



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 4, 2023 – 02:07 PM EST

PDB ID : 8F65
Title : Crystal structure of VACV D13 in complex with BBL030900
Authors : Subedi, B.P.; Garriga, D.; Coulibaly, F.
Deposited on : 2022-11-16
Resolution : 2.85 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

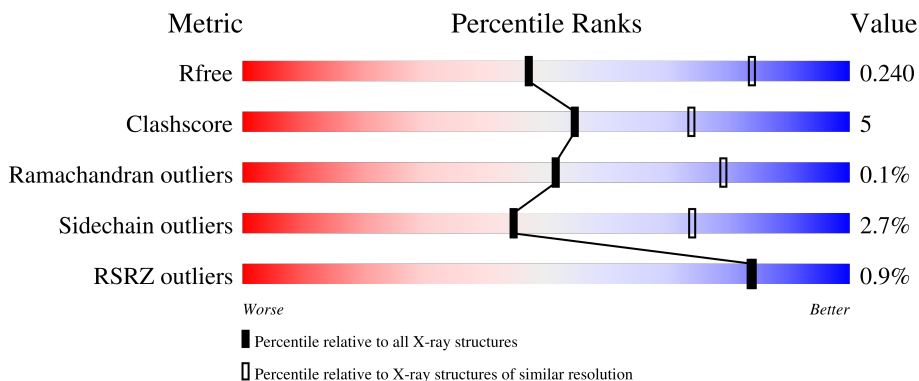
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


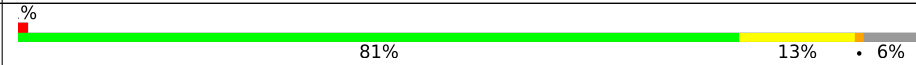
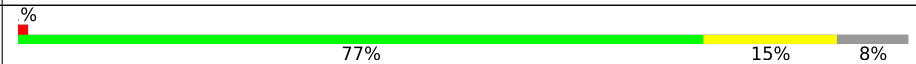
The reported resolution of this entry is 2.85 Å.

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	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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	576	 82% 11% • 6%
1	B	576	 81% 13% • 6%
1	C	576	 77% 15% 8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	FMT	A	601	-	-	X	-
2	FMT	A	607	-	-	X	-
2	FMT	C	609	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 13277 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 Scaffold protein D13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	543	4308	2752	705	836	15	0	5	0
1	B	542	4250	2719	689	829	13	0	0	0
1	C	532	4203	2692	678	820	13	0	3	0

There are 75 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-24	MET	-	initiating methionine	UNP P68440
A	-23	SER	-	expression tag	UNP P68440
A	-22	TYR	-	expression tag	UNP P68440
A	-21	TYR	-	expression tag	UNP P68440
A	-20	HIS	-	expression tag	UNP P68440
A	-19	HIS	-	expression tag	UNP P68440
A	-18	HIS	-	expression tag	UNP P68440
A	-17	HIS	-	expression tag	UNP P68440
A	-16	HIS	-	expression tag	UNP P68440
A	-15	HIS	-	expression tag	UNP P68440
A	-14	ASP	-	expression tag	UNP P68440
A	-13	TYR	-	expression tag	UNP P68440
A	-12	ASP	-	expression tag	UNP P68440
A	-11	ILE	-	expression tag	UNP P68440
A	-10	PRO	-	expression tag	UNP P68440
A	-9	THR	-	expression tag	UNP P68440
A	-8	THR	-	expression tag	UNP P68440
A	-7	GLU	-	expression tag	UNP P68440
A	-6	ASN	-	expression tag	UNP P68440
A	-5	LEU	-	expression tag	UNP P68440
A	-4	TYR	-	expression tag	UNP P68440
A	-3	PHE	-	expression tag	UNP P68440
A	-2	GLN	-	expression tag	UNP P68440

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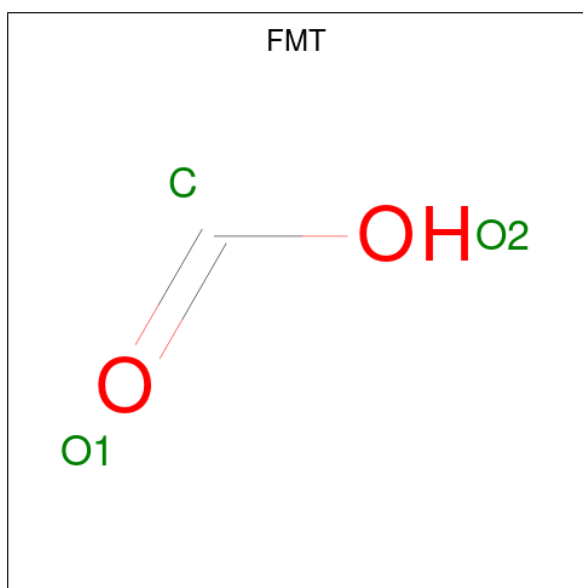
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P68440
A	0	ALA	-	expression tag	UNP P68440
B	-24	MET	-	initiating methionine	UNP P68440
B	-23	SER	-	expression tag	UNP P68440
B	-22	TYR	-	expression tag	UNP P68440
B	-21	TYR	-	expression tag	UNP P68440
B	-20	HIS	-	expression tag	UNP P68440
B	-19	HIS	-	expression tag	UNP P68440
B	-18	HIS	-	expression tag	UNP P68440
B	-17	HIS	-	expression tag	UNP P68440
B	-16	HIS	-	expression tag	UNP P68440
B	-15	HIS	-	expression tag	UNP P68440
B	-14	ASP	-	expression tag	UNP P68440
B	-13	TYR	-	expression tag	UNP P68440
B	-12	ASP	-	expression tag	UNP P68440
B	-11	ILE	-	expression tag	UNP P68440
B	-10	PRO	-	expression tag	UNP P68440
B	-9	THR	-	expression tag	UNP P68440
B	-8	THR	-	expression tag	UNP P68440
B	-7	GLU	-	expression tag	UNP P68440
B	-6	ASN	-	expression tag	UNP P68440
B	-5	LEU	-	expression tag	UNP P68440
B	-4	TYR	-	expression tag	UNP P68440
B	-3	PHE	-	expression tag	UNP P68440
B	-2	GLN	-	expression tag	UNP P68440
B	-1	GLY	-	expression tag	UNP P68440
B	0	ALA	-	expression tag	UNP P68440
C	-24	MET	-	initiating methionine	UNP P68440
C	-23	SER	-	expression tag	UNP P68440
C	-22	TYR	-	expression tag	UNP P68440
C	-21	TYR	-	expression tag	UNP P68440
C	-20	HIS	-	expression tag	UNP P68440
C	-19	HIS	-	expression tag	UNP P68440
C	-18	HIS	-	expression tag	UNP P68440
C	-17	HIS	-	expression tag	UNP P68440
C	-16	HIS	-	expression tag	UNP P68440
C	-15	HIS	-	expression tag	UNP P68440
C	-14	ASP	-	expression tag	UNP P68440
C	-13	TYR	-	expression tag	UNP P68440
C	-12	ASP	-	expression tag	UNP P68440
C	-11	ILE	-	expression tag	UNP P68440
C	-10	PRO	-	expression tag	UNP P68440

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-9	THR	-	expression tag	UNP P68440
C	-8	THR	-	expression tag	UNP P68440
C	-7	GLU	-	expression tag	UNP P68440
C	-6	ASN	-	expression tag	UNP P68440
C	-5	LEU	-	expression tag	UNP P68440
C	-4	TYR	-	expression tag	UNP P68440
C	-3	PHE	-	expression tag	UNP P68440
C	-2	GLN	-	expression tag	UNP P68440
C	-1	GLY	-	expression tag	UNP P68440
C	0	ALA	-	expression tag	UNP P68440

- Molecule 2 is FORMIC ACID (three-letter code: FMT) (formula: CH₂O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 3 1 2	0	0
2	A	1	Total C O 3 1 2	0	0
2	A	1	Total C O 3 1 2	0	0
2	A	1	Total C O 3 1 2	0	0
2	A	1	Total C O 3 1 2	0	0
2	A	1	Total C O 3 1 2	0	0

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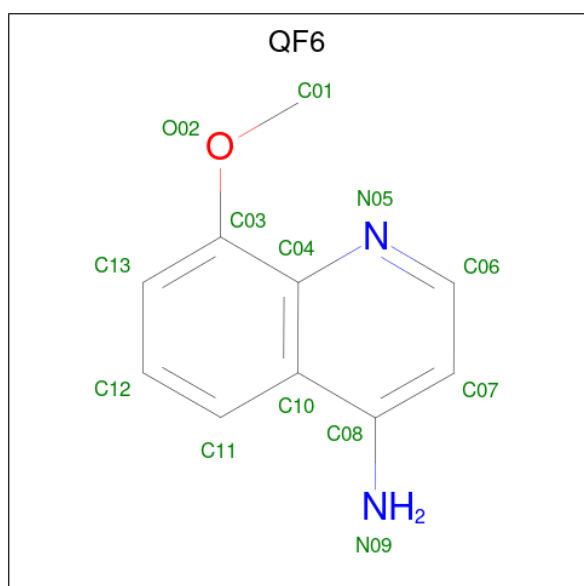
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total 3	C 1	O 2	0	0
2	A	1	Total 3	C 1	O 2	0	0
2	A	1	Total 3	C 1	O 2	0	0
2	A	1	Total 3	C 1	O 2	0	0
2	A	1	Total 3	C 1	O 2	0	0
2	A	1	Total 3	C 1	O 2	0	0
2	A	1	Total 3	C 1	O 2	0	0
2	B	1	Total 3	C 1	O 2	0	0
2	B	1	Total 3	C 1	O 2	0	0
2	B	1	Total 3	C 1	O 2	0	0
2	B	1	Total 3	C 1	O 2	0	0
2	B	1	Total 3	C 1	O 2	0	0
2	B	1	Total 3	C 1	O 2	0	0
2	B	1	Total 3	C 1	O 2	0	0
2	B	1	Total 3	C 1	O 2	0	0
2	B	1	Total 3	C 1	O 2	0	0
2	C	1	Total 3	C 1	O 2	0	0
2	C	1	Total 3	C 1	O 2	0	0
2	C	1	Total 3	C 1	O 2	0	0
2	C	1	Total 3	C 1	O 2	0	0
2	C	1	Total 3	C 1	O 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			3	1	2		
2	C	1	Total	C	O	0	0
			3	1	2		
2	C	1	Total	C	O	0	0
			3	1	2		
2	C	1	Total	C	O	0	0
			3	1	2		

- Molecule 3 is 8-methoxyquinolin-4-amine (three-letter code: QF6) (formula: C₁₀H₁₀N₂O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			13	10	2	1		

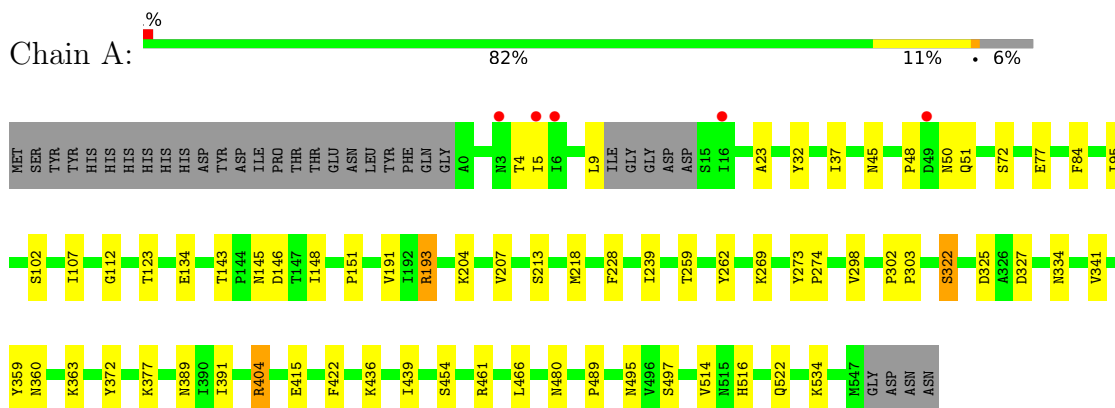
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	150	Total	O	0	0
			150	150		
4	B	122	Total	O	0	0
			122	122		
4	C	138	Total	O	0	0
			138	138		

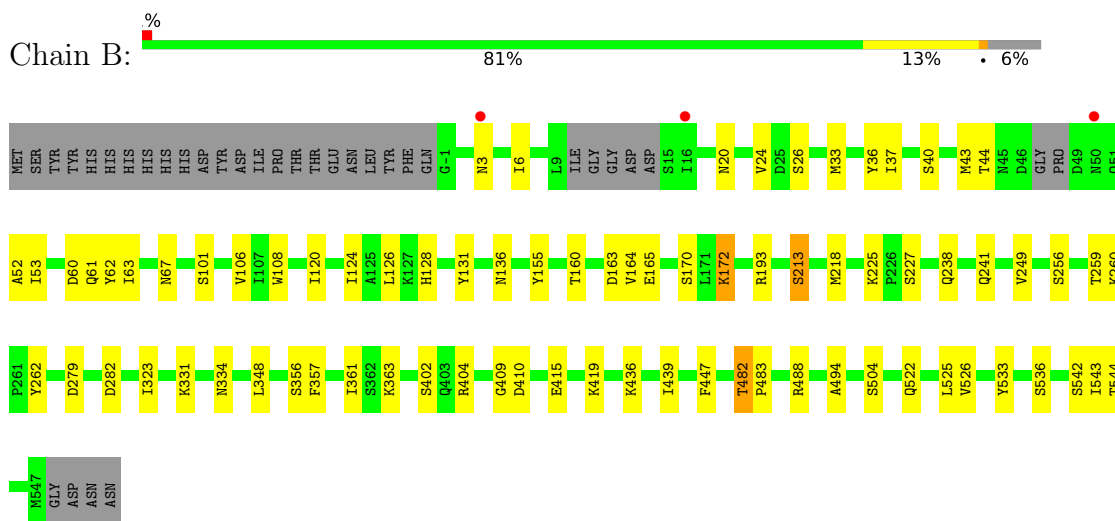
3 Residue-property plots [i](#)

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.

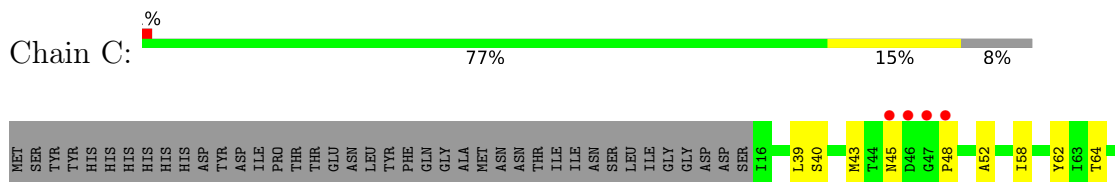
• Molecule 1: Scaffold protein D13

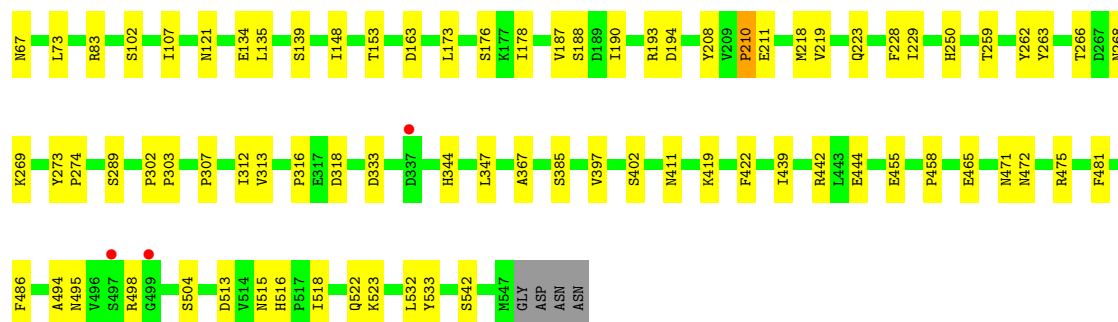


• Molecule 1: Scaffold protein D13



• Molecule 1: Scaffold protein D13





4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	191.08Å 191.08Å 254.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.26 – 2.85 47.77 – 2.85	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.26-2.85) 94.1 (47.77-2.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.56 (at 2.86Å)	Xtrriage
Refinement program	PHENIX 1.20rc4_4425, BUSTER	Depositor
R, R_{free}	0.173 , 0.239 0.174 , 0.240	Depositor DCC
R_{free} test set	3255 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	50.2	Xtrriage
Anisotropy	0.058	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13277	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.60% 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: FMT, QF6

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.46	0/4407	0.62	0/6000
1	B	0.43	0/4341	0.59	0/5911
1	C	0.43	0/4300	0.61	1/5857 (0.0%)
All	All	0.44	0/13048	0.61	1/17768 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	481	PHE	C-N-CA	5.68	135.91	121.70

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	4308	0	4243	43	0
1	B	4250	0	4163	49	0
1	C	4203	0	4123	53	0
2	A	39	0	13	5	0
2	B	24	0	8	1	0
2	C	30	0	10	2	0
3	B	13	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	150	0	0	3	0
4	B	122	0	0	1	0
4	C	138	0	0	0	0
All	All	13277	0	12560	132	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 5.

All (132) 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:259:THR:HG21	1:A:439:ILE:HD11	1.62	0.81
1:B:415:GLU:N	1:B:415:GLU:OE2	2.15	0.79
1:A:143:THR:HG22	1:A:145:ASN:H	1.54	0.71
1:A:134:GLU:OE2	1:B:436:LYS:NZ	2.25	0.69
1:C:83[A]:ARG:NH1	1:C:194:ASP:OD2	2.26	0.69
1:B:43:MET:HE1	1:B:52:ALA:HB1	1.76	0.68
1:C:266:THR:HG22	1:C:268:ASN:H	1.58	0.68
1:A:461:ARG:NH1	4:A:702:HOH:O	2.29	0.66
1:B:60:ASP:HB3	1:B:63:ILE:HD12	1.77	0.66
1:C:266:THR:CG2	1:C:268:ASN:H	2.09	0.64
1:A:360:ASN:HB3	1:A:363:LYS:HG3	1.79	0.63
1:A:84:PHE:HB3	1:A:148:ILE:HG21	1.81	0.63
1:B:356:SER:OG	1:B:357:PHE:N	2.28	0.62
1:A:495:ASN:OD1	1:A:497:SER:HB3	2.01	0.61
1:A:5:ILE:HG23	1:B:33:MET:HG2	1.82	0.61
1:A:415:GLU:OE1	1:A:415:GLU:N	2.33	0.60
1:A:514:VAL:HG13	2:A:607:FMT:H	1.85	0.59
1:A:262:TYR:HB3	1:A:522:GLN:HB2	1.83	0.58
1:A:228:PHE:CE1	1:A:534:LYS:HG3	2.39	0.57
1:B:24:VAL:HG21	1:C:486:PHE:HD1	1.70	0.56
1:B:533:TYR:CD2	1:B:543:ILE:HD12	2.40	0.56
1:C:259:THR:HG21	1:C:439:ILE:HD11	1.87	0.56
1:B:419:LYS:NZ	4:B:703:HOH:O	2.39	0.56
1:B:24:VAL:HG21	1:C:486:PHE:CD1	2.41	0.55
1:B:67:ASN:HB2	1:B:163:ASP:OD1	2.06	0.55
1:B:238:GLN:HB2	1:B:241:GLN:HG2	1.89	0.55
1:C:367:ALA:HB3	1:C:385:SER:HB2	1.88	0.55
1:C:442:ARG:NH2	1:C:444:GLU:OE2	2.37	0.55
1:A:239:ILE:HD11	1:A:522:GLN:HB3	1.90	0.54
1:C:273:TYR:CD1	1:C:274:PRO:HD2	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:356:SER:HG	1:B:357:PHE:H	1.54	0.54
1:C:43:MET:HE1	1:C:52:ALA:HB1	1.90	0.54
1:C:307:PRO:HG2	1:C:411:ASN:HA	1.89	0.53
1:C:64:THR:HG21	1:C:219:VAL:HB	1.91	0.53
1:C:83[A]:ARG:NH2	1:C:419:LYS:HG2	2.23	0.53
1:C:250:HIS:HA	1:C:494:ALA:HB2	1.92	0.52
1:B:6:ILE:HA	1:C:218:MET:SD	2.50	0.52
1:B:259:THR:HG21	1:B:439:ILE:HD11	1.93	0.51
1:C:269:LYS:HG3	1:C:422:PHE:CE1	2.46	0.51
1:C:102:SER:HB3	1:C:107:ILE:HD11	1.91	0.51
1:B:170:SER:O	1:B:172:LYS:NZ	2.43	0.51
1:A:143:THR:HB	1:A:146:ASP:HB2	1.91	0.51
1:C:173:LEU:O	1:C:176:SER:OG	2.23	0.51
1:B:249:VAL:O	1:B:494:ALA:HB2	2.11	0.50
1:B:128:HIS:NE2	1:C:465:GLU:OE1	2.31	0.50
1:C:471:ASN:OD1	1:C:471:ASN:N	2.45	0.50
1:C:495:ASN:HD22	1:C:498:ARG:CB	2.25	0.50
1:A:102:SER:HB3	1:A:107:ILE:HD11	1.93	0.50
1:B:3:ASN:HA	1:B:6:ILE:HD11	1.93	0.49
1:A:514:VAL:H	2:A:607:FMT:H	1.76	0.49
1:C:269:LYS:HG3	1:C:422:PHE:CD1	2.48	0.49
1:B:131:TYR:HA	1:B:136:ASN:OD1	2.13	0.49
1:B:20:ASN:ND2	1:C:223:GLN:HB3	2.28	0.48
1:A:298:VAL:HG22	1:A:341:VAL:HG22	1.96	0.48
1:A:77:GLU:HB2	1:A:151:PRO:HG3	1.94	0.48
1:A:404[A]:ARG:HE	1:A:404[A]:ARG:HB3	1.55	0.47
1:A:48:PRO:O	1:A:51:GLN:HB2	2.14	0.47
1:A:9:LEU:HD23	1:B:218:MET:HG2	1.96	0.47
1:B:101:SER:HB3	1:B:106:VAL:HA	1.97	0.47
1:B:260:LYS:HG3	1:B:526:VAL:HG21	1.97	0.47
1:A:23:ALA:HB1	1:B:225:LYS:HB2	1.97	0.46
1:B:6:ILE:HG22	1:C:218:MET:SD	2.55	0.46
1:A:436:LYS:NZ	1:C:134:GLU:OE2	2.42	0.46
1:C:208:TYR:O	1:C:210:PRO:HD3	2.15	0.46
1:A:363:LYS:NZ	4:A:710:H0H:O	2.48	0.46
1:C:263:TYR:O	1:C:266:THR:HB	2.16	0.46
1:C:455:GLU:OE2	1:C:475:ARG:HD3	2.15	0.46
1:A:191:VAL:HG12	1:A:193[A]:ARG:HD2	1.98	0.46
1:C:516:HIS:HB3	2:C:609:FMT:C	2.46	0.46
1:C:316:PRO:HB2	1:C:318[B]:ASP:OD1	2.15	0.45
1:C:518:ILE:HG22	2:C:609:FMT:H	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:58:ILE:HD12	1:C:178:ILE:HG22	1.98	0.45
1:A:325:ASP:HB2	4:A:710:HOH:O	2.17	0.45
1:A:273:TYR:CD1	1:A:274:PRO:HD2	2.52	0.44
1:A:404[B]:ARG:H	1:A:404[B]:ARG:HG3	1.42	0.44
1:C:187:VAL:HG23	1:C:190:ILE:HD11	2.00	0.44
1:C:302:PRO:HA	1:C:303:PRO:HD3	1.87	0.44
1:A:516:HIS:HB3	2:A:601:FMT:C	2.47	0.44
1:C:139:SER:HB2	1:C:148:ILE:HG22	2.00	0.44
1:A:302:PRO:HA	1:A:303:PRO:HD3	1.89	0.44
1:C:229:ILE:HG23	1:C:533:TYR:HB2	1.99	0.44
1:C:62:TYR:O	1:C:218:MET:HG2	2.17	0.44
1:B:172:LYS:HB2	1:B:172:LYS:HE2	1.76	0.43
1:A:480:ASN:ND2	1:A:489:PRO:HB3	2.33	0.43
1:B:120:ILE:HD12	1:B:126:LEU:HD11	2.00	0.43
1:C:67:ASN:HB2	1:C:163:ASP:OD1	2.19	0.43
1:B:43:MET:HE2	1:B:43:MET:HB3	1.91	0.43
1:C:266:THR:HG22	1:C:268:ASN:N	2.29	0.43
1:A:377:LYS:HD3	1:A:377:LYS:HA	1.79	0.43
1:C:303:PRO:HG2	1:C:312:ILE:HG21	1.99	0.43
1:A:359:TYR:CE1	2:A:601:FMT:H	2.54	0.43
1:B:348:LEU:HA	1:B:409:GLY:O	2.19	0.43
1:A:37:ILE:O	1:A:213:SER:HA	2.19	0.42
1:C:83[A]:ARG:HA	1:C:139:SER:O	2.19	0.42
1:A:334:ASN:O	1:A:372:TYR:HB3	2.18	0.42
1:B:262:TYR:HD2	1:B:522:GLN:NE2	2.17	0.42
1:B:37:ILE:O	1:B:213:SER:HA	2.20	0.42
1:B:279:ASP:HB2	1:B:282:ASP:OD2	2.20	0.42
1:B:348:LEU:HD23	1:B:410:ASP:HA	2.02	0.42
1:C:135:LEU:HD23	1:C:135:LEU:HA	1.89	0.42
1:A:269:LYS:HG3	1:A:422:PHE:CE1	2.55	0.42
1:B:44:THR:HG23	1:B:53:ILE:HB	2.02	0.41
1:A:466:LEU:HD12	1:A:466:LEU:HA	1.88	0.41
1:B:323:ILE:O	1:B:363:LYS:HE3	2.21	0.41
1:B:402:SER:OG	1:B:404:ARG:NH1	2.52	0.41
1:B:482:THR:HG22	1:B:483:PRO:HD2	2.02	0.41
1:C:513:ASP:OD1	1:C:515:ASN:HB2	2.20	0.41
1:B:334:ASN:O	1:B:334:ASN:ND2	2.54	0.41
1:C:228:PHE:HB3	1:C:532:LEU:HD11	2.02	0.41
1:C:439:ILE:HD12	1:C:523:LYS:HD2	2.02	0.41
1:A:32:TYR:HA	1:A:218:MET:O	2.20	0.41
1:A:322:SER:OG	1:A:327:ASP:OD1	2.31	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:ILE:H	1:B:124:ILE:HG12	1.65	0.41
1:C:73:LEU:O	1:C:153:THR:HA	2.21	0.41
1:C:523:LYS:HB2	1:C:523:LYS:HE2	1.80	0.41
1:B:108:TRP:CZ3	1:B:160:THR:HG22	2.55	0.41
1:B:260:LYS:HE3	1:B:260:LYS:HB3	1.77	0.41
1:B:361:ILE:HD12	1:B:361:ILE:HA	1.91	0.41
1:A:95:ILE:O	1:A:112:GLY:HA3	2.20	0.41
1:B:36:TYR:CE2	2:B:602:FMT:H	2.56	0.41
1:C:269:LYS:O	1:C:397:VAL:HG23	2.21	0.41
1:B:155:TYR:OH	1:C:458:PRO:HD2	2.21	0.41
1:B:525:LEU:HD12	1:B:525:LEU:HA	1.85	0.41
1:C:39:LEU:O	1:C:211:GLU:HG3	2.20	0.41
1:C:262:TYR:HD2	1:C:522:GLN:NE2	2.19	0.40
1:A:9:LEU:CD2	1:B:218:MET:HG2	2.51	0.40
1:B:62:TYR:CD1	1:B:218:MET:HE1	2.56	0.40
1:A:359:TYR:CZ	2:A:601:FMT:H	2.56	0.40
1:B:164:VAL:HG12	1:B:165:GLU:HG2	2.04	0.40
1:A:389:ASN:OD1	1:A:391:ILE:HB	2.22	0.40
1:C:344:HIS:CE1	1:C:347:LEU:HA	2.56	0.40
1:C:513:ASP:HB3	1:C:516:HIS:CD2	2.57	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	544/576 (94%)	516 (95%)	28 (5%)	0	100	100
1	B	536/576 (93%)	507 (95%)	29 (5%)	0	100	100
1	C	533/576 (92%)	501 (94%)	30 (6%)	2 (0%)	34	62
All	All	1613/1728 (93%)	1524 (94%)	87 (5%)	2 (0%)	51	79

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	48	PRO
1	C	210	PRO

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	493/528 (93%)	480 (97%)	13 (3%)	46	75
1	B	483/528 (92%)	467 (97%)	16 (3%)	38	68
1	C	480/528 (91%)	468 (98%)	12 (2%)	47	76
All	All	1456/1584 (92%)	1415 (97%)	41 (3%)	44	73

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	THR
1	A	45	ASN
1	A	50	ASN
1	A	72	SER
1	A	123	THR
1	A	193[A]	ARG
1	A	193[B]	ARG
1	A	204	LYS
1	A	207	VAL
1	A	322	SER
1	A	404[A]	ARG
1	A	404[B]	ARG
1	A	454	SER
1	B	26	SER
1	B	40	SER
1	B	61	GLN
1	B	172	LYS
1	B	193	ARG
1	B	213	SER

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Mol	Chain	Res	Type
1	B	227	SER
1	B	256	SER
1	B	331	LYS
1	B	447	PHE
1	B	482	THR
1	B	488	ARG
1	B	504	SER
1	B	536	SER
1	B	542	SER
1	B	544	THR
1	C	40	SER
1	C	45	ASN
1	C	121	ASN
1	C	188	SER
1	C	193	ARG
1	C	289	SER
1	C	313	VAL
1	C	333	ASP
1	C	402	SER
1	C	472	ASN
1	C	504	SER
1	C	542	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	ASN
1	B	334	ASN
1	B	522	GLN
1	C	104	ASN
1	C	145	ASN
1	C	238	GLN
1	C	334	ASN
1	C	495	ASN
1	C	522	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

32 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	FMT	C	608	-	2,2,2	0.59	0	1,1,1	0.06	0
2	FMT	A	608	-	2,2,2	0.70	0	1,1,1	0.27	0
2	FMT	C	603	-	2,2,2	0.75	0	1,1,1	0.37	0
2	FMT	B	608	-	2,2,2	0.63	0	1,1,1	0.13	0
2	FMT	A	606	-	2,2,2	0.63	0	1,1,1	0.08	0
2	FMT	A	605	-	2,2,2	0.70	0	1,1,1	0.31	0
2	FMT	B	601	-	2,2,2	0.75	0	1,1,1	0.38	0
3	QF6	B	609	-	14,14,14	0.21	0	19,19,19	0.28	0
2	FMT	A	604	-	2,2,2	0.66	0	1,1,1	0.25	0
2	FMT	B	603	-	2,2,2	0.64	0	1,1,1	0.27	0
2	FMT	A	610	-	2,2,2	0.74	0	1,1,1	0.40	0
2	FMT	A	602	-	2,2,2	0.79	0	1,1,1	0.47	0
2	FMT	C	606	-	2,2,2	0.61	0	1,1,1	0.06	0
2	FMT	A	609	-	2,2,2	0.64	0	1,1,1	0.15	0
2	FMT	C	601	-	2,2,2	0.58	0	1,1,1	0.16	0
2	FMT	C	605	-	2,2,2	0.70	0	1,1,1	0.40	0
2	FMT	C	602	-	2,2,2	0.64	0	1,1,1	0.16	0
2	FMT	C	607	-	2,2,2	0.68	0	1,1,1	0.28	0
2	FMT	C	604	-	2,2,2	0.70	0	1,1,1	0.44	0
2	FMT	B	605	-	2,2,2	0.63	0	1,1,1	0.29	0
2	FMT	C	610	-	2,2,2	0.61	0	1,1,1	0.09	0
2	FMT	A	611	-	2,2,2	0.63	0	1,1,1	0.06	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FMT	A	601	-	2,2,2	0.64	0	1,1,1	0.29	0
2	FMT	A	612	-	2,2,2	0.68	0	1,1,1	0.22	0
2	FMT	A	613	-	2,2,2	0.67	0	1,1,1	0.23	0
2	FMT	A	607	-	2,2,2	0.57	0	1,1,1	0.06	0
2	FMT	B	604	-	2,2,2	0.65	0	1,1,1	0.14	0
2	FMT	B	606	-	2,2,2	0.64	0	1,1,1	0.25	0
2	FMT	A	603	-	2,2,2	0.70	0	1,1,1	0.51	0
2	FMT	C	609	-	2,2,2	0.62	0	1,1,1	0.04	0
2	FMT	B	607	-	2,2,2	0.65	0	1,1,1	0.06	0
2	FMT	B	602	-	2,2,2	0.67	0	1,1,1	0.19	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	QF6	B	609	-	-	0/2/2/2	0/2/2/2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

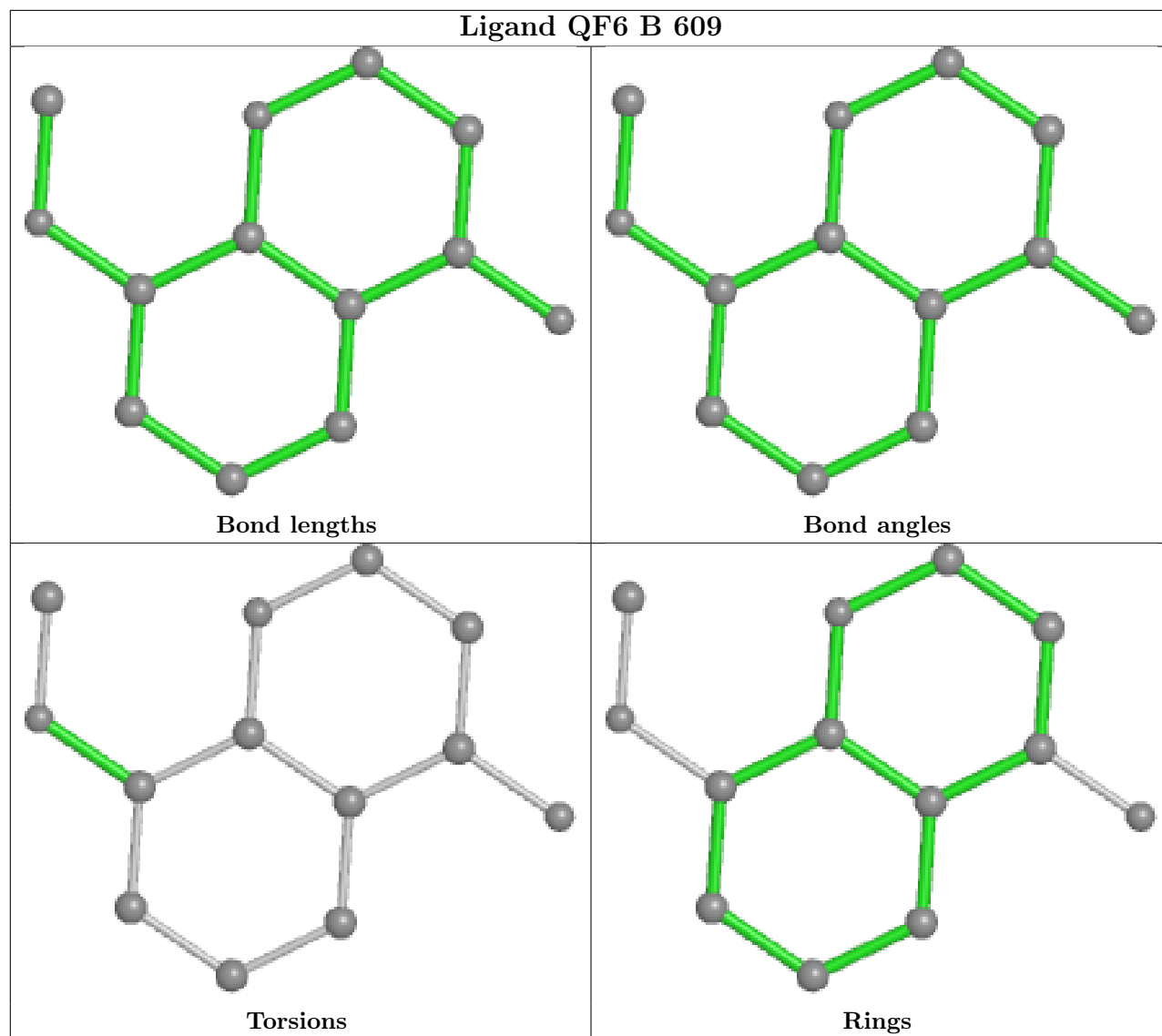
There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	FMT	3	0
2	A	607	FMT	2	0
2	C	609	FMT	2	0
2	B	602	FMT	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.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	543/576 (94%)	-0.33	5 (0%) 84 84	28, 43, 76, 110	0
1	B	542/576 (94%)	-0.27	3 (0%) 89 89	33, 48, 81, 109	0
1	C	532/576 (92%)	-0.33	7 (1%) 77 76	31, 45, 74, 117	0
All	All	1617/1728 (93%)	-0.31	15 (0%) 84 84	28, 46, 79, 117	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	337	ASP	3.7
1	C	47	GLY	3.5
1	A	3	ASN	3.3
1	C	48	PRO	2.8
1	C	497	SER	2.6
1	A	49	ASP	2.5
1	C	499	GLY	2.4
1	B	16	ILE	2.4
1	A	5	ILE	2.2
1	C	46	ASP	2.1
1	A	16	ILE	2.1
1	B	50	ASN	2.1
1	A	6	ILE	2.0
1	B	3	ASN	2.0
1	C	45	ASN	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](#)

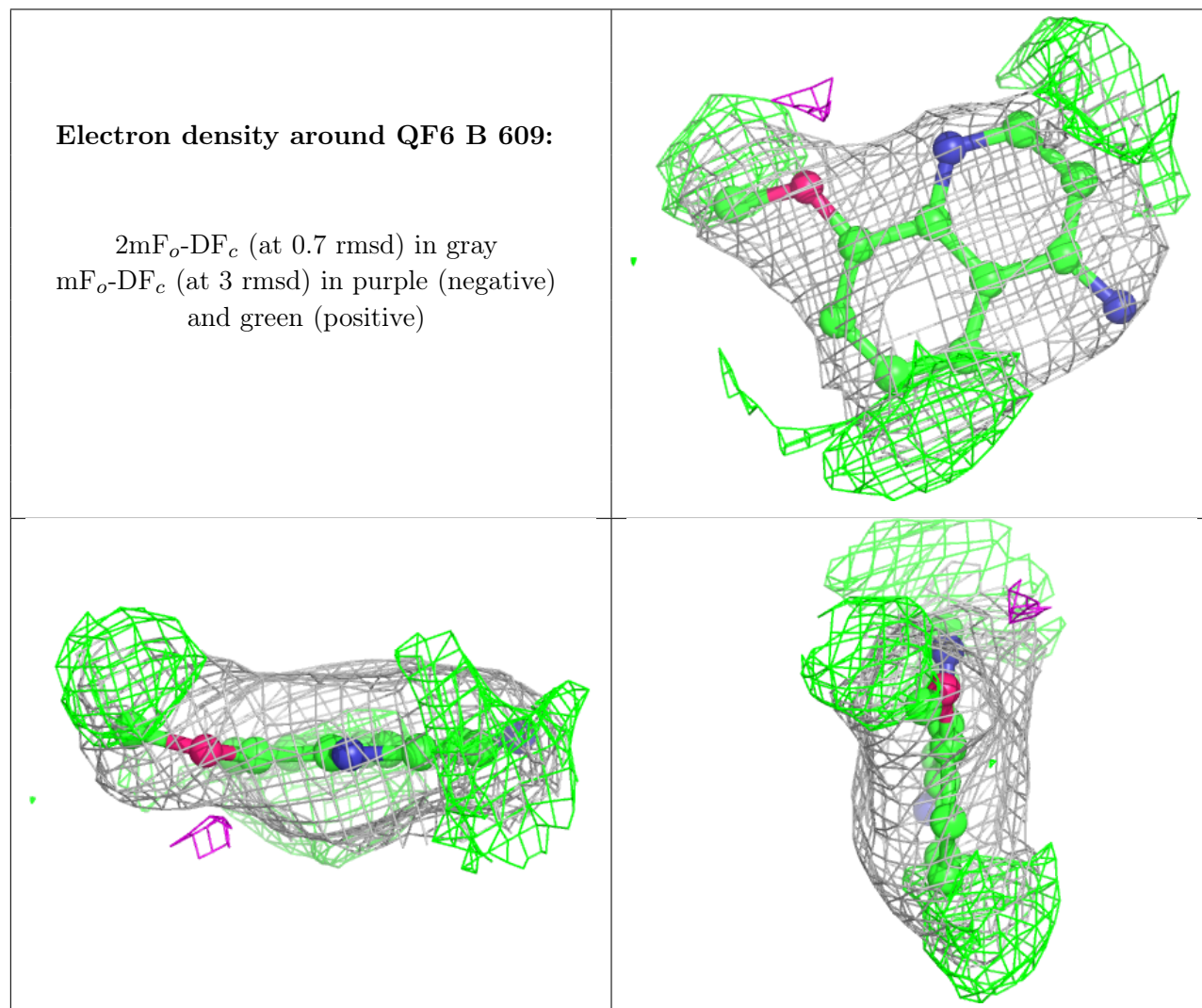
There are no monosaccharides in this entry.

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(Å ²)	Q<0.9
2	FMT	A	607	3/3	0.73	0.29	53,53,61,63	0
2	FMT	A	606	3/3	0.79	0.25	59,59,66,72	0
2	FMT	B	605	3/3	0.80	0.28	50,50,59,65	0
2	FMT	A	601	3/3	0.82	0.22	44,44,51,56	0
2	FMT	C	610	3/3	0.85	0.23	62,62,64,69	0
3	QF6	B	609	13/13	0.85	0.24	63,97,102,104	0
2	FMT	A	612	3/3	0.86	0.45	80,80,81,82	0
2	FMT	A	609	3/3	0.86	0.30	58,58,59,59	0
2	FMT	C	606	3/3	0.87	0.22	60,60,62,67	0
2	FMT	B	604	3/3	0.89	0.28	54,54,61,64	0
2	FMT	B	608	3/3	0.90	0.17	57,57,66,66	0
2	FMT	C	602	3/3	0.91	0.19	45,45,46,46	0
2	FMT	C	609	3/3	0.91	0.21	41,41,50,55	0
2	FMT	C	601	3/3	0.92	0.15	46,46,61,65	0
2	FMT	C	608	3/3	0.93	0.16	44,44,48,52	0
2	FMT	B	607	3/3	0.93	0.28	71,71,75,76	0
2	FMT	A	613	3/3	0.94	0.23	51,51,54,54	0
2	FMT	A	610	3/3	0.94	0.17	36,36,39,39	0
2	FMT	A	608	3/3	0.94	0.21	55,55,56,60	0
2	FMT	A	605	3/3	0.95	0.15	41,41,42,45	0
2	FMT	A	603	3/3	0.96	0.20	40,40,42,44	0
2	FMT	B	602	3/3	0.97	0.20	42,42,47,47	0
2	FMT	B	603	3/3	0.97	0.19	43,43,46,48	0
2	FMT	C	603	3/3	0.97	0.25	39,39,42,44	0
2	FMT	A	604	3/3	0.97	0.14	38,38,40,41	0
2	FMT	C	605	3/3	0.98	0.14	40,40,43,49	0
2	FMT	B	601	3/3	0.98	0.18	43,43,44,47	0
2	FMT	C	607	3/3	0.98	0.23	56,56,57,59	0
2	FMT	A	602	3/3	0.98	0.17	35,35,37,40	0
2	FMT	B	606	3/3	0.98	0.18	50,50,52,54	0
2	FMT	A	611	3/3	0.98	0.12	46,46,58,61	0
2	FMT	C	604	3/3	0.98	0.15	37,37,39,49	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.