



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:10 AM UTC

PDB ID : 4BE1 / pdb_00004be1
Title : PFV intasome with inhibitor XZ-116
Authors : Hare, S.; Cherepanov, P.
Deposited on : 2012-10-08
Resolution : 2.71 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

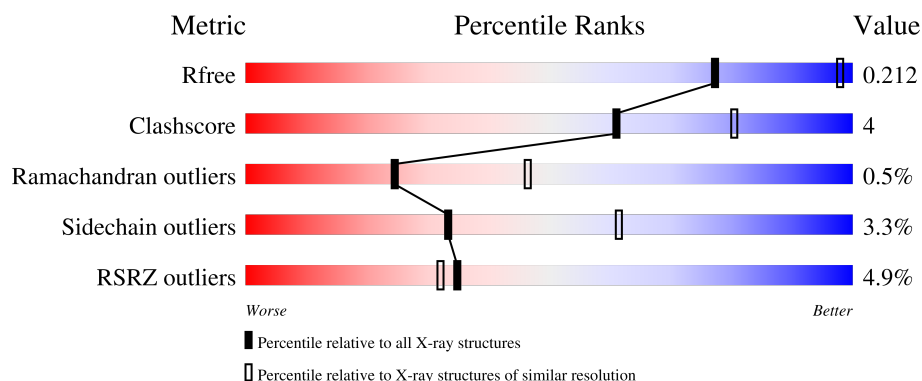
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.71 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4348 (2.74-2.70)
Clashscore	190562	4665 (2.74-2.70)
Ramachandran outliers	187476	4584 (2.74-2.70)
Sidechain outliers	187428	4585 (2.74-2.70)
RSRZ outliers	180081	4348 (2.74-2.70)

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	395	2% 77% 15% 7%
1	B	395	5% 42% 5% 53%
2	C	19	5% 68% 26% 5%
3	D	17	12% 76% 18% 6%

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 5369 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 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
			1417	919	230	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			

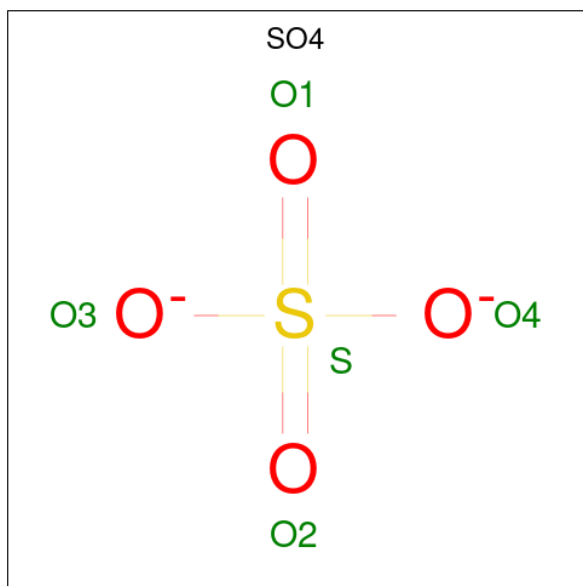
- Molecule 3 is a DNA chain called 17 NUCLEOTIDE PREPROCESSED PFV DONOR DNA (TRANSFERRED STRAND).

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 (CCD ID: ZN) (formula: Zn).

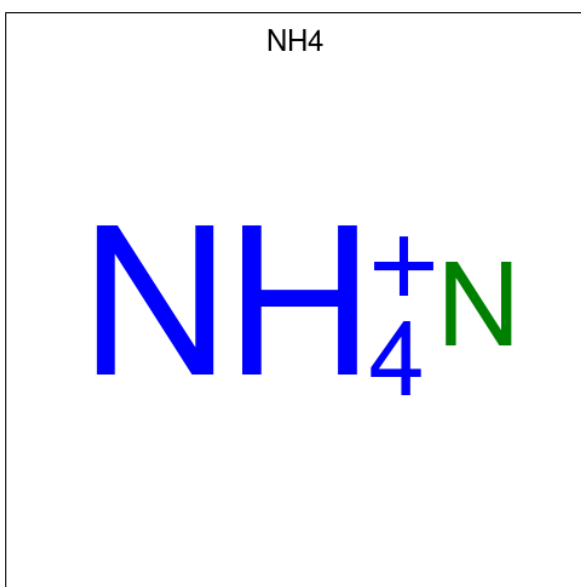
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



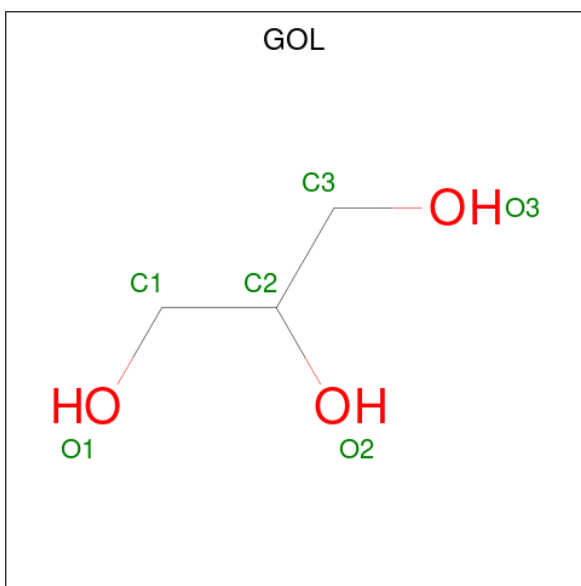
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is AMMONIUM ION (CCD ID: NH4) (formula: H₄N).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	N		0	0
			1	1			

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

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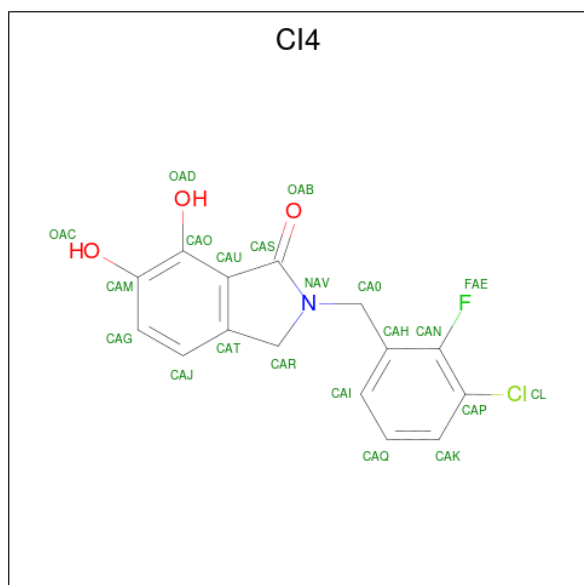
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 8 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	2	Total	Mg	0	0
			2	2		
8	B	1	Total	Mg	0	0
			1	1		

- Molecule 9 is 2-(3-chloro-2-fluorobenzyl)-6,7-dihydroxy-2,3-dihydro-1H-isoindol-1-one (CCD ID: CI4) (formula: C₁₅H₁₁ClFNO₃).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
9	A	1	Total	C	Cl	F	N	O	0	0
			21	15	1	1	1	3		

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	163	Total	O	0	0
			163	163		
10	B	37	Total	O	0	0
			37	37		

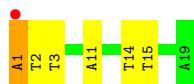
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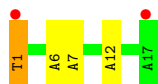
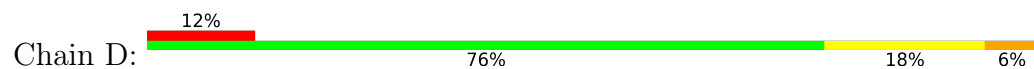
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	C	18	Total	O	0	0
			18	18		
10	D	17	Total	O	0	0
			17	17		

- Molecule 1: INTEGRASE





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



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	161.03Å 161.03Å 123.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.92 – 2.71 50.92 – 2.71	Depositor EDS
% Data completeness (in resolution range)	97.9 (50.92-2.71) 97.9 (50.92-2.71)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.73Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.180 , 0.213 0.178 , 0.212	Depositor DCC
R_{free} test set	2215 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	63.1	Xtriage
Anisotropy	0.293	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5369	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.31% 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: NH4, ZN, MG, SO4, CI4, GOL

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	0.90	3/2999 (0.1%)	1.02	5/4094 (0.1%)
1	B	0.86	0/1457	1.01	1/1996 (0.1%)
2	C	0.46	0/433	1.21	5/667 (0.7%)
3	D	0.40	0/387	1.19	4/595 (0.7%)
All	All	0.83	3/5276 (0.1%)	1.05	15/7352 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	213	HIS	CG-CD2	5.78	1.42	1.35
1	A	62	HIS	CG-CD2	5.72	1.42	1.35
1	A	357	HIS	CG-CD2	5.03	1.41	1.35

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1	DT	P-O3'-C3'	7.29	131.14	120.20
2	C	1	DA	P-O3'-C3'	6.43	129.84	120.20
2	C	11	DA	C4'-C3'-O3'	-6.28	100.57	110.00
1	A	31	PHE	N-CA-C	6.24	119.53	109.86
3	D	12	DA	C4'-C3'-O3'	-6.18	100.73	110.00
1	A	356	ASP	N-CA-C	6.13	118.98	109.60
3	D	6	DA	O3'-P-O5'	-5.98	95.02	104.00
2	C	15	DT	C4'-C3'-O3'	-5.79	101.32	110.00
3	D	7	DA	O3'-P-O5'	-5.79	95.32	104.00
2	C	14	DT	C4'-C3'-O3'	-5.57	101.65	110.00
1	A	348	ASN	CA-C-N	-5.52	114.13	119.87
1	A	348	ASN	C-N-CA	-5.52	114.13	119.87
1	A	347	LEU	N-CA-C	5.28	117.72	111.33
2	C	1	DA	O3'-P-O5'	-5.16	96.26	104.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	217	SER	N-CA-C	-5.01	100.85	108.76

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	30	0
1	B	1417	0	1373	9	0
2	C	387	0	218	3	0
3	D	345	0	193	1	0
4	A	1	0	0	0	0
5	A	10	0	0	0	0
5	B	5	0	0	1	0
6	A	1	0	0	0	0
7	A	18	0	24	3	0
7	B	6	0	8	0	0
8	A	2	0	0	0	0
8	B	1	0	0	0	0
9	A	21	0	9	0	0
10	A	163	0	0	2	1
10	B	37	0	0	0	0
10	C	18	0	0	0	0
10	D	17	0	0	0	0
All	All	5369	0	4770	42	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 (42) 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:258:SER:O	1:B:261:LEU:O	2.11	0.68
2:C:1:DA:H2''	2:C:2:DT:H5'	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:162:SER:HB2	7:A:1378:GOL:H2	1.79	0.64
1:A:69:ARG:HD2	1:A:70:GLU:OE2	2.01	0.61
1:A:51:GLN:HA	1:A:54:ARG:HD2	1.83	0.59
1:A:137:GLN:HG2	7:A:1379:GOL:H32	1.83	0.59
1:A:12:LEU:HD23	1:A:37:VAL:HG21	1.84	0.58
1:B:127:ILE:HA	1:B:144:VAL:O	2.04	0.57
1:A:219:LYS:NZ	10:A:2090:HOH:O	2.40	0.55
1:A:161:PRO:O	1:A:189:ALA:HB2	2.07	0.54
1:A:221:GLU:HA	1:A:221:GLU:OE1	2.08	0.53
1:A:108:ALA:O	1:A:314:SER:HA	2.08	0.53
1:A:73:LEU:HD22	1:A:86:ARG:CZ	2.40	0.52
1:A:136:SER:HA	7:A:1379:GOL:H2	1.92	0.52
1:A:246:LEU:HB2	1:A:247:PRO:HD3	1.92	0.51
1:A:292:LEU:O	1:A:296:GLN:HG3	2.11	0.51
1:A:315:TRP:CE2	1:A:371:PRO:HD3	2.48	0.48
1:A:97:GLN:NE2	1:A:339:LYS:HG2	2.29	0.48
1:A:324:GLN:HG2	1:A:340:PRO:HA	1.94	0.48
10:A:2070:HOH:O	1:B:266:HIS:HD2	1.96	0.48
1:A:111:PRO:O	1:A:350:ARG:HD3	2.13	0.47
1:A:112:ILE:O	1:A:307:THR:HG22	2.15	0.47
1:A:325:GLU:HA	1:A:369:LEU:HD23	1.97	0.46
1:A:348:ASN:HB2	1:A:349:PRO:CD	2.45	0.46
1:A:115:PRO:HG3	2:C:3:DT:C2	2.50	0.46
1:A:356:ASP:C	1:A:356:ASP:OD1	2.58	0.46
1:B:141:TYR:CZ	1:B:161:PRO:HD3	2.51	0.46
1:B:280:ASN:O	1:B:281:GLN:C	2.58	0.45
1:A:13:ASP:C	1:A:15:LEU:H	2.25	0.45
1:A:84:ASN:HB3	1:A:87:LYS:HG3	1.99	0.44
2:C:1:DA:C2'	2:C:2:DT:H5'	2.46	0.44
1:A:122:PHE:O	1:A:179:PRO:HA	2.17	0.43
1:A:295:LEU:HD21	1:A:299:ARG:NH2	2.33	0.43
1:B:222:ARG:NE	5:B:1300:SO4:O4	2.41	0.43
1:A:97:GLN:HE21	1:A:339:LYS:HG2	1.83	0.43
3:D:1:DT:O2	3:D:1:DT:O4'	2.38	0.42
1:B:243:TYR:C	1:B:243:TYR:CD1	2.97	0.42
1:A:12:LEU:HD23	1:A:37:VAL:CG2	2.50	0.42
1:A:344:LEU:HD11	1:A:355:LEU:HB2	2.02	0.42
1:B:231:LEU:HD23	1:B:231:LEU:HA	1.95	0.41
1:A:183:HIS:HA	1:A:207:GLU:O	2.20	0.41
1:B:134:PRO:HB2	1:B:239:PRO:O	2.21	0.41

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:A:2012:HOH:O	10:A:2109:HOH:O[8_554]	2.08	0.12

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	366/395 (93%)	345 (94%)	18 (5%)	3 (1%)	16	35
1	B	182/395 (46%)	172 (94%)	10 (6%)	0	100	100
All	All	548/790 (69%)	517 (94%)	28 (5%)	3 (0%)	24	46

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	201	GLU
1	A	14	GLN
1	A	359	GLY

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	327/354 (92%)	316 (97%)	11 (3%)	32	60
1	B	153/354 (43%)	148 (97%)	5 (3%)	33	61
All	All	480/708 (68%)	464 (97%)	16 (3%)	33	61

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	LYS
1	A	40	SER
1	A	55	GLN
1	A	137	GLN
1	A	178	ILE
1	A	219	LYS
1	A	284	LEU
1	A	318	VAL
1	A	351	THR
1	A	356	ASP
1	A	366	ILE
1	B	117	ARG
1	B	164	SER
1	B	235	LEU
1	B	256	THR
1	B	274	SER

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

Mol	Chain	Res	Type
1	A	255	ASN
1	A	296	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 13 ligands modelled in this entry, 4 are monoatomic and 1 is modelled with single atom - leaving 8 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	$\# Z > 2$
9	CI4	A	1383	8	23,23,23	1.42	2 (8%)	30,34,34	1.99	5 (16%)
7	GOL	A	1378	-	5,5,5	0.52	0	5,5,5	0.79	0
7	GOL	A	1379	-	5,5,5	0.23	0	5,5,5	0.84	0
7	GOL	A	1380	-	5,5,5	0.41	0	5,5,5	0.42	0
5	SO4	A	1384	-	4,4,4	0.55	0	6,6,6	0.25	0
5	SO4	B	1300	-	4,4,4	0.57	0	6,6,6	0.34	0
5	SO4	A	1376	-	4,4,4	0.50	0	6,6,6	0.21	0
7	GOL	B	1301	-	5,5,5	0.46	0	5,5,5	0.35	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
9	CI4	A	1383	8	-	0/4/16/16	0/3/3/3
7	GOL	A	1378	-	-	4/4/4/4	-
7	GOL	A	1379	-	-	3/4/4/4	-
7	GOL	A	1380	-	-	1/4/4/4	-
7	GOL	B	1301	-	-	0/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	1383	CI4	CAU-CAS	-4.67	1.41	1.49
9	A	1383	CI4	CAR-CAT	2.59	1.53	1.50

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	1383	CI4	CAU-CAS-NAV	7.37	111.73	106.40
9	A	1383	CI4	CAN-CAP-CL	-3.41	117.25	119.44
9	A	1383	CI4	CAK-CAP-CAN	3.35	121.35	119.37
9	A	1383	CI4	CAR-CAT-CAU	-3.12	107.69	109.96
9	A	1383	CI4	OAB-CAS-CAU	-2.12	125.93	129.15

There are no chirality outliers.

All (8) torsion outliers are listed below:

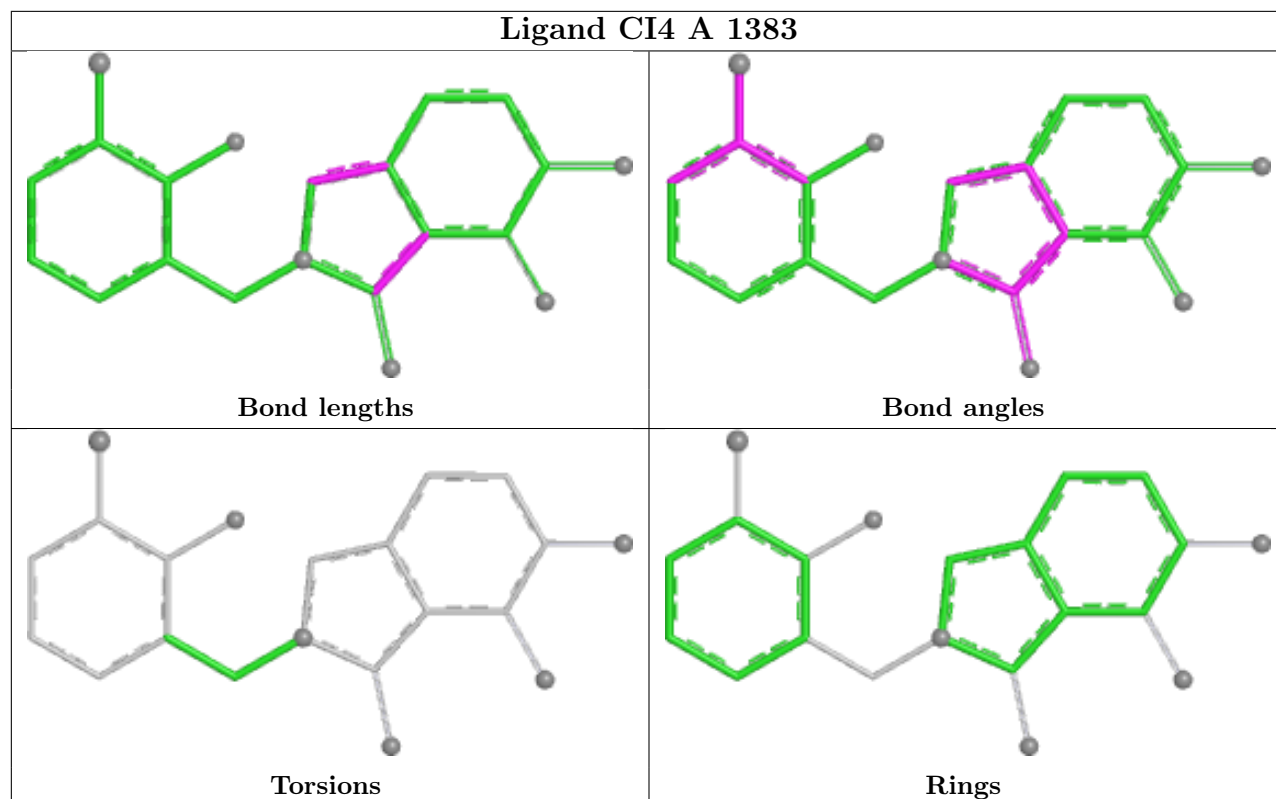
Mol	Chain	Res	Type	Atoms
7	A	1378	GOL	C1-C2-C3-O3
7	A	1379	GOL	O2-C2-C3-O3
7	A	1378	GOL	O1-C1-C2-C3
7	A	1379	GOL	O1-C1-C2-C3
7	A	1379	GOL	C1-C2-C3-O3
7	A	1378	GOL	O1-C1-C2-O2
7	A	1378	GOL	O2-C2-C3-O3
7	A	1380	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	1378	GOL	1	0
7	A	1379	GOL	2	0
5	B	1300	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%)	-0.45	7 (1%) 66 63	37, 53, 96, 136	0
1	B	184/395 (46%)	0.19	19 (10%) 12 10	46, 62, 143, 175	0
2	C	19/19 (100%)	-0.49	1 (5%) 32 28	44, 57, 88, 123	0
3	D	17/17 (100%)	-0.59	2 (11%) 9 8	45, 51, 113, 114	0
All	All	588/826 (71%)	-0.26	29 (4%) 35 32	37, 56, 114, 175	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	8	LEU	6.3
1	B	288	ARG	6.2
1	B	281	GLN	5.3
1	B	295	LEU	4.8
2	C	1	DA	4.2
1	B	299	ARG	4.0
1	B	297	GLU	3.8
1	B	279	ALA	3.7
1	A	358	LEU	3.6
1	B	286	LEU	3.6
1	A	10	ALA	3.6
1	B	117	ARG	3.2
1	B	292	LEU	3.2
1	B	213	HIS	3.1
1	B	296	GLN	3.1
1	B	298	ILE	2.9
1	A	357	HIS	2.9
1	B	234	LEU	2.9
1	B	257	TYR	2.9
1	B	287	THR	2.6
3	D	1	DT	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	235	LEU	2.4
3	D	17	DA	2.4
1	A	29	THR	2.3
1	B	283	THR	2.3
1	B	261	LEU	2.3
1	B	294	LEU	2.3
1	A	19	HIS	2.2
1	A	355	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

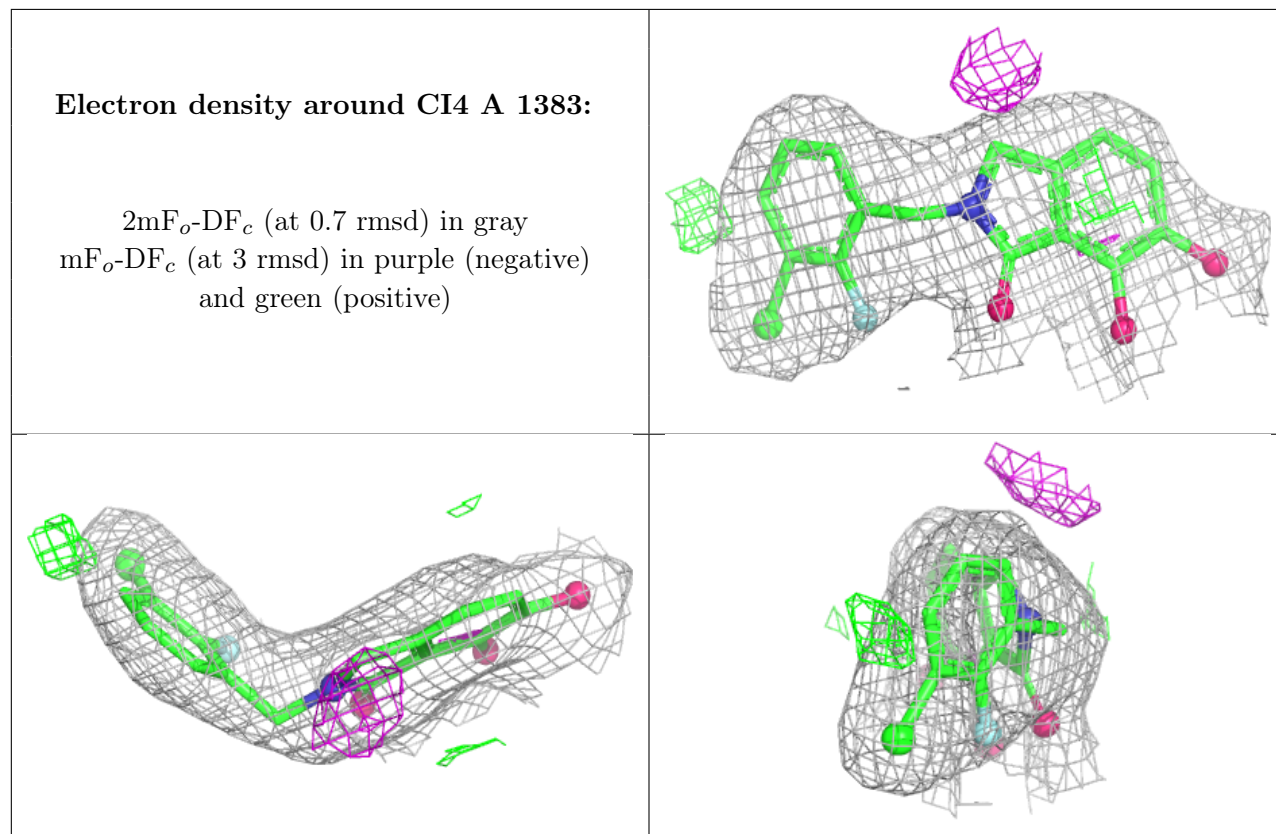
6.4 Ligands ⓘ

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
7	GOL	A	1378	6/6	0.88	0.21	77,89,95,95	0
7	GOL	A	1379	6/6	0.92	0.21	69,77,80,89	0
5	SO4	B	1300	5/5	0.93	0.19	78,82,92,100	0
7	GOL	A	1380	6/6	0.95	0.14	76,89,100,105	0
8	MG	B	1302	1/1	0.96	0.07	88,88,88,88	0
5	SO4	A	1376	5/5	0.97	0.10	90,93,95,98	0
6	NH4	A	1377	1/1	0.97	0.23	32,32,32,32	0
7	GOL	B	1301	6/6	0.97	0.14	57,69,75,79	0
5	SO4	A	1384	5/5	0.97	0.17	80,85,87,93	0
9	CI4	A	1383	21/21	0.97	0.07	38,48,53,55	0
8	MG	A	1382	1/1	1.00	0.04	53,53,53,53	0
4	ZN	A	393	1/1	1.00	0.01	51,51,51,51	0
8	MG	A	1381	1/1	1.00	0.02	47,47,47,47	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.