



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 08:37 pm BST

PDB ID : 4BE0
Title : PFV intasome with inhibitor XZ-115
Authors : Hare, S.; Cherepanov, P.
Deposited on : 2012-10-08
Resolution : 2.68 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

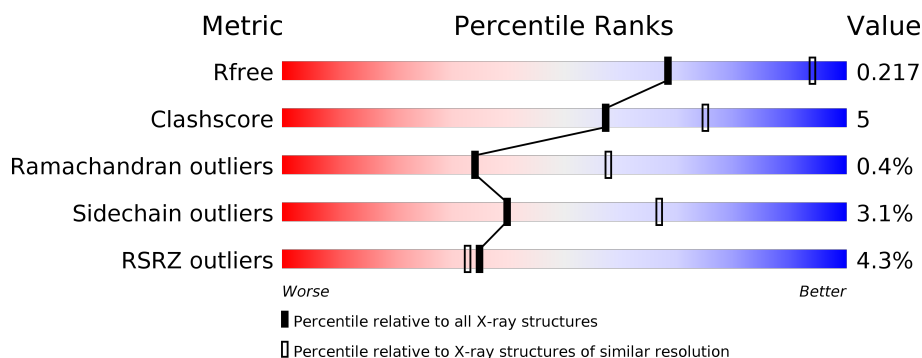
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.68 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	3863 (2.70-2.66)
Clashscore	141614	4210 (2.70-2.66)
Ramachandran outliers	138981	4141 (2.70-2.66)
Sidechain outliers	138945	4141 (2.70-2.66)
RSRZ outliers	127900	3780 (2.70-2.66)

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	395	<div> <div>3%</div> <div> <div></div> <div>79%</div> <div>13%</div> <div>7%</div> </div> </div>
1	B	395	<div> <div>3%</div> <div> <div></div> <div>40%</div> <div>7%</div> <div>53%</div> </div> </div>
2	C	19	<div> <div></div> <div> <div>63%</div> <div>21%</div> <div>16%</div> </div> </div>
3	D	17	<div> <div></div> <div> <div>82%</div> <div>18%</div> </div> </div>

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 5372 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 PFV INTEGRASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	368	Total	C	N	O	S	0	0	0
			2920	1872	512	532	4			
1	B	184	Total	C	N	O	S	0	0	0
			1423	922	233	267	1			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P14350
A	-1	PRO	-	expression tag	UNP P14350
A	0	GLY	-	expression tag	UNP P14350
A	217	SER	GLY	variant	UNP P14350
A	218	GLY	SER	variant	UNP P14350
B	-2	GLY	-	expression tag	UNP P14350
B	-1	PRO	-	expression tag	UNP P14350
B	0	GLY	-	expression tag	UNP P14350
B	217	SER	GLY	variant	UNP P14350
B	218	GLY	SER	variant	UNP P14350

- Molecule 2 is a DNA chain called 19 NUCLEOTIDE PREPROCESSED PFV DONOR DNA (NON-TRANSFERRED STRAND).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	19	Total	C	N	O	P	0	0	0
			387	187	68	114	18			

- Molecule 3 is a DNA chain called 17 NUCLEOTIDE PREPROCESSED PFV DONOR DNA (TRANSFERRED STRAND).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	17	Total	C	N	O	P	0	0	0
			345	166	65	98	16			

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).

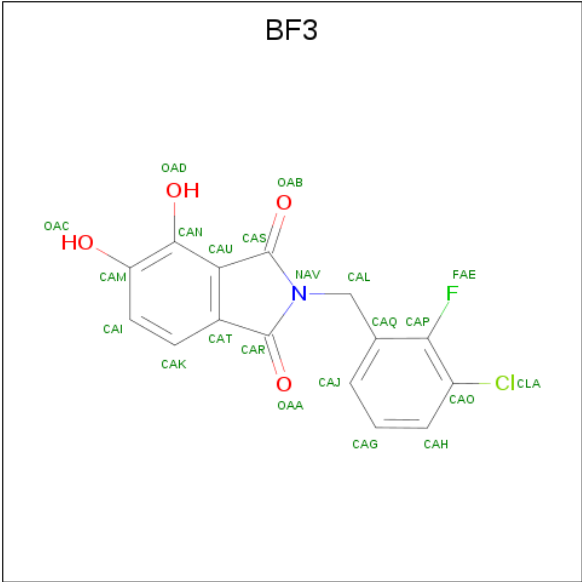


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	2	Total	Mg	0	0
			2	2		

- Molecule 8 is 2-(3-chloro-2-fluorobenzyl)-4,5-dihydroxy-1H-isoindole-1,3(2H)-dione (three-letter code: BF3) (formula: C₁₅H₉ClFNO₄).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
8	A	1	Total	C	Cl	F	N	O	0	0
			22	15	1	1	1	4		


- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	155	Total	O	0	0
			155	155		
9	B	31	Total	O	0	0
			31	31		
9	C	26	Total	O	0	0
			26	26		
9	D	15	Total	O	0	0
			15	15		

- Molecule 1: PFV INTEGRASE



- Molecule 3: 17 NUCLEOTIDE PREPROCESSED PFV DONOR DNA (TRANSFERRED STRAND)

Chain D:  82% 18%



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	160.12Å 160.12Å 124.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.22 – 2.68 39.22 – 2.68	Depositor EDS
% Data completeness (in resolution range)	98.7 (39.22-2.68) 98.8 (39.22-2.68)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.183 , 0.218 0.183 , 0.217	Depositor DCC
R_{free} test set	2287 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	63.9	Xtriage
Anisotropy	0.203	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 40.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5372	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.55% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: BF3, GOL, MG, ZN, SO4

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.65	1/2999 (0.0%)	0.74	1/4094 (0.0%)
1	B	0.66	2/1463 (0.1%)	0.71	0/2003
2	C	0.54	0/433	1.09	3/667 (0.4%)
3	D	0.49	0/387	1.04	2/595 (0.3%)
All	All	0.63	3/5282 (0.1%)	0.80	6/7359 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	154	TRP	CD2-CE2	5.89	1.48	1.41
1	A	81	TRP	CD2-CE2	5.09	1.47	1.41
1	B	242	TRP	CD2-CE2	5.01	1.47	1.41

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1	DT	P-O3'-C3'	7.94	129.22	119.70
2	C	1	DA	P-O3'-C3'	7.35	128.52	119.70
2	C	11	DA	P-O3'-C3'	6.47	127.47	119.70
3	D	7	DA	O3'-P-O5'	-5.99	92.62	104.00
1	A	69	ARG	NE-CZ-NH2	-5.49	117.56	120.30
2	C	2	DT	P-O3'-C3'	5.43	126.21	119.70

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	2920	0	2945	34	0
1	B	1423	0	1384	12	0
2	C	387	0	218	5	0
3	D	345	0	193	1	0
4	A	1	0	0	0	0
5	A	10	0	0	1	0
5	B	5	0	0	1	0
6	A	24	0	32	3	0
6	B	6	0	8	1	0
7	A	2	0	0	0	0
8	A	22	0	7	0	0
9	A	155	0	0	5	0
9	B	31	0	0	0	0
9	C	26	0	0	0	0
9	D	15	0	0	0	0
All	All	5372	0	4787	49	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 (49) 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:137:GLN:H	6:A:1378:GOL:H2	1.35	0.89
1:B:258:SER:O	1:B:261:LEU:O	1.98	0.80
1:A:137:GLN:N	6:A:1378:GOL:H2	2.03	0.73
1:A:219:LYS:HE3	2:C:6:DC:OP1	1.92	0.70
1:B:264:THR:OG1	1:B:267:GLN:HG3	1.98	0.63
1:A:162:SER:HB2	6:A:1377:GOL:H2	1.85	0.59
3:D:17:DA:N3	3:D:17:DA:H5"	2.16	0.58
1:A:69:ARG:HD2	1:A:70:GLU:OE2	2.04	0.58
1:A:161:PRO:O	1:A:189:ALA:HB2	2.06	0.55
1:A:156:TYR:OH	1:A:173:LEU:HB2	2.07	0.55
1:A:19:HIS:HD2	9:A:2001:HOH:O	1.89	0.54
1:A:315:TRP:CE2	1:A:371:PRO:HD3	2.42	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:GLN:HG3	1:A:21:ILE:HD11	1.90	0.54
1:A:111:PRO:O	1:A:350:ARG:HD3	2.08	0.54
2:C:1:DA:H2"	2:C:2:DT:H5'	1.89	0.53
1:B:212:TYR:O	1:B:214:PRO:HD3	2.09	0.53
1:B:283:THR:C	1:B:285:ASP:H	2.14	0.51
1:A:82:TRP:HB2	1:A:83:PRO:HD2	1.94	0.50
1:B:127:ILE:HA	1:B:144:VAL:O	2.11	0.50
1:A:108:ALA:O	1:A:314:SER:HA	2.13	0.49
1:A:81:TRP:CE3	1:A:82:TRP:HA	2.48	0.49
1:A:219:LYS:NZ	9:A:2082:HOH:O	2.46	0.48
1:A:9:ASP:O	1:A:13:ASP:HB2	2.13	0.48
1:A:97:GLN:OE1	1:A:339:LYS:HG2	2.14	0.47
1:A:325:GLU:HA	1:A:369:LEU:HD23	1.96	0.47
1:A:246:LEU:HB2	1:A:247:PRO:HD3	1.97	0.46
1:A:115:PRO:HG3	2:C:3:DT:C2	2.51	0.46
1:B:222:ARG:NE	5:B:1300:SO4:O4	2.46	0.46
1:A:73:LEU:HD22	1:A:86:ARG:CZ	2.46	0.45
1:A:357:HIS:HD2	9:A:2134:HOH:O	2.00	0.45
1:B:216:SER:HB2	1:B:217:SER:H	1.52	0.45
1:A:221:GLU:O	1:A:224:ASN:HB2	2.18	0.44
1:A:135:PRO:HG3	1:A:140:LEU:HD21	1.99	0.44
1:A:122:PHE:O	1:A:179:PRO:HA	2.17	0.44
1:A:295:LEU:HD21	1:A:299:ARG:NH2	2.32	0.43
1:A:357:HIS:HB2	5:A:1376:SO4:O4	2.18	0.43
1:A:222:ARG:HD3	2:C:5:DT:O2	2.19	0.43
1:A:357:HIS:CD2	9:A:2134:HOH:O	2.72	0.42
1:A:73:LEU:O	1:A:73:LEU:HG	2.18	0.42
1:A:221:GLU:OE1	1:A:221:GLU:HA	2.19	0.42
1:A:348:ASN:HB2	1:A:349:PRO:CD	2.49	0.42
2:C:10:DG:H2"	2:C:11:DA:C8	2.54	0.42
1:A:98:GLN:OE1	9:A:2022:HOH:O	2.21	0.42
1:B:267:GLN:HB3	6:B:1301:GOL:H31	2.02	0.41
1:A:337:TRP:CG	1:A:370:LYS:HE2	2.56	0.41
1:B:223:LYS:HA	1:B:223:LYS:HD3	1.88	0.41
1:B:212:TYR:C	1:B:214:PRO:HD3	2.41	0.41
1:B:258:SER:HA	1:B:259:PRO:HD2	1.93	0.41
1:B:161:PRO:O	1:B:189:ALA:HB2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	366/395 (93%)	346 (94%)	18 (5%)	2 (0%)	29	52
1	B	182/395 (46%)	170 (93%)	12 (7%)	0	100	100
All	All	548/790 (69%)	516 (94%)	30 (6%)	2 (0%)	34	58

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	201	GLU
1	A	14	GLN

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	327/354 (92%)	320 (98%)	7 (2%)	53	78
1	B	154/354 (44%)	146 (95%)	8 (5%)	23	46
All	All	481/708 (68%)	466 (97%)	15 (3%)	40	67

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	GLN
1	A	149	MET
1	A	307	THR
1	A	356	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	360	ASN
1	A	364	VAL
1	A	366	ILE
1	B	117	ARG
1	B	164	SER
1	B	235	LEU
1	B	274	SER
1	B	277	PRO
1	B	282	ASP
1	B	286	LEU
1	B	288	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	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](#)

Of 12 ligands modelled in this entry, 3 are monoatomic - leaving 9 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
6	GOL	B	1301	-	5,5,5	0.40	0	5,5,5	0.27	0
6	GOL	A	1377	-	5,5,5	0.40	0	5,5,5	0.69	0
5	SO4	B	1300	-	4,4,4	0.38	0	6,6,6	0.23	0
6	GOL	A	1384	-	5,5,5	0.38	0	5,5,5	0.48	0
6	GOL	A	1379	-	5,5,5	0.36	0	5,5,5	0.52	0
5	SO4	A	1383	-	4,4,4	0.53	0	6,6,6	0.33	0
8	BF3	A	1382	7	24,24,24	2.21	4 (16%)	35,36,36	2.44	10 (28%)
6	GOL	A	1378	-	5,5,5	0.44	0	5,5,5	1.37	1 (20%)
5	SO4	A	1376	-	4,4,4	0.43	0	6,6,6	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
6	GOL	B	1301	-	-	0/4/4/4	-
6	GOL	A	1377	-	-	4/4/4/4	-
6	GOL	A	1384	-	-	2/4/4/4	-
6	GOL	A	1379	-	-	0/4/4/4	-
8	BF3	A	1382	7	-	0/4/20/20	0/3/3/3
6	GOL	A	1378	-	-	3/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	1382	BF3	CAU-CAS	-5.38	1.39	1.49
8	A	1382	BF3	CAT-CAR	-5.28	1.40	1.48
8	A	1382	BF3	CAR-NAV	-4.90	1.33	1.39
8	A	1382	BF3	CAS-NAV	-4.90	1.33	1.39

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	1382	BF3	CAP-CAO-CLA	-7.43	116.64	119.50
8	A	1382	BF3	CAU-CAS-NAV	6.20	108.81	105.56
8	A	1382	BF3	CAT-CAR-NAV	5.27	109.62	105.88
8	A	1382	BF3	CAJ-CAQ-CAP	4.17	119.46	116.43
8	A	1382	BF3	CAS-NAV-CAR	-3.33	109.31	112.03
8	A	1382	BF3	CAU-CAT-CAR	-2.97	105.32	108.32
8	A	1382	BF3	OAB-CAS-CAU	-2.58	125.42	129.09

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	1382	BF3	CAK-CAT-CAR	2.57	133.82	129.63
6	A	1378	GOL	O3-C3-C2	-2.37	98.85	110.20
8	A	1382	BF3	CAH-CAO-CLA	2.30	123.01	118.41
8	A	1382	BF3	CAN-CAU-CAS	2.08	133.16	129.48

There are no chirality outliers.

All (9) torsion outliers are listed below:

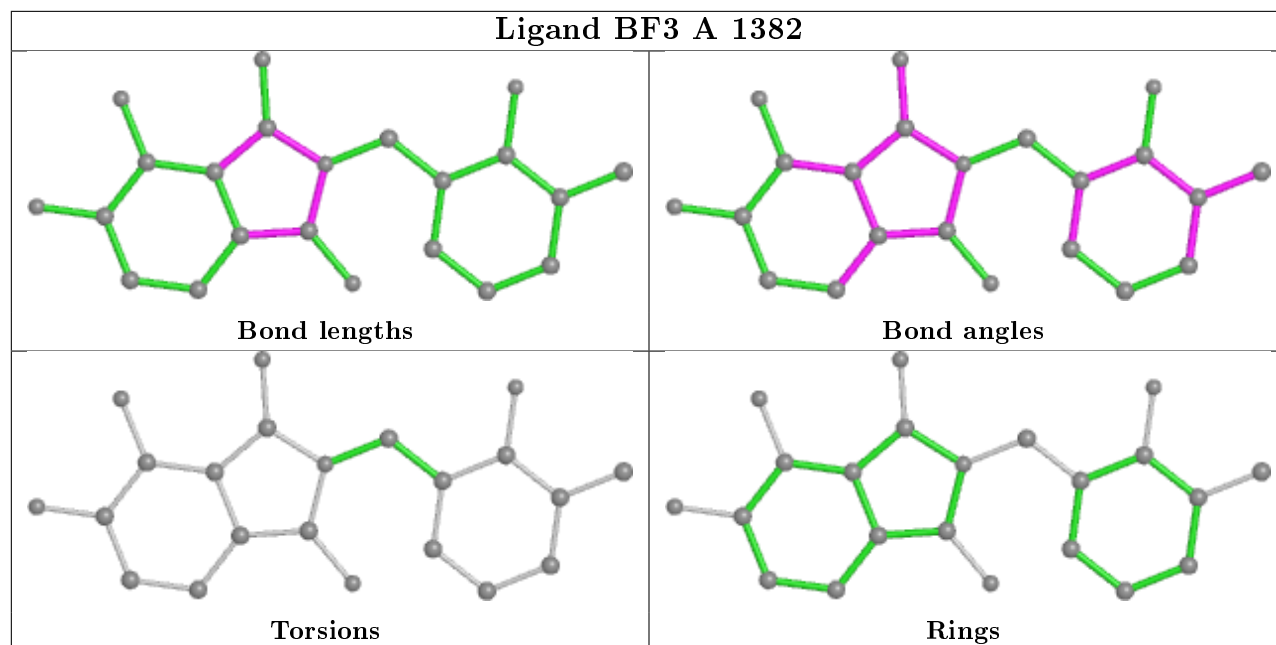
Mol	Chain	Res	Type	Atoms
6	A	1377	GOL	O1-C1-C2-O2
6	A	1377	GOL	O1-C1-C2-C3
6	A	1384	GOL	O1-C1-C2-O2
6	A	1384	GOL	O1-C1-C2-C3
6	A	1378	GOL	C1-C2-C3-O3
6	A	1378	GOL	O2-C2-C3-O3
6	A	1377	GOL	O2-C2-C3-O3
6	A	1377	GOL	C1-C2-C3-O3
6	A	1378	GOL	O1-C1-C2-C3

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	1301	GOL	1	0
6	A	1377	GOL	1	0
5	B	1300	SO4	1	0
6	A	1378	GOL	2	0
5	A	1376	SO4	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 ⓘ

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	368/395 (93%)	-0.13	12 (3%) 46 45	38, 52, 95, 135	0
1	B	184/395 (46%)	-0.09	13 (7%) 16 13	40, 61, 138, 167	0
2	C	19/19 (100%)	-0.27	0 100 100	43, 56, 85, 114	0
3	D	17/17 (100%)	-0.42	0 100 100	43, 50, 110, 112	0
All	All	588/826 (71%)	-0.13	25 (4%) 35 33	38, 54, 113, 167	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	296	GLN	4.4
1	A	17	GLN	4.3
1	B	280	ASN	3.9
1	B	279	ALA	3.3
1	B	292	LEU	3.3
1	A	8	LEU	3.2
1	B	293	SER	3.0
1	B	295	LEU	2.9
1	A	19	HIS	2.9
1	A	16	LEU	2.5
1	B	288	ARG	2.4
1	A	9	ASP	2.4
1	A	10	ALA	2.3
1	A	21	ILE	2.2
1	A	15	LEU	2.2
1	B	294	LEU	2.1
1	A	18	GLY	2.1
1	A	358	LEU	2.1
1	B	286	LEU	2.1
1	A	29	THR	2.1
1	B	299	ARG	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	13	ASP	2.1
1	B	260	VAL	2.1
1	B	298	ILE	2.0
1	B	297	GLU	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 carbohydrates in this entry.

6.4 Ligands [i](#)

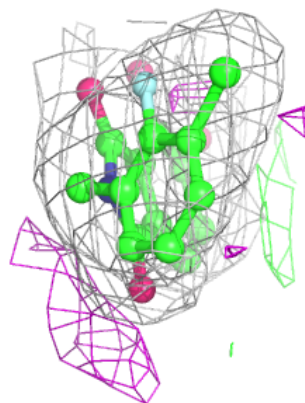
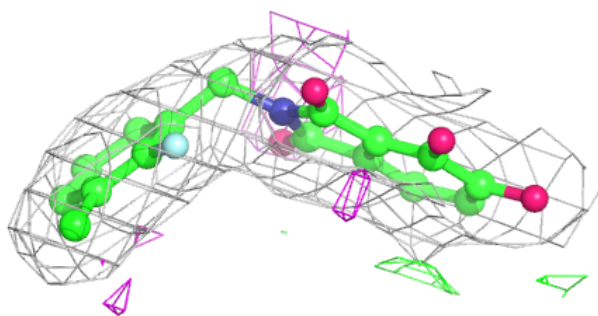
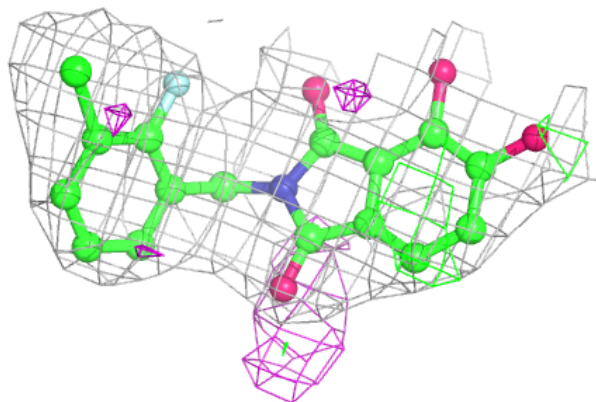
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	SO4	A	1383	5/5	0.82	0.33	59,65,74,75	5
6	GOL	A	1379	6/6	0.86	0.22	84,90,93,94	0
6	GOL	A	1384	6/6	0.88	0.24	69,92,94,104	0
6	GOL	A	1377	6/6	0.89	0.19	69,81,93,96	0
6	GOL	B	1301	6/6	0.89	0.30	61,70,75,86	0
6	GOL	A	1378	6/6	0.92	0.16	61,66,82,94	0
5	SO4	B	1300	5/5	0.96	0.17	87,92,96,101	0
5	SO4	A	1376	5/5	0.96	0.15	85,91,93,94	0
8	BF3	A	1382	22/22	0.97	0.13	38,48,52,52	0
7	MG	A	1381	1/1	0.99	0.08	44,44,44,44	0
7	MG	A	1380	1/1	0.99	0.06	40,40,40,40	0
4	ZN	A	393	1/1	1.00	0.12	46,46,46,46	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 BF3 A 1382:

$2mF_o - DF_c$ (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.