



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:25 AM GMT

PDB ID : 3OYN  
Title : Crystal structure of the PFV N224H mutant intasome bound to magnesium and the INSTI MK2048  
Authors : Hare, S.; Cherepanov, P.  
Deposited on : 2010-09-23  
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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

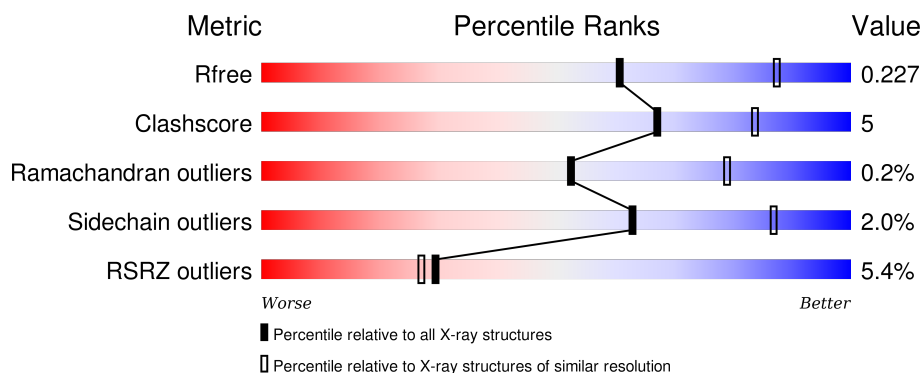
# 1 Overall quality at a glance

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))
$R_{free}$	91344	2780 (2.70-2.66)
Clashscore	102246	3138 (2.70-2.66)
Ramachandran outliers	100387	3089 (2.70-2.66)
Sidechain outliers	100360	3089 (2.70-2.66)
RSRZ outliers	91569	2789 (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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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>81%</div> <div>12%</div> <div>7%</div> </div>
1	B	395	<div> <div>5%</div> <div>41%</div> <div>6%</div> <div>53%</div> </div>
2	C	19	<div> <div>53%</div> <div>42%</div> <div>5%</div> </div>
3	D	17	<div> <div>59%</div> <div>29%</div> <div>12%</div> </div>

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
6	GOL	A	803	-	-	-	X
6	GOL	A	805	-	-	-	X

## 2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 5368 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
			2922	1874	513	531	4			
1	B	184	Total	C	N	O	S	0	0	0
			1450	941	236	272	1			

There are 12 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
A	224	HIS	ASN	ENGINEERED MUTATION	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
B	224	HIS	ASN	ENGINEERED MUTATION	UNP P14350

- Molecule 2 is a DNA chain called DNA (5'-D(\*AP\*TP\*TP\*GP\*TP\*CP\*AP\*TP\*GP\*GP\*AP\*AP\*TP\*TP\*TP\*CP\*GP\*CP\*A)-3').

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 DNA (5'-D(\*TP\*GP\*CP\*GP\*AP\*AP\*AP\*TP\*TP\*CP\*CP\*AP\*TP\*GP\*AP\*CP\*A)-3').

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: O<sub>4</sub>S).



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: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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		

- Molecule 7 is AMMONIUM ION (three-letter code: NH4) (formula: H<sub>4</sub>N).

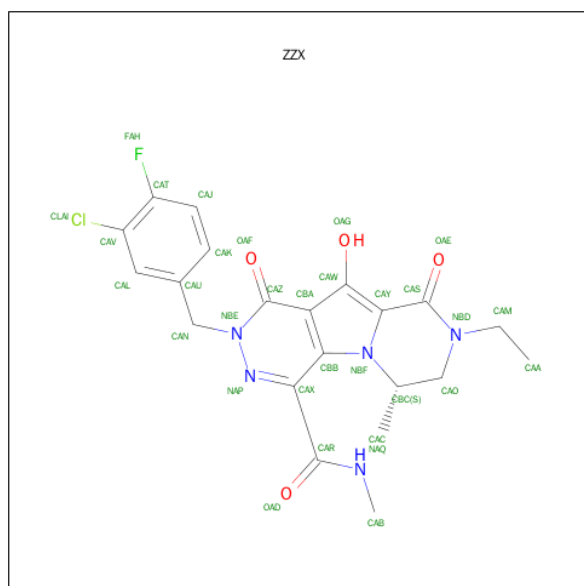


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total N 1 1	0	0

- Molecule 8 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 9 is (6S)-2-(3-CHLORO-4-FLUOROBENZYL)-8-ETHYL-10-HYDROXY-N,6-DIMETHYL-1,9-DIOXO-1,2,6,7,8,9-HEXAHYDROPYRAZINO[1',2':1,5]PYRROLO[2,3-D]PYRIDAZINE-4-CARBOXAMIDE (three-letter code: ZZX) (formula: C<sub>21</sub>H<sub>21</sub>ClFN<sub>5</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
9	A	1	Total	C	Cl	F	N	O	0	0
			32	21	1	1	5	4		

- Molecule 10 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	118	Total O 118 118	0	0
10	B	30	Total O 30 30	0	0
10	C	28	Total O 28 28	0	0

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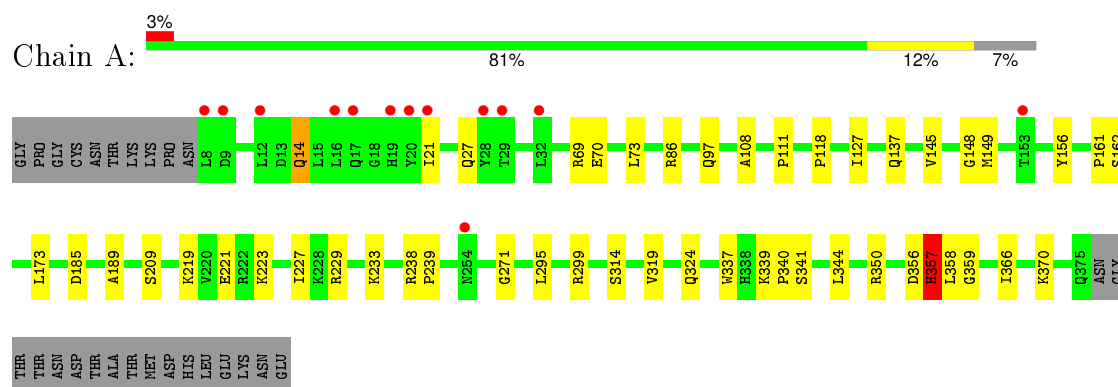
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	D	13	Total	O	0	0
			13	13		



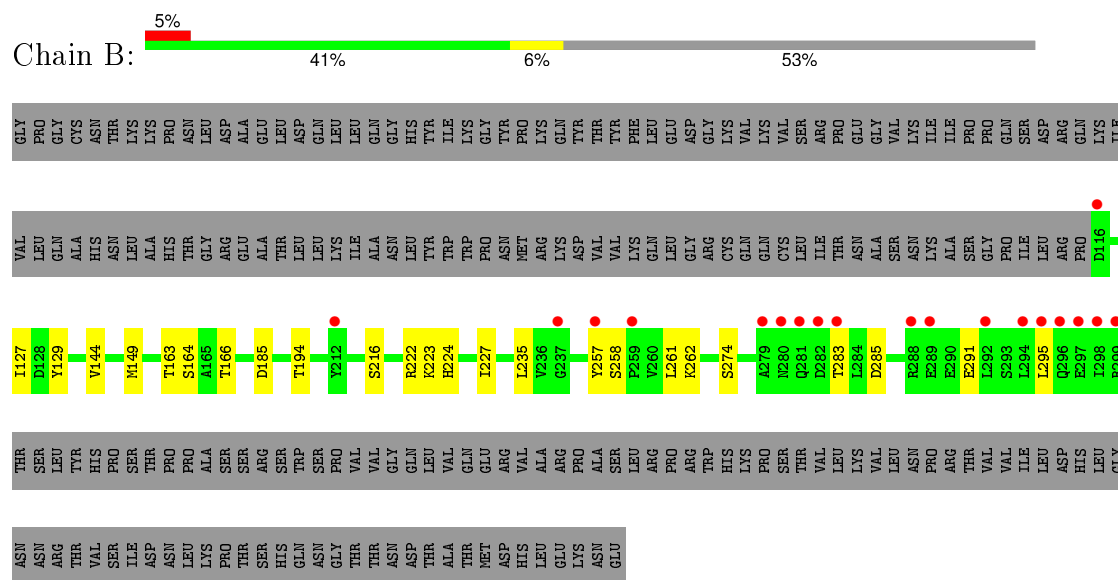
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: PFV integrase



#### • Molecule 1: PFV integrase

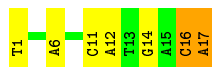


#### • Molecule 2: DNA (5'-D(\*AP\*TP\*TP\*GP\*TP\*CP\*AP\*TP\*GP\*GP\*AP\*AP\*TP\*TP\*TP\*CP\*GP\*CP\*A)-3')



- Molecule 3: DNA (5'-D(\*TP\*GP\*CP\*GP\*AP\*AP\*AP\*TP\*TP\*CP\*CP\*AP\*TP\*GP\*AP\*CP\*A)-3')

Chain D: 59% 29% 12%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	160.67Å 160.67Å 123.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.97 – 2.68 38.18 – 2.68	Depositor EDS
% Data completeness (in resolution range)	97.2 (38.97-2.68) 97.2 (38.18-2.68)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.81 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.208 , 0.231 0.202 , 0.227	Depositor DCC
$R_{free}$ test set	2253 reflections (5.34%)	DCC
Wilson B-factor (Å <sup>2</sup> )	69.6	Xtriage
Anisotropy	0.299	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 46.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 44515 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5368	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: GOL, MG, ZZX, ZN, NH4, 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	0.52	0/3002	0.65	0/4098
1	B	0.57	0/1491	0.62	0/2039
2	C	1.03	0/433	1.79	11/667 (1.6%)
3	D	1.11	1/387 (0.3%)	1.73	8/595 (1.3%)
All	All	0.65	1/5313 (0.0%)	0.93	19/7399 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	17	DA	N9-C4	5.34	1.41	1.37

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	8	DT	O4'-C4'-C3'	-8.76	100.74	106.00
3	D	17	DA	O4'-C1'-N9	7.56	113.29	108.00
2	C	12	DA	O4'-C1'-N9	-7.30	102.89	108.00
3	D	11	DC	O4'-C1'-N1	7.25	113.08	108.00
3	D	1	DT	P-O3'-C3'	7.18	128.32	119.70
3	D	14	DG	O4'-C1'-N9	-6.97	103.12	108.00
2	C	3	DT	C5-C4-O4	-6.79	120.14	124.90
3	D	16	DC	O4'-C4'-C3'	-6.34	101.96	104.50
3	D	12	DA	O4'-C1'-N9	-6.23	103.64	108.00
2	C	17	DG	O4'-C1'-N9	-6.06	103.75	108.00
2	C	1	DA	P-O3'-C3'	6.00	126.91	119.70
2	C	5	DT	P-O3'-C3'	5.67	126.50	119.70
2	C	2	DT	C4-C5-C7	5.56	122.34	119.00
2	C	8	DT	P-O3'-C3'	-5.42	113.20	119.70
3	D	6	DA	O4'-C1'-N9	-5.38	104.23	108.00
2	C	13	DT	O4'-C1'-N1	-5.30	104.29	108.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	17	DG	P-O3'-C3'	5.26	126.01	119.70
2	C	5	DT	O4'-C1'-N1	-5.18	104.37	108.00
3	D	6	DA	P-O3'-C3'	5.10	125.82	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	2922	0	2946	31	0
1	B	1450	0	1432	11	0
2	C	387	0	218	2	0
3	D	345	0	193	3	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	2	0
7	A	1	0	0	0	0
8	A	2	0	0	0	0
9	A	32	0	20	2	0
10	A	118	0	0	2	1
10	B	30	0	0	0	0
10	C	28	0	0	1	0
10	D	13	0	0	0	1
All	All	5368	0	4841	48	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 5.

All (48) 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:802:GOL:H2	1.45	0.81
1:A:219:LYS:HE3	2:C:6:DC:OP1	1.91	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:358:LEU:N	1:A:359:GLY:HA2	2.06	0.69
5:A:399:SO4:O2	10:A:515:HOH:O	2.11	0.68
1:A:97:GLN:OE1	10:A:491:HOH:O	2.11	0.68
3:D:16:DC:H3'	3:D:17:DA:H5'	1.80	0.63
2:C:1:DA:N7	10:C:319:HOH:O	2.29	0.63
1:B:258:SER:O	1:B:261:LEU:O	2.16	0.63
9:A:398:ZZX:HANA	3:D:17:DA:H3'	1.82	0.61
1:B:291:GLU:O	1:B:295:LEU:HG	2.01	0.59
1:A:295:LEU:HD21	1:A:299:ARG:NH2	2.20	0.56
1:B:283:THR:C	1:B:285:ASP:H	2.10	0.55
3:D:16:DC:H3'	3:D:17:DA:C5'	2.37	0.54
1:A:73:LEU:HD22	1:A:86:ARG:CZ	2.38	0.54
1:B:257:TYR:HE1	1:B:262:LYS:O	1.91	0.53
1:A:356:ASP:O	1:A:357:HIS:HB3	2.09	0.52
1:A:229:ARG:O	1:A:233:LYS:HG3	2.10	0.52
1:A:162:SER:HB2	6:A:801:GOL:H2	1.92	0.52
1:A:108:ALA:O	1:A:314:SER:HA	2.11	0.51
1:A:341:SER:HB3	1:A:356:ASP:HA	1.93	0.50
1:B:129:TYR:HH	1:B:166:THR:HG1	1.58	0.49
1:A:337:TRP:CG	1:A:370:LYS:HE2	2.47	0.49
1:A:185:ASP:HA	1:A:209:SER:HB2	1.95	0.49
1:A:97:GLN:NE2	1:A:339:LYS:HG2	2.28	0.48
1:A:223:LYS:O	1:A:227:ILE:HG13	2.13	0.48
1:A:156:TYR:OH	1:A:173:LEU:HB2	2.14	0.48
1:B:163:THR:HG21	1:B:194:THR:HB	1.95	0.48
1:A:324:GLN:HG2	1:A:340:PRO:HA	1.97	0.47
9:A:398:ZZX:OAD	9:A:398:ZZX:HBC	2.15	0.46
1:A:161:PRO:O	1:A:189:ALA:HB2	2.16	0.46
1:A:69:ARG:HD2	1:A:70:GLU:OE2	2.16	0.45
1:A:127:ILE:HG22	1:A:145:VAL:HG13	1.99	0.45
1:A:238:ARG:HB3	1:A:239:PRO:HD2	1.97	0.45
1:A:337:TRP:CD2	1:A:370:LYS:HE2	2.51	0.45
1:A:319:VAL:HG13	1:A:344:LEU:O	2.17	0.45
1:A:295:LEU:HD21	1:A:299:ARG:CZ	2.47	0.44
1:B:185:ASP:C	1:B:185:ASP:OD2	2.56	0.44
1:A:358:LEU:N	1:A:359:GLY:CA	2.79	0.44
1:A:271:GLY:HA2	1:A:295:LEU:HD13	1.99	0.43
1:B:223:LYS:HA	1:B:223:LYS:HD3	1.85	0.42
1:B:222:ARG:NE	5:B:393:SO4:O3	2.48	0.42
1:A:356:ASP:O	1:A:357:HIS:CB	2.68	0.41
1:A:118:PRO:HG2	1:A:148:GLY:HA3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:GLN:HB3	1:A:21:ILE:HD11	2.02	0.41
1:A:111:PRO:O	1:A:350:ARG:HD3	2.20	0.41
1:B:127:ILE:HA	1:B:144:VAL:O	2.21	0.41
1:B:224:HIS:HA	1:B:227:ILE:HD12	2.02	0.40
1:A:221:GLU:OE1	1:A:221:GLU:HA	2.22	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)
10:A:422:HOH:O	10:D:501:HOH:O[8_554]	2.07	0.13

## 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%)	353 (96%)	12 (3%)	1 (0%)	46	73
1	B	182/395 (46%)	177 (97%)	5 (3%)	0	100	100
All	All	548/790 (69%)	530 (97%)	17 (3%)	1 (0%)	52	79

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	357	HIS

### 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%)	322 (98%)	5 (2%)	72	91
1	B	161/354 (46%)	156 (97%)	5 (3%)	47	76
All	All	488/708 (69%)	478 (98%)	10 (2%)	63	87

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
1	A	27	GLN
1	A	149	MET
1	A	357	HIS
1	A	366	ILE
1	B	149	MET
1	B	164	SER
1	B	216	SER
1	B	235	LEU
1	B	274	SER

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

Mol	Chain	Res	Type
1	B	224	HIS

### 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 carbohydrates in this entry.



## 5.6 Ligand geometry

Of 12 ligands modelled in this entry, 1 is modelled with single atom and 3 are monoatomic - 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
5	SO4	A	394	-	4,4,4	0.31	0	6,6,6	0.33	0
9	ZZX	A	398	8	28,35,35	1.27	3 (10%)	25,53,53	2.46	10 (40%)
5	SO4	A	399	-	4,4,4	0.69	0	6,6,6	0.50	0
6	GOL	A	801	-	5,5,5	0.35	0	5,5,5	0.52	0
6	GOL	A	802	-	5,5,5	0.38	0	5,5,5	0.45	0
6	GOL	A	803	-	5,5,5	0.38	0	5,5,5	0.19	0
6	GOL	A	805	-	5,5,5	0.41	0	5,5,5	0.28	0
5	SO4	B	393	-	4,4,4	0.15	0	6,6,6	0.23	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
5	SO4	A	394	-	-	0/0/0/0	0/0/0/0
9	ZZX	A	398	8	-	0/10/28/28	0/3/4/4
5	SO4	A	399	-	-	0/0/0/0	0/0/0/0
6	GOL	A	801	-	-	0/4/4/4	0/0/0/0
6	GOL	A	802	-	-	0/4/4/4	0/0/0/0
6	GOL	A	803	-	-	0/4/4/4	0/0/0/0
6	GOL	A	805	-	-	0/4/4/4	0/0/0/0
5	SO4	B	393	-	-	0/0/0/0	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	398	ZZX	CBA-CBB	-2.67	1.39	1.42
9	A	398	ZZX	CAX-CAR	2.23	1.54	1.50
9	A	398	ZZX	CAS-NBD	3.42	1.41	1.35

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	398	ZZX	CAC-CBC-NBF	-7.00	104.88	111.25
9	A	398	ZZX	CAV-CAL-CAU	-5.60	116.75	120.36
9	A	398	ZZX	FAH-CAT-CAV	-2.97	116.39	119.01
9	A	398	ZZX	CAN-CAU-CAL	-2.94	115.51	120.30
9	A	398	ZZX	OAE-CAS-NBD	-2.80	120.15	123.30
9	A	398	ZZX	CAB-NAQ-CAR	-2.60	118.70	121.80
9	A	398	ZZX	OAD-CAR-CAX	-2.56	117.55	121.01
9	A	398	ZZX	CBB-CAX-NAP	-2.22	119.10	122.33
9	A	398	ZZX	CAK-CAU-CAL	2.02	121.53	118.55
9	A	398	ZZX	CAL-CAV-CAT	3.32	122.67	119.71

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	398	ZZX	2	0
5	A	399	SO4	1	0
6	A	801	GOL	1	0
6	A	802	GOL	1	0
5	B	393	SO4	1	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	368/395 (93%)	0.15	13 (3%) 48 47	43, 58, 104, 132	0
1	B	184/395 (46%)	0.23	19 (10%) 9 6	50, 69, 120, 130	0
2	C	19/19 (100%)	-0.42	0 100 100	45, 60, 90, 115	0
3	D	17/17 (100%)	-0.43	0 100 100	50, 56, 86, 109	0
All	All	588/826 (71%)	0.14	32 (5%) 29 27	43, 61, 112, 132	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	8	LEU	8.5
1	B	280	ASN	4.6
1	A	9	ASP	4.5
1	A	19	HIS	4.4
1	A	29	THR	4.3
1	A	17	GLN	4.1
1	A	21	ILE	3.9
1	B	296	GLN	3.9
1	B	298	ILE	3.4
1	B	281	GLN	3.3
1	B	279	ALA	3.2
1	B	116	ASP	3.0
1	B	288	ARG	2.9
1	B	295	LEU	2.8
1	A	28	TYR	2.8
1	B	297	GLU	2.8
1	A	16	LEU	2.8
1	A	20	TYR	2.7
1	A	12	LEU	2.6
1	B	212	TYR	2.6
1	B	294	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	259	PRO	2.5
1	B	283	THR	2.4
1	A	254	ASN	2.4
1	B	237	GLY	2.4
1	B	282	ASP	2.3
1	B	292	LEU	2.3
1	B	289	GLU	2.3
1	B	299	ARG	2.2
1	B	257	TYR	2.1
1	A	153	THR	2.1
1	A	32	LEU	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
6	GOL	A	803	6/6	0.94	0.41	4.19	66,67,69,69	0
6	GOL	A	805	6/6	0.90	0.26	3.56	86,86,87,88	0
5	SO4	A	399	5/5	0.92	0.20	0.49	69,73,76,77	0
4	ZN	A	393	1/1	1.00	0.16	0.38	56,56,56,56	0
9	ZZX	A	398	32/32	0.98	0.15	-0.26	42,48,52,54	0
5	SO4	B	393	5/5	0.93	0.14	-0.70	105,105,106,106	0
8	MG	A	397	1/1	1.00	0.14	-0.75	33,33,33,33	0
6	GOL	A	802	6/6	0.81	0.20	-	68,74,75,76	0
6	GOL	A	801	6/6	0.83	0.20	-	88,90,90,90	0
8	MG	A	396	1/1	0.99	0.11	-	39,39,39,39	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
5	SO4	A	394	5/5	0.98	0.08	-	70,72,74,75	0
7	NH4	A	395	1/1	0.92	0.55	-	44,44,44,44	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.