



# Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 04:11 am BST

PDB ID : 3OYM  
Title : Crystal structure of the PFV N224H mutant intasome bound to manganese  
Authors : Hare, S.; Cherepanov, P.  
Deposited on : 2010-09-23  
Resolution : 2.02 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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

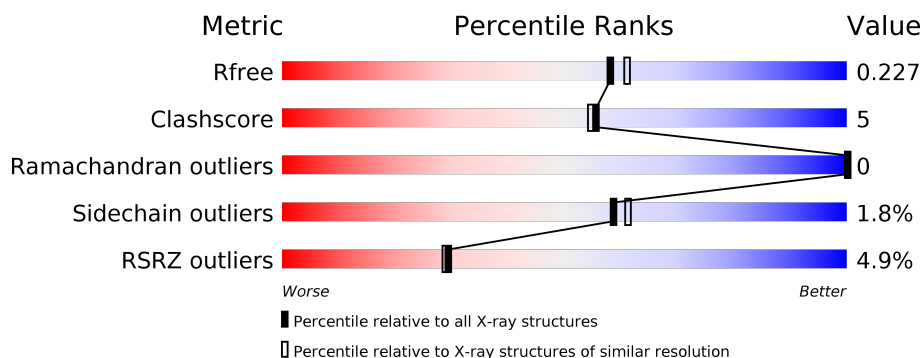
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.02 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	10434 (2.04-2.00)
Clashscore	141614	11643 (2.04-2.00)
Ramachandran outliers	138981	11493 (2.04-2.00)
Sidechain outliers	138945	11492 (2.04-2.00)
RSRZ outliers	127900	10220 (2.04-2.00)

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 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>4%</div> <div>83%</div> <div>10%</div> <div>7%</div> </div>
1	B	395	<div> <div>3%</div> <div>43%</div> <div>53%</div> </div>
2	C	19	<div> <div>5%</div> <div>58%</div> <div>21%</div> <div>21%</div> </div>
3	D	17	<div> <div>6%</div> <div>59%</div> <div>41%</div> </div>

## 2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 5821 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	5	0
			2966	1904	524	534	4			
1	B	184	Total	C	N	O	S	0	3	0
			1469	954	242	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			

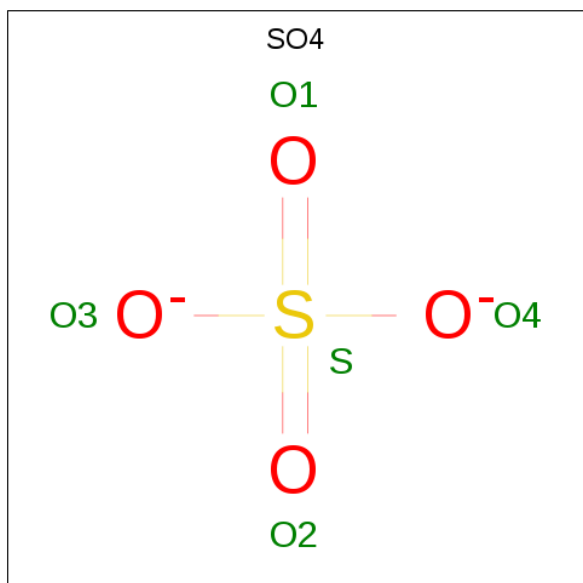
- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 5 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total	Mn	0	0
			2	2		
5	A	1	Total	Mn	0	0
			1	1		

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



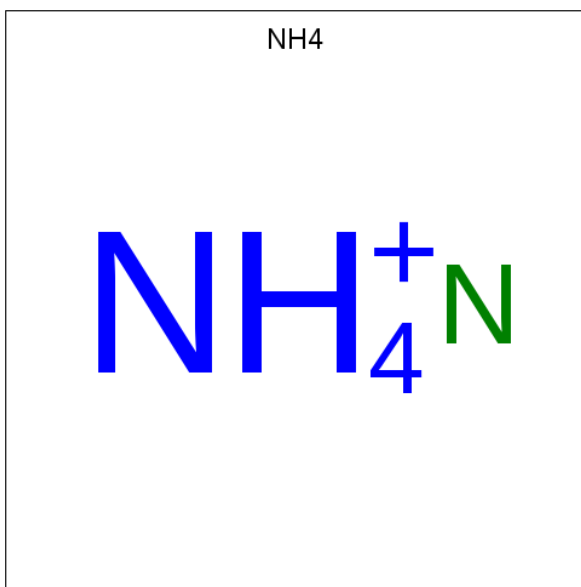
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



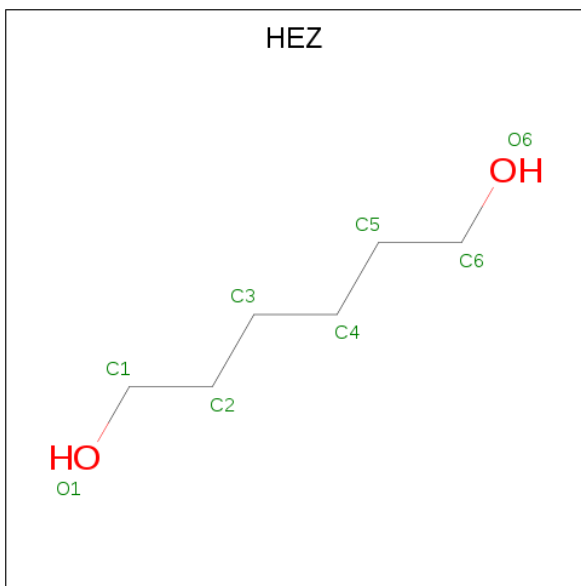
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	B	1	Total	C	O	0	0
			6	3	3		
7	B	1	Total	C	O	0	0
			6	3	3		
7	C	1	Total	C	O	0	0
			6	3	3		

- Molecule 8 is AMMONIUM ION (three-letter code: NH4) (formula:  $H_4N$ ).



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

- Molecule 9 is HEXANE-1,6-DIOL (three-letter code: HEZ) (formula:  $C_6H_{14}O_2$ ).



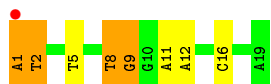
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	1	Total	C	O	0	0
			8	6	2		

- Molecule 10 is water.

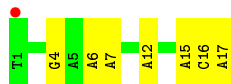
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	304	Total 304	O 304	0	0
10	B	146	Total 146	O 146	0	0
10	C	72	Total 72	O 72	0	0
10	D	74	Total 74	O 74	0	0







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



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	158.05Å 158.05Å 122.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.89 – 2.02 39.89 – 2.02	Depositor EDS
% Data completeness (in resolution range)	98.5 (39.89-2.02) 98.5 (39.89-2.02)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.78 (at 2.01Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.197 , 0.228 0.198 , 0.227	Depositor DCC
$R_{free}$ test set	5035 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.0	Xtriage
Anisotropy	0.385	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 47.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5821	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% 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, 0.333 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: GOL, ZN, MN, HEZ, 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	0.51	0/3061	0.62	2/4174 (0.0%)
1	B	0.54	0/1519	0.56	0/2075
2	C	0.95	0/433	1.70	10/667 (1.5%)
3	D	0.91	0/387	1.59	5/595 (0.8%)
All	All	0.60	0/5400	0.87	17/7511 (0.2%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	12	DA	O4'-C1'-N9	-10.84	100.42	108.00
1	A	69	ARG	NE-CZ-NH2	-9.53	115.54	120.30
1	A	69	ARG	NE-CZ-NH1	8.01	124.30	120.30
3	D	6	DA	O4'-C1'-N9	-7.02	103.08	108.00
3	D	7	DA	O4'-C1'-N9	-6.66	103.34	108.00
2	C	9	DG	P-O5'-C5'	6.23	130.87	120.90
3	D	17	DA	O4'-C1'-N9	6.15	112.30	108.00
2	C	8	DT	O4'-C1'-N1	-5.90	103.87	108.00
2	C	8	DT	O4'-C4'-C3'	-5.80	102.18	104.50
2	C	2	DT	N3-C2-O2	-5.80	118.82	122.30
2	C	1	DA	O4'-C1'-N9	5.77	112.04	108.00
2	C	16	DC	O4'-C1'-N1	-5.75	103.97	108.00
2	C	2	DT	N3-C4-O4	5.66	123.30	119.90
3	D	4	DG	P-O3'-C3'	5.59	126.41	119.70
2	C	5	DT	O4'-C1'-N1	-5.46	104.18	108.00
2	C	16	DC	P-O3'-C3'	5.35	126.12	119.70
2	C	12	DA	O4'-C1'-N9	-5.32	104.27	108.00

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	2966	0	3025	26	0
1	B	1469	0	1475	12	0
2	C	387	0	218	11	0
3	D	345	0	193	3	0
4	A	1	0	0	0	0
5	A	1	0	0	0	0
5	B	2	0	0	0	0
6	A	10	0	0	0	0
6	B	5	0	0	0	0
7	A	12	0	16	3	0
7	B	12	0	16	1	0
7	C	6	0	8	0	0
8	A	1	0	0	0	0
9	B	8	0	14	1	0
10	A	304	0	0	8	0
10	B	146	0	0	1	0
10	C	72	0	0	7	0
10	D	74	0	0	3	0
All	All	5821	0	4965	52	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 (52) 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)
2:C:1:DA:H4'	2:C:2:DT:H5'	1.24	1.12
1:B:117[B]:ARG:HH11	1:B:117[B]:ARG:HG2	1.13	1.06
1:B:117[B]:ARG:HH11	1:B:117[B]:ARG:CG	1.70	1.05
2:C:9:DG:P	10:C:92:HOH:O	2.27	0.90
1:A:97[A]:GLN:NE2	1:A:339:LYS:HE2	1.87	0.89
2:C:1:DA:C8	2:C:2:DT:H71	2.10	0.86
1:A:97[A]:GLN:HE22	1:A:339:LYS:HE2	1.41	0.86
2:C:1:DA:H2	10:C:319:HOH:O	1.62	0.83
1:A:303:TYR:HD2	10:A:689:HOH:O	1.61	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:15:DA:O3'	10:D:31:HOH:O	1.98	0.82
1:A:162:SER:HB2	7:A:801:GOL:H2	1.63	0.81
1:B:117[B]:ARG:NH1	1:B:117[B]:ARG:CG	2.41	0.75
1:A:136:SER:HA	7:A:398:GOL:H12	1.67	0.74
3:D:16:DC:H5'	10:D:31:HOH:O	1.88	0.74
1:A:120:LYS:HE3	10:A:497:HOH:O	1.88	0.72
2:C:8:DT:O3'	10:C:92:HOH:O	2.07	0.72
2:C:1:DA:H8	2:C:2:DT:C7	2.02	0.72
1:B:117[B]:ARG:NH1	1:B:117[B]:ARG:HG2	1.92	0.72
2:C:11:DA:OP2	10:C:517:HOH:O	2.11	0.69
1:A:304:HIS:ND1	10:A:522:HOH:O	2.27	0.67
2:C:1:DA:C2	10:C:319:HOH:O	2.42	0.67
2:C:1:DA:C8	2:C:2:DT:C7	2.76	0.67
1:A:174:THR:HB	1:A:178:ILE:HD13	1.78	0.65
1:A:341:SER:HB2	1:A:355:LEU:O	1.98	0.64
1:A:69:ARG:HD2	1:A:70:GLU:OE2	2.01	0.60
2:C:9:DG:C5'	10:C:92:HOH:O	2.50	0.59
1:B:228:LYS:HG2	9:B:397:HEZ:H12	1.85	0.59
2:C:9:DG:OP1	10:C:92:HOH:O	2.17	0.56
1:B:221:GLU:OE2	7:B:395:GOL:H31	2.06	0.55
1:A:137:GLN:H	7:A:398:GOL:H2	1.71	0.55
1:A:73:LEU:HD22	1:A:86:ARG:CZ	2.43	0.49
1:B:130:ILE:HD12	1:B:144:VAL:HG21	1.95	0.48
1:A:108:ALA:O	1:A:314:SER:HA	2.13	0.48
1:A:73:LEU:O	1:A:73:LEU:HG	2.13	0.48
1:B:117[B]:ARG:HH11	1:B:117[B]:ARG:HG3	1.70	0.47
1:B:119:GLN:HG3	10:B:482:HOH:O	2.14	0.46
1:A:223:LYS:NZ	10:A:501:HOH:O	2.47	0.45
1:B:235:LEU:HD21	1:B:242:TRP:HB3	1.99	0.45
1:A:348:ASN:HB2	1:A:349:PRO:CD	2.47	0.44
1:A:98:GLN:OE1	10:A:465:HOH:O	2.21	0.44
1:A:122:PHE:O	1:A:179:PRO:HA	2.18	0.44
1:B:141:TYR:CE2	1:B:161:PRO:HD3	2.52	0.43
3:D:16:DC:P	10:D:31:HOH:O	2.75	0.43
1:A:326:ARG:HD3	10:A:625:HOH:O	2.19	0.43
1:A:375:GLN:O	10:A:586:HOH:O	2.21	0.43
1:A:67:THR:HB	1:A:71:ALA:HB3	2.02	0.42
1:A:344:LEU:HD11	1:A:355:LEU:HB2	2.01	0.41
1:A:219:LYS:NZ	10:A:411:HOH:O	2.50	0.41
1:A:145:VAL:O	1:A:153:THR:HA	2.21	0.41
1:A:155:LEU:HD13	1:A:246:LEU:HD22	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:LEU:HD13	1:A:86:ARG:HG3	2.03	0.40
1:B:119:GLN:HG2	1:B:149:MET:HB2	2.03	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	371/395 (94%)	366 (99%)	5 (1%)	0	100	100
1	B	185/395 (47%)	184 (100%)	1 (0%)	0	100	100
All	All	556/790 (70%)	550 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 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	335/354 (95%)	328 (98%)	7 (2%)	53	55
1	B	165/354 (47%)	163 (99%)	2 (1%)	71	75
All	All	500/708 (71%)	491 (98%)	9 (2%)	59	61

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	36	LYS
1	A	51	GLN
1	A	149	MET
1	A	173	LEU
1	A	262	LYS
1	A	351	THR
1	B	116	ASP
1	B	274	SER

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

Mol	Chain	Res	Type
1	A	60	GLN
1	A	98	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 14 ligands modelled in this entry, 4 are monoatomic and 1 is modelled with single atom - 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	SO4	A	395	-	4,4,4	0.19	0	6,6,6	0.25	0
9	HEZ	B	397	-	7,7,7	0.39	0	6,6,6	0.38	0
7	GOL	A	398	-	5,5,5	0.47	0	5,5,5	0.99	0
6	SO4	A	397	-	4,4,4	0.14	0	6,6,6	0.27	0
7	GOL	B	395	-	5,5,5	0.35	0	5,5,5	0.67	0
7	GOL	C	20	-	5,5,5	0.52	0	5,5,5	1.20	1 (20%)
7	GOL	B	803	-	5,5,5	0.39	0	5,5,5	0.13	0
7	GOL	A	801	-	5,5,5	0.45	0	5,5,5	0.48	0
6	SO4	B	396	-	4,4,4	0.15	0	6,6,6	0.09	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	HEZ	B	397	-	-	2/5/5/5	-
7	GOL	A	398	-	-	0/4/4/4	-
7	GOL	B	395	-	-	2/4/4/4	-
7	GOL	C	20	-	-	3/4/4/4	-
7	GOL	B	803	-	-	2/4/4/4	-
7	GOL	A	801	-	-	2/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	20	GOL	C3-C2-C1	-2.52	101.91	111.70

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	B	395	GOL	O1-C1-C2-C3
7	C	20	GOL	C1-C2-C3-O3
7	B	803	GOL	O1-C1-C2-C3
7	A	801	GOL	C1-C2-C3-O3
7	B	395	GOL	O1-C1-C2-O2
9	B	397	HEZ	C2-C3-C4-C5
9	B	397	HEZ	C1-C2-C3-C4

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Mol	Chain	Res	Type	Atoms
7	C	20	GOL	O2-C2-C3-O3
7	B	803	GOL	O1-C1-C2-O2
7	A	801	GOL	O2-C2-C3-O3
7	C	20	GOL	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	B	397	HEZ	1	0
7	A	398	GOL	2	0
7	B	395	GOL	1	0
7	A	801	GOL	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%)	-0.04	17 (4%) 32 32	23, 33, 63, 83	0
1	B	184/395 (46%)	-0.04	10 (5%) 25 25	22, 34, 56, 83	0
2	C	19/19 (100%)	-0.49	1 (5%) 26 26	25, 36, 55, 74	0
3	D	17/17 (100%)	-0.63	1 (5%) 22 21	28, 31, 51, 68	0
All	All	588/826 (71%)	-0.07	29 (4%) 29 29	22, 33, 62, 83	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	358	LEU	6.9
1	A	8	LEU	6.2
1	B	278	PHE	6.1
1	A	9	ASP	5.5
1	A	359	GLY	5.4
1	B	116	ASP	4.7
1	B	298	ILE	4.6
1	B	295	LEU	4.5
1	B	279	ALA	4.1
1	A	19	HIS	4.1
1	A	357	HIS	4.0
1	A	17	GLN	3.7
1	B	299	ARG	3.6
1	B	296	GLN	3.4
1	A	10	ALA	2.9
1	A	16	LEU	2.8
1	A	29	THR	2.6
1	A	14	GLN	2.5
1	B	281	GLN	2.4
1	A	319	VAL	2.4
1	A	30	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	21	ILE	2.4
2	C	1	DA	2.3
1	A	13	ASP	2.3
1	A	35	GLY	2.2
1	A	32	LEU	2.1
1	B	280	ASN	2.1
1	B	237	GLY	2.1
3	D	1	DT	2.1

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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	B-factors(Å <sup>2</sup> )	Q<0.9
9	HEZ	B	397	8/8	0.74	0.20	58,58,61,62	0
7	GOL	A	398	6/6	0.78	0.18	55,56,56,57	0
7	GOL	B	803	6/6	0.79	0.23	62,64,65,68	0
7	GOL	A	801	6/6	0.82	0.14	58,59,60,60	0
7	GOL	B	395	6/6	0.86	0.16	49,51,51,51	0
6	SO4	B	396	5/5	0.92	0.16	98,98,98,98	0
7	GOL	C	20	6/6	0.95	0.13	30,32,37,42	0
6	SO4	A	397	5/5	0.96	0.11	64,64,65,65	0
8	NH4	A	396	1/1	0.96	0.39	31,31,31,31	0
6	SO4	A	395	5/5	0.98	0.08	57,57,59,59	0
5	MN	B	394	1/1	0.99	0.07	36,36,36,36	0
5	MN	A	394	1/1	1.00	0.04	36,36,36,36	0
4	ZN	A	393	1/1	1.00	0.09	29,29,29,29	0
5	MN	B	393	1/1	1.00	0.03	39,39,39,39	0

## 6.5 Other polymers

There are no such residues in this entry.