



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 04:27 am GMT

PDB ID : 3S3O
Title : Crystal structure of the Prototype Foamy Virus (PFV) N224H mutant in-tosome in complex with magnesium and Dolutegravir (S/GSK1349572)
Authors : Hare, S.; Cherepanov, P.
Deposited on : 2011-05-18
Resolution : 2.55 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

<http://wwpdb.org/validation/2016/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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

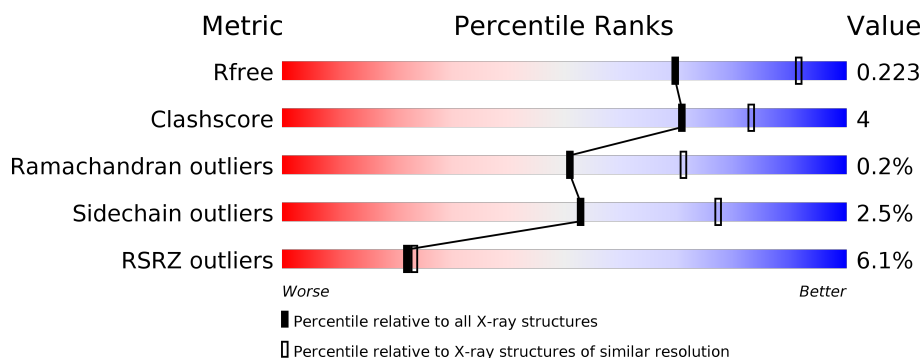
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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}	100719	4993 (2.58-2.50)
Clashscore	112137	5755 (2.58-2.50)
Ramachandran outliers	110173	5652 (2.58-2.50)
Sidechain outliers	110143	5654 (2.58-2.50)
RSRZ outliers	101464	5026 (2.58-2.50)

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. 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% 10% 7%</div> </div>
1	B	395	<div> <div>5%</div> <div>41% 5% 53%</div> </div>
2	C	19	<div> <div>32% 68%</div> </div>
3	D	17	<div> <div>59% 29% 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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	GOL	A	803	-	-	-	X

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 5396 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	1	0
			2934	1881	517	532	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	ENGINEERED MUTATION	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	VARIANT	UNP P14350

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	17	Total	C	N	O	P	0	1	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).



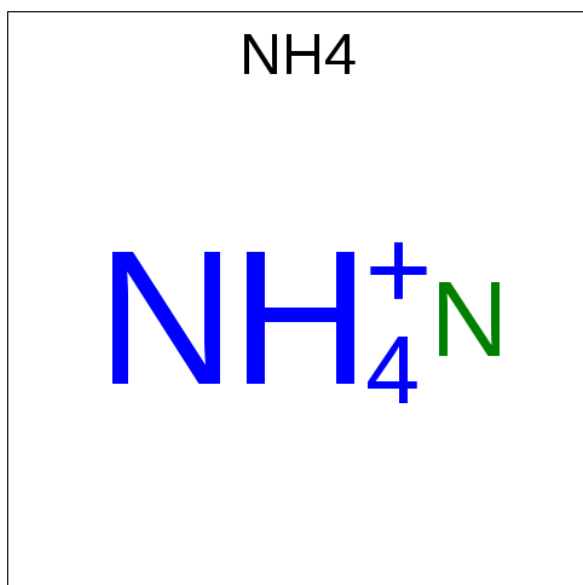
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		

- 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		

- Molecule 7 is AMMONIUM ION (three-letter code: NH4) (formula: H₄N).

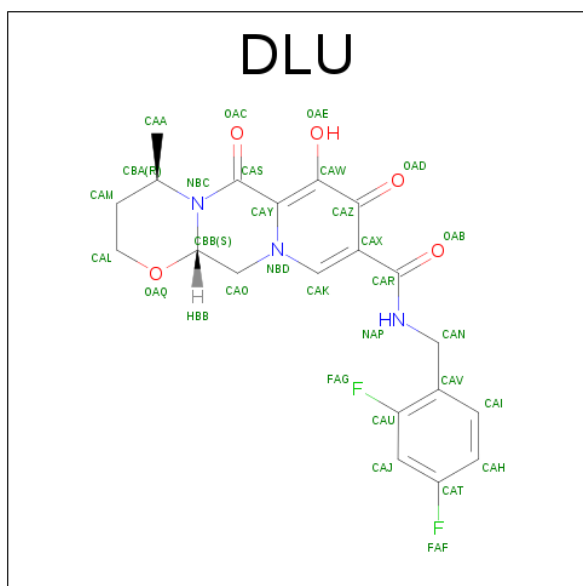


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total N 1 1	0	0

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

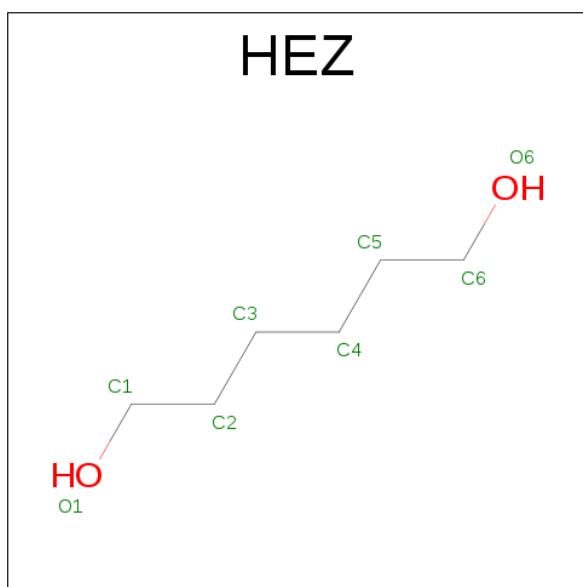
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	2	Total Mg 2 2	0	0

- Molecule 9 is (4R,12AS)-N-(2,4-DIFLUOROBENZYL)-7-HYDROXY-4-METHYL-6,8-DIOXO-3,4,6,8,12,12A-HEXAHYDRO-2H-PYRIDO[1',2':4,5]PYRAZINO[2,1-B][1,3]OXAZINE-9-CARBOXAMIDE (three-letter code: DLU) (formula: $C_{20}H_{19}F_2N_3O_5$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	A	1	Total 30	C 20	F 2	N 3	O 5	0	0

- Molecule 10 is HEXANE-1,6-DIOL (three-letter code: HEZ) (formula: $C_6H_{14}O_2$).



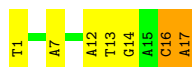
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	B	1	Total	C	O	0	0
			8	6	2		

- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	117	Total	O	0	0
			117	117		
11	B	36	Total	O	0	0
			36	36		
11	C	28	Total	O	0	0
			28	28		
11	D	18	Total	O	0	0
			18	18		

- Molecule 3: 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, α , β , γ	160.44Å 160.44Å 123.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.14 – 2.55 39.14 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.1 (39.14-2.55) 99.1 (39.14-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 2.54Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.208 , 0.230 0.201 , 0.223	Depositor DCC
R_{free} test set	2639 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	62.2	Xtriage
Anisotropy	0.271	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 39.7	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	0.95	EDS
Total number of atoms	5396	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.57% 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: GOL, ZN, MG, HEZ, NH4, SO4, DLU

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	0.56	0/3017	0.65	0/4117
1	B	0.56	0/1491	0.59	0/2039
2	C	1.03	0/433	1.82	11/667 (1.6%)
3	D	1.06	0/387	1.76	8/595 (1.3%)
All	All	0.66	0/5328	0.94	19/7418 (0.3%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	5	DT	O4'-C1'-N1	-11.12	100.22	108.00
2	C	7	DA	O4'-C1'-N9	-9.03	101.68	108.00
3	D	1	DT	P-O3'-C3'	7.60	128.82	119.70
2	C	1	DA	P-O3'-C3'	7.49	128.69	119.70
2	C	12	DA	O4'-C1'-N9	-7.27	102.91	108.00
3	D	13	DT	O4'-C1'-N1	-6.88	103.19	108.00
3	D	16	DC	O4'-C1'-N1	-6.64	103.35	108.00
3	D	12	DA	O4'-C1'-N9	-6.28	103.61	108.00
2	C	2	DT	P-O3'-C3'	5.94	126.83	119.70
2	C	13	DT	O4'-C1'-N1	-5.77	103.96	108.00
3	D	17[A]	DA	P-O5'-C5'	-5.73	111.73	120.90
2	C	2	DT	N3-C2-O2	-5.70	118.88	122.30
3	D	7	DA	O4'-C1'-N9	-5.63	104.06	108.00
3	D	14	DG	O4'-C1'-N9	-5.48	104.17	108.00
2	C	17	DG	O4'-C1'-N9	-5.46	104.18	108.00
3	D	17[A]	DA	O4'-C1'-N9	5.19	111.63	108.00
2	C	3	DT	N3-C4-O4	5.13	122.98	119.90
2	C	8	DT	C6-C5-C7	-5.11	119.83	122.90
2	C	18	DC	P-O3'-C3'	5.02	125.72	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	2934	0	2965	27	0
1	B	1450	0	1432	11	0
2	C	387	0	218	2	0
3	D	345	0	193	2	0
4	A	1	0	0	0	0
5	A	10	0	0	0	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	30	0	18	0	0
10	B	8	0	14	3	0
11	A	117	0	0	0	0
11	B	36	0	0	0	0
11	C	28	0	0	0	0
11	D	18	0	0	0	0
All	All	5396	0	4872	41	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 4.

All (41) 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	1.40	0.84
1:A:162:SER:HB2	6:A:801:GOL:H2	1.60	0.82
1:B:258:SER:O	1:B:261:LEU:O	2.08	0.71
1:A:69:ARG:HD2	1:A:70:GLU:OE2	1.93	0.67
1:A:358:LEU:N	1:A:359:GLY:HA2	2.10	0.66
1:A:356:ASP:O	1:A:357:HIS:HB3	1.96	0.66
1:A:73:LEU:HD13	1:A:86:ARG:HG3	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:LYS:HE2	1:A:95:ARG:HH21	1.66	0.60
1:A:219:LYS:HE3	2:C:6:DC:OP1	2.02	0.60
3:D:16:DC:H3'	3:D:17[A]:DA:H5'	1.83	0.60
1:B:250:GLN:HE22	10:B:394:HEZ:H12	1.70	0.56
1:A:341:SER:HB2	1:A:355:LEU:O	2.07	0.55
1:B:250:GLN:HE22	10:B:394:HEZ:C1	2.21	0.54
1:B:286:LEU:H	1:B:286:LEU:HD12	1.78	0.48
1:A:341:SER:CB	1:A:356:ASP:HA	2.44	0.48
1:A:315:TRP:CE2	1:A:371:PRO:HD3	2.48	0.48
1:A:174:THR:HB	1:A:178:ILE:HD13	1.95	0.47
1:A:76:ILE:HG22	1:A:85:MET:CE	2.44	0.47
1:B:130:ILE:HD12	1:B:144:VAL:HG21	1.96	0.47
1:A:337:TRP:CG	1:A:370:LYS:HE2	2.50	0.47
1:B:283:THR:C	1:B:285:ASP:H	2.18	0.47
3:D:16:DC:H3'	3:D:17[A]:DA:C5'	2.47	0.45
1:A:36:LYS:HE2	1:A:47:ILE:HD13	1.99	0.44
1:A:161:PRO:O	1:A:189:ALA:HB2	2.18	0.44
1:A:225:SER:O	1:A:229:ARG:HB2	2.17	0.44
1:A:156:TYR:HA	1:A:157:PRO:HD2	1.82	0.43
2:C:10:DG:H2''	2:C:11:DA:C8	2.54	0.43
1:B:184:SER:O	1:B:208:PHE:HA	2.19	0.43
1:A:91:LYS:HE2	1:A:95:ARG:NH2	2.34	0.42
1:A:43:GLU:H	1:A:43:GLU:CD	2.23	0.42
1:B:222:ARG:NE	5:B:393:SO4:O3	2.42	0.42
1:A:82:TRP:HB2	1:A:83:PRO:HD2	2.01	0.42
1:B:127:ILE:HA	1:B:144:VAL:O	2.20	0.41
1:A:12:LEU:HD23	1:A:37:VAL:HG21	2.01	0.41
1:A:141:TYR:CE2	1:A:161:PRO:HD3	2.56	0.41
1:A:356:ASP:O	1:A:357:HIS:CB	2.67	0.41
1:A:358:LEU:H	1:A:359:GLY:HA2	1.83	0.41
1:A:111:PRO:O	1:A:350:ARG:HD3	2.20	0.41
1:B:155:LEU:O	10:B:394:HEZ:H12	2.20	0.40
1:B:161:PRO:O	1:B:189:ALA:HB2	2.21	0.40
1:A:341:SER:HB3	1:A:356:ASP:HA	2.02	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	367/395 (93%)	355 (97%)	11 (3%)	1 (0%)	44	64
1	B	182/395 (46%)	176 (97%)	6 (3%)	0	100	100
All	All	549/790 (70%)	531 (97%)	17 (3%)	1 (0%)	51	71

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	329/354 (93%)	324 (98%)	5 (2%)	70	87
1	B	161/354 (46%)	154 (96%)	7 (4%)	33	56
All	All	490/708 (69%)	478 (98%)	12 (2%)	53	78

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ASP
1	A	149	MET
1	A	219	LYS
1	A	312	SER
1	A	357	HIS

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Mol	Chain	Res	Type
1	B	149	MET
1	B	164	SER
1	B	215	GLN
1	B	240	THR
1	B	274	SER
1	B	286	LEU
1	B	298	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	304	HIS
1	B	250	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 13 ligands modelled in this entry, 1 is modelled with single atom and 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
5	SO4	A	394	-	4,4,4	0.14	0	6,6,6	0.29	0
5	SO4	A	398	-	4,4,4	0.25	0	6,6,6	0.20	0
9	DLU	A	399	8	27,33,33	1.03	1 (3%)	33,49,49	1.93	10 (30%)
6	GOL	A	801	-	5,5,5	0.45	0	5,5,5	0.61	0
6	GOL	A	802	-	5,5,5	0.50	0	5,5,5	0.50	0
6	GOL	A	803	-	5,5,5	0.34	0	5,5,5	0.33	0
6	GOL	A	805	-	5,5,5	0.41	0	5,5,5	0.38	0
5	SO4	B	393	-	4,4,4	0.19	0	6,6,6	0.23	0
10	HEZ	B	394	-	7,7,7	0.50	0	6,6,6	0.86	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
5	SO4	A	398	-	-	0/0/0/0	0/0/0/0
9	DLU	A	399	8	-	0/9/35/35	1/3/4/4
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
10	HEZ	B	394	-	-	0/5/5/5	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	399	DLU	CAK-NBD	2.06	1.41	1.36

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	399	DLU	OAB-CAR-CAX	-3.73	116.31	121.66
9	A	399	DLU	CAO-NBD-CAK	-3.32	115.72	118.12
9	A	399	DLU	CAA-CBA-CAM	-3.32	106.10	112.70
9	A	399	DLU	CAK-NBD-CAY	-2.78	120.48	122.27
9	A	399	DLU	CAN-NAP-CAR	-2.12	116.43	121.82
9	A	399	DLU	CAN-CAV-CAI	-2.03	116.45	121.00
9	A	399	DLU	OAQ-CAL-CAM	2.54	115.82	110.71
9	A	399	DLU	CAX-CAR-NAP	2.70	120.16	116.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	399	DLU	CBA-NBC-CAS	3.00	123.55	118.25
9	A	399	DLU	CAO-NBD-CAY	4.22	123.79	118.68

There are no chirality outliers.

There are no torsion outliers.

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	A	399	DLU	CAL-CAM-CBA-CBB-NBC-OAQ

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	801	GOL	1	0
6	A	802	GOL	1	0
5	B	393	SO4	1	0
10	B	394	HEZ	3	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, 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%)	0.12	17 (4%) 33 36	37, 50, 95, 122	0
1	B	184/395 (46%)	0.27	19 (10%) 7 7	44, 61, 137, 146	0
2	C	19/19 (100%)	-0.32	0 100 100	38, 55, 81, 104	0
3	D	17/17 (100%)	-0.48	0 100 100	43, 48, 71, 103	0
All	All	588/826 (71%)	0.14	36 (6%) 22 23	37, 54, 115, 146	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	280	ASN	6.1
1	B	295	LEU	5.5
1	B	279	ALA	5.3
1	B	296	GLN	5.1
1	B	281	GLN	5.0
1	A	15	LEU	4.7
1	A	16	LEU	4.1
1	A	21	ILE	4.0
1	A	17	GLN	3.9
1	B	283	THR	3.9
1	A	8	LEU	3.7
1	A	14	GLN	3.7
1	B	298	ILE	3.7
1	B	116	ASP	3.5
1	A	19	HIS	3.4
1	A	20	TYR	3.2
1	B	237	GLY	3.1
1	B	292	LEU	3.1
1	A	29	THR	3.0
1	B	297	GLU	3.0
1	A	30	TYR	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	18	GLY	2.8
1	B	282	ASP	2.8
1	B	294	LEU	2.7
1	A	9	ASP	2.6
1	B	260	VAL	2.6
1	B	293	SER	2.6
1	A	359	GLY	2.6
1	B	289	GLU	2.6
1	A	358	LEU	2.4
1	B	288	ARG	2.4
1	B	215	GLN	2.3
1	A	31	PHE	2.3
1	A	32	LEU	2.2
1	B	212	TYR	2.1
1	A	265	PRO	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, 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	LLDF	B-factors(Å ²)	Q<0.9
6	GOL	A	803	6/6	0.81	0.26	2.32	74,75,75,78	0
6	GOL	A	805	6/6	0.91	0.18	1.46	86,88,88,90	0
10	HEZ	B	394	8/8	0.87	0.20	1.30	60,62,64,65	0
5	SO4	A	398	5/5	0.95	0.14	-0.26	84,86,86,86	0
5	SO4	B	393	5/5	0.95	0.15	-0.48	86,86,87,88	0
4	ZN	A	393	1/1	1.00	0.12	-0.88	46,46,46,46	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
9	DLU	A	399	30/30	0.96	0.12	-1.00	29,36,40,40	0
8	MG	A	397	1/1	0.98	0.08	-1.86	35,35,35,35	0
8	MG	A	396	1/1	0.98	0.06	-	35,35,35,35	0
5	SO4	A	394	5/5	0.97	0.12	-	74,76,78,78	0
6	GOL	A	802	6/6	0.69	0.27	-	66,72,76,76	0
7	NH4	A	395	1/1	0.94	0.48	-	32,32,32,32	0
6	GOL	A	801	6/6	0.70	0.23	-	78,81,81,81	0

6.5 Other polymers [i](#)

There are no such residues in this entry.