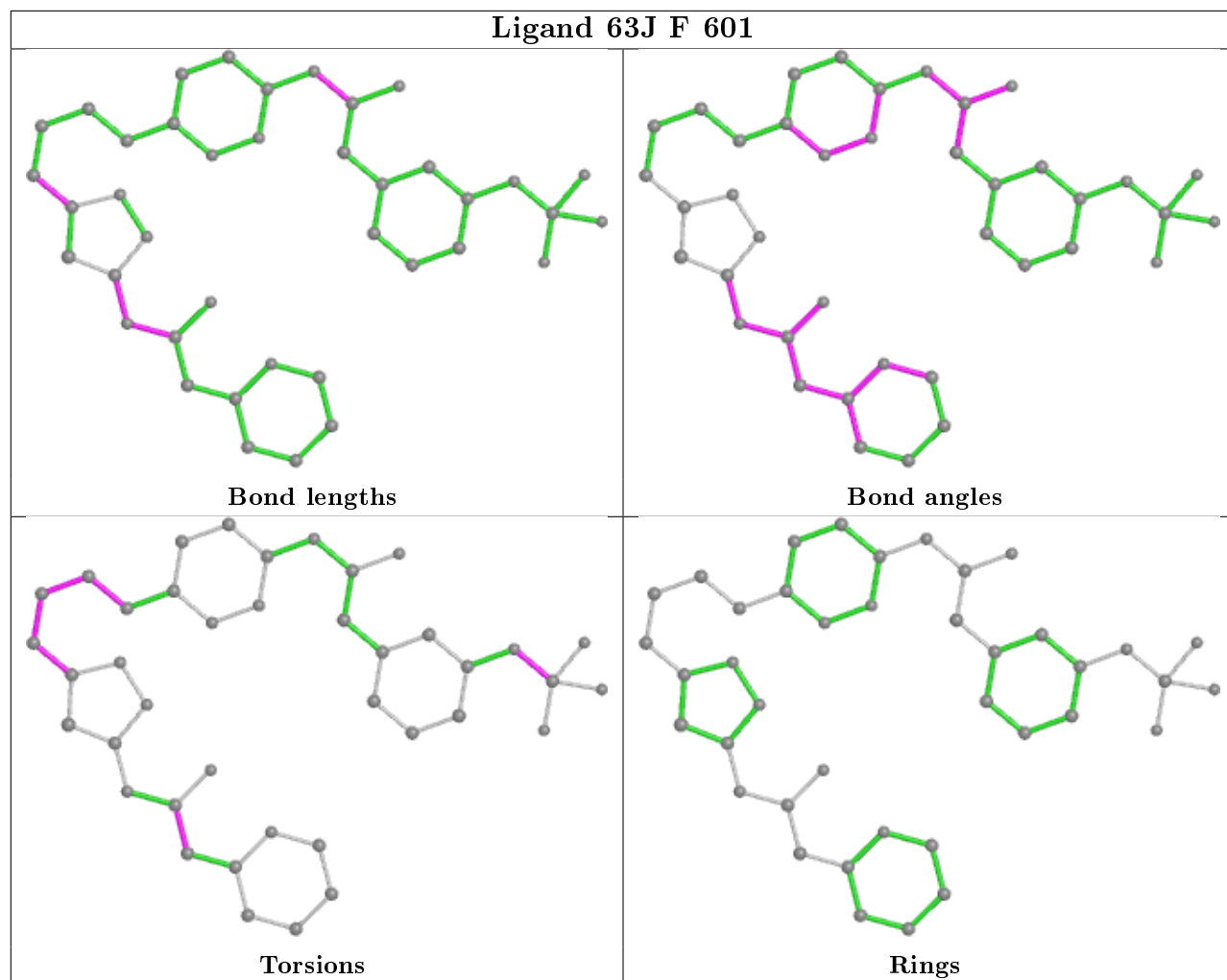
Mol	Chain	Res	Type	Atoms
2	F	601	63J	F35-C34-O33-C31
2	C	601	63J	C19-C01-C02-C03
2	F	602	63J	C01-C02-C03-C04
2	F	602	63J	N08-C09-C10-C11
2	F	601	63J	F37-C34-O33-C31
2	F	601	63J	O17-C09-C10-C11
2	F	601	63J	N08-C09-C10-C11
2	F	601	63J	F36-C34-O33-C31
2	C	601	63J	C32-C31-O33-C34
2	D	601	63J	C01-C02-C03-C04
2	C	601	63J	C01-C02-C03-C04
2	F	602	63J	C32-C31-O33-C34
2	C	601	63J	C30-C31-O33-C34
2	F	602	63J	C30-C31-O33-C34

There are no ring outliers.

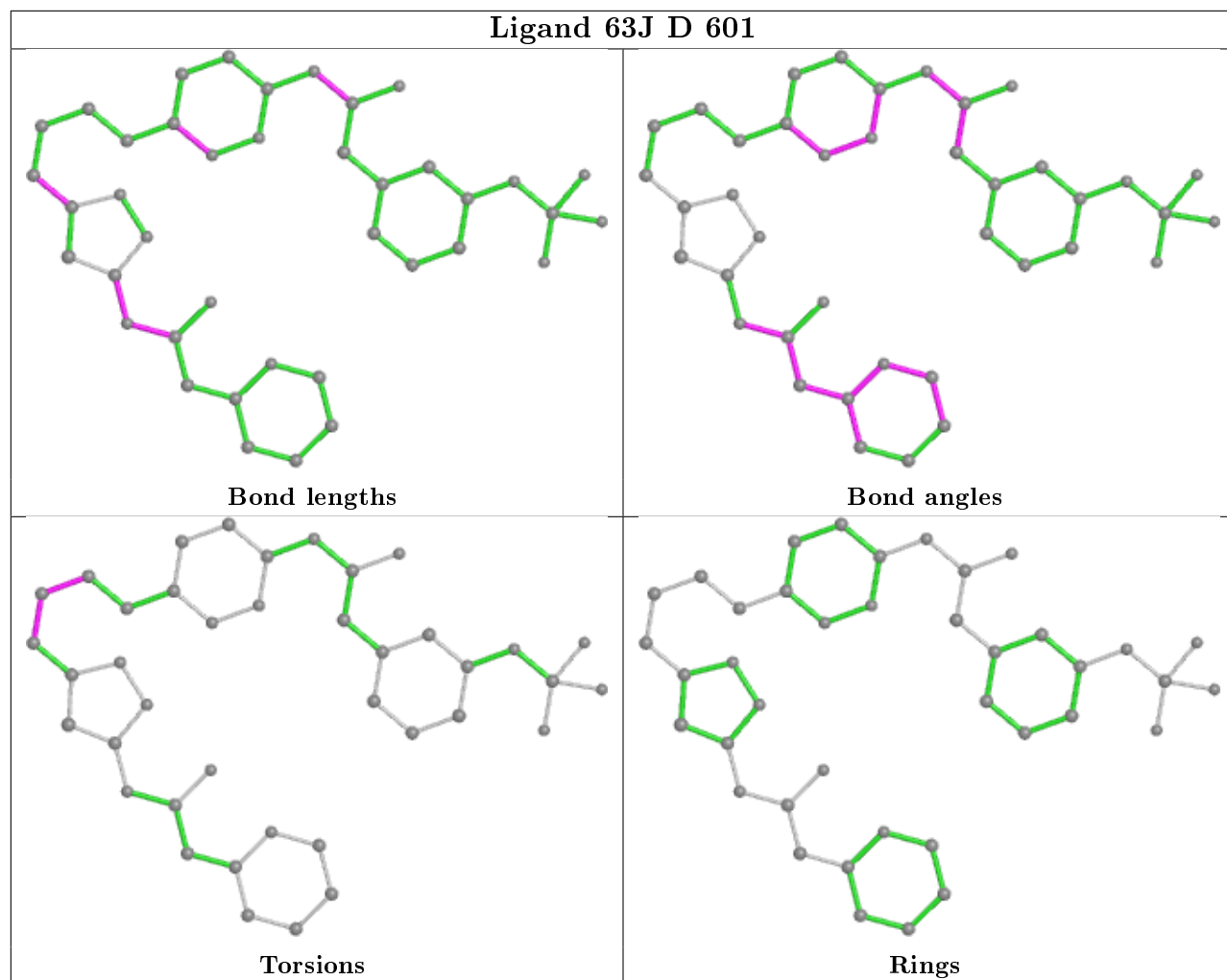
No monomer is involved in short contacts.

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.

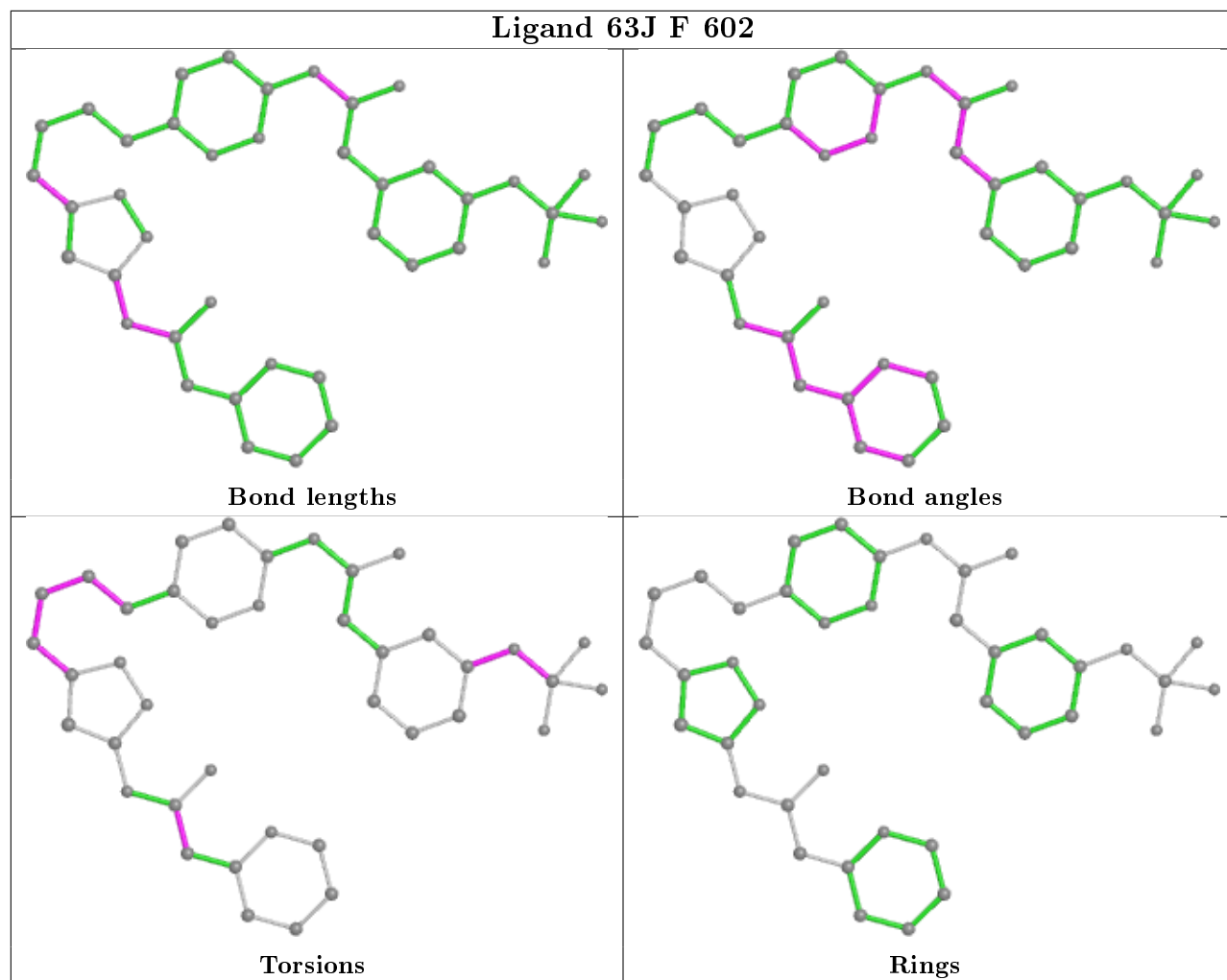
Ligand 63J F 601

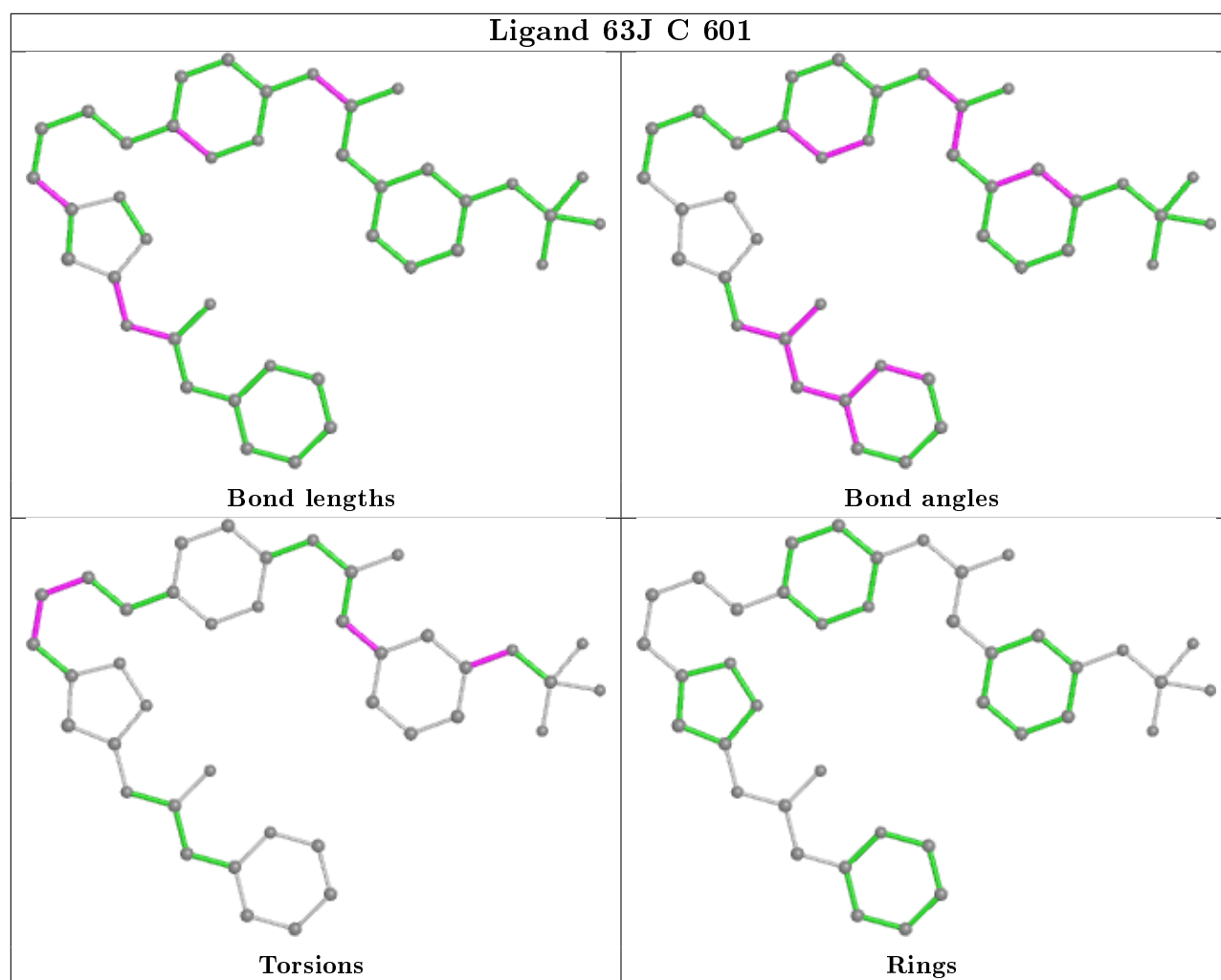


Ligand 63J D 601



Ligand 63J F 602





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	312/333 (93%)	0.25	6 (1%) 66 71	34, 46, 68, 88	0
1	B	311/333 (93%)	0.38	3 (0%) 82 85	34, 45, 66, 87	0
1	C	311/333 (93%)	0.29	2 (0%) 89 91	34, 46, 67, 87	0
1	D	311/333 (93%)	0.26	2 (0%) 89 91	34, 46, 66, 86	0
1	E	311/333 (93%)	0.31	5 (1%) 72 75	33, 45, 66, 85	0
1	F	311/333 (93%)	0.26	3 (0%) 82 85	34, 46, 67, 88	0
1	G	311/333 (93%)	0.38	4 (1%) 77 80	33, 45, 67, 89	0
1	H	311/333 (93%)	0.32	5 (1%) 72 75	33, 45, 67, 86	0
All	All	2489/2664 (93%)	0.31	30 (1%) 79 82	33, 45, 67, 89	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	316	LEU	4.7
1	H	319	ASN	3.3
1	G	524	LEU	3.1
1	A	524	LEU	3.0
1	F	315	GLY	2.8
1	B	361	PHE	2.8
1	F	269	ASP	2.8
1	A	315	GLY	2.7
1	E	319	ASN	2.6
1	G	223	ASP	2.5
1	A	221	ILE	2.5
1	C	315	GLY	2.5
1	E	348	GLY	2.5
1	D	269	ASP	2.4
1	E	524	LEU	2.4
1	A	269	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	225	MET	2.4
1	H	225	MET	2.3
1	H	524	LEU	2.3
1	B	225	MET	2.2
1	E	527	PHE	2.2
1	G	319	ASN	2.2
1	E	221	ILE	2.2
1	G	525	CYS	2.2
1	H	221	ILE	2.2
1	B	319	ASN	2.1
1	F	525	CYS	2.1
1	H	269	ASP	2.1
1	D	223	ASP	2.1
1	C	221	ILE	2.1

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 carbohydrates in this entry.

6.4 Ligands [i](#)

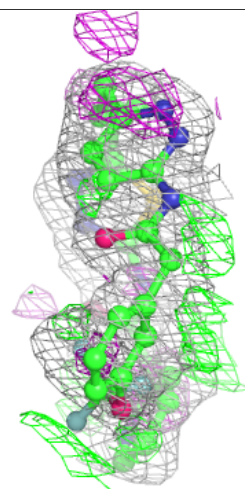
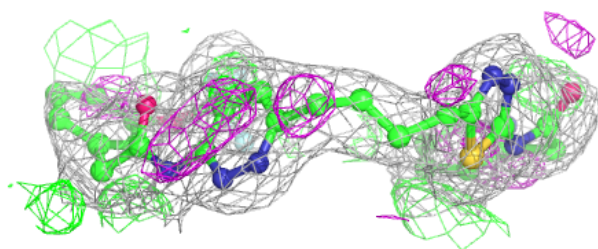
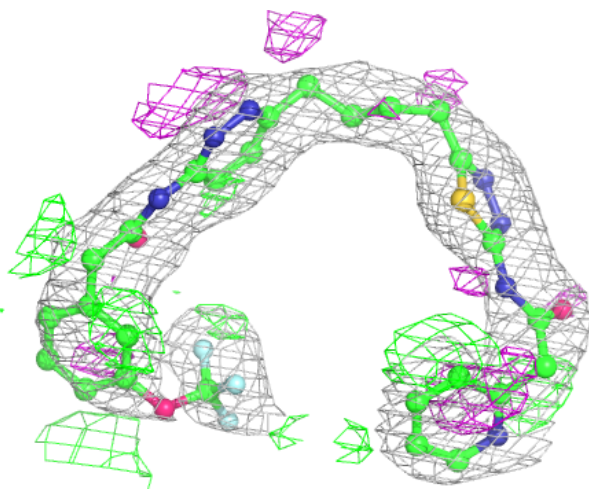
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	63J	F	602	40/40	0.89	0.21	44,64,80,83	5
2	63J	D	601	40/40	0.90	0.20	43,57,73,76	5
2	63J	F	601	40/40	0.90	0.21	43,65,79,80	4
2	63J	C	601	40/40	0.91	0.21	44,58,76,78	5

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.

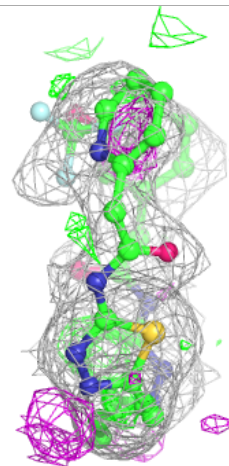
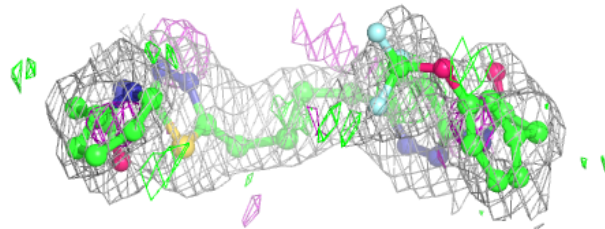
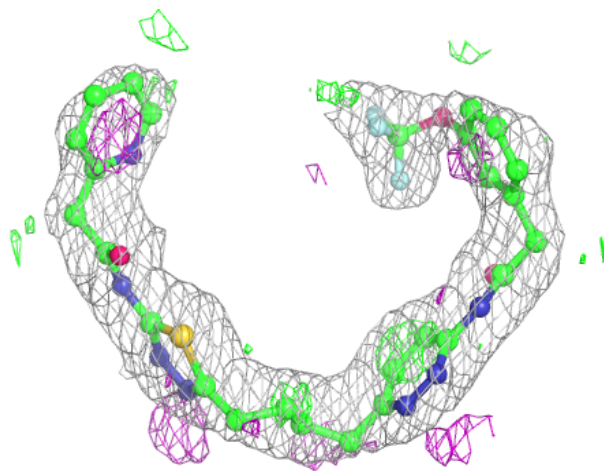
Electron density around 63J F 602:

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



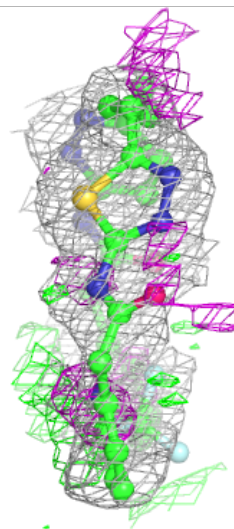
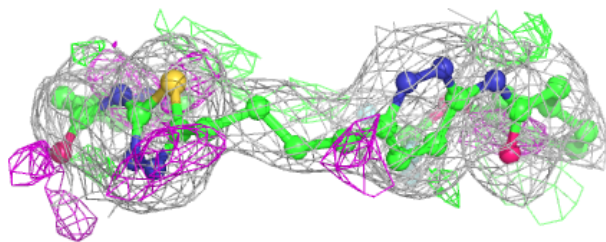
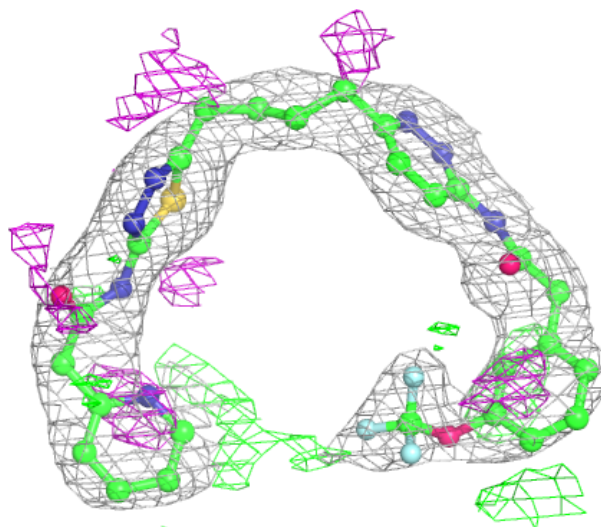
Electron density around 63J D 601:

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



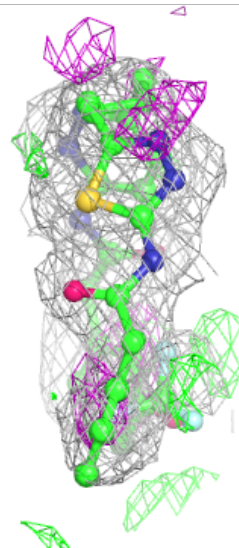
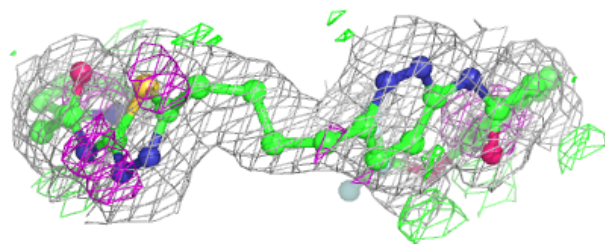
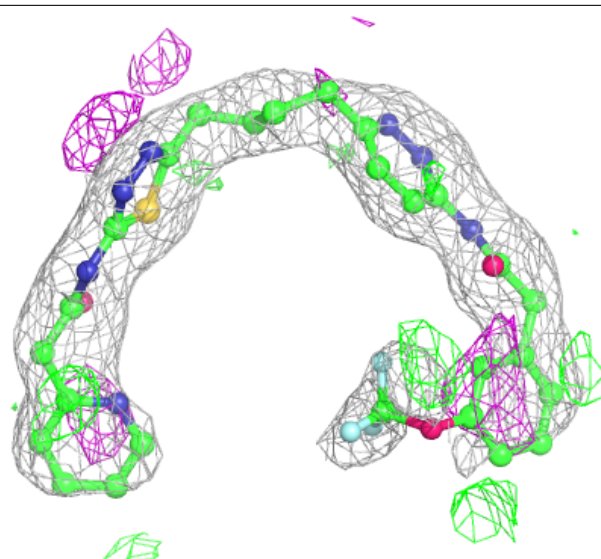
Electron density around 63J F 601:

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



Electron density around 63J 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.