



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:23 PM UTC

PDB ID : 3VQ4 / pdb_00003vq4
Title : Fragments bound to HIV-1 integrase
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Deposited on : 2012-03-20
Resolution : 1.90 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
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

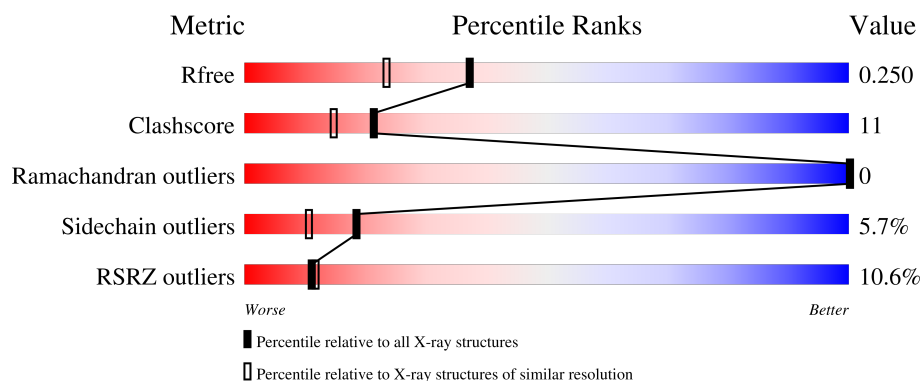
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	158	9% 71% 18% •• 9%
1	B	158	9% 66% 20% • 11%

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
3	SO4	A	303	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 2306 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 POL polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	144	Total	C	N	O	S	0	0	0
			1100	696	192	208	4			
1	B	140	Total	C	N	O	S	0	0	0
			1075	680	188	203	4			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	56	SER	CYS	engineered mutation	UNP Q72498
A	131	ASP	TRP	engineered mutation	UNP Q72498
A	139	ASP	PHE	engineered mutation	UNP Q72498
A	185	HIS	PHE	engineered mutation	UNP Q72498
B	56	SER	CYS	engineered mutation	UNP Q72498
B	131	ASP	TRP	engineered mutation	UNP Q72498
B	139	ASP	PHE	engineered mutation	UNP Q72498
B	185	HIS	PHE	engineered mutation	UNP Q72498

- Molecule 2 is CADMIUM ION (CCD ID: CD) (formula: Cd).

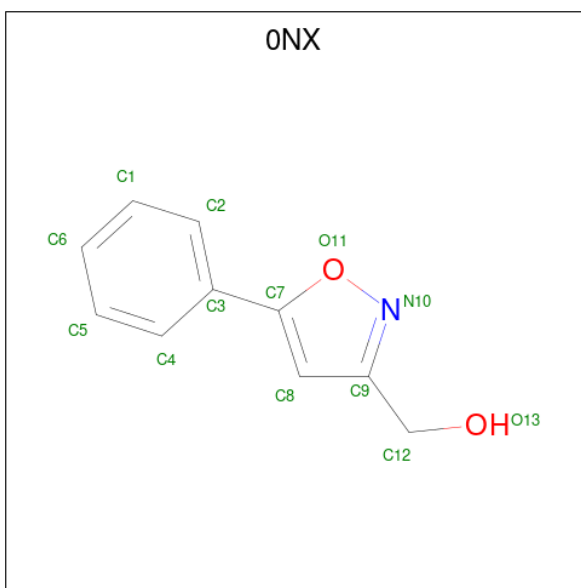
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Cd	0	0
			2	2		
2	B	2	Total	Cd	0	0
			2	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 4 is (5-phenyl-1,2-oxazol-3-yl)methanol (CCD ID: 0NX) (formula: C₁₀H₉NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			13	10	1	2		
4	A	1	Total	C	N	O	0	0
			13	10	1	2		
4	B	1	Total	C	N	O	0	0
			13	10	1	2		

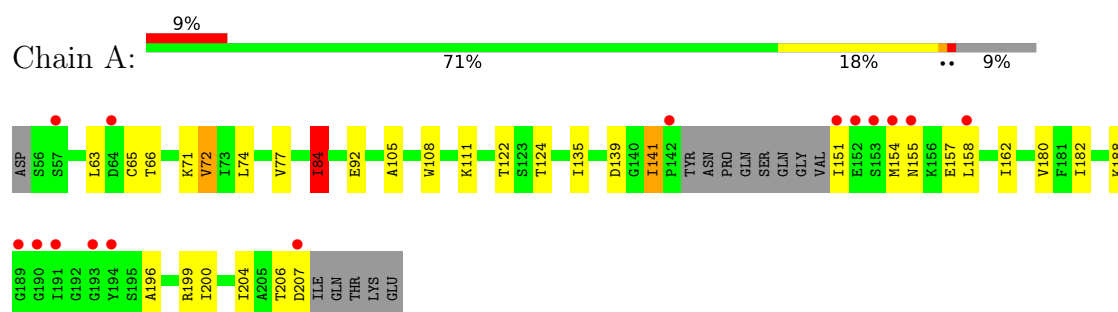
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	40	Total 40	O 40	0	0
5	B	43	Total 43	O 43	0	0

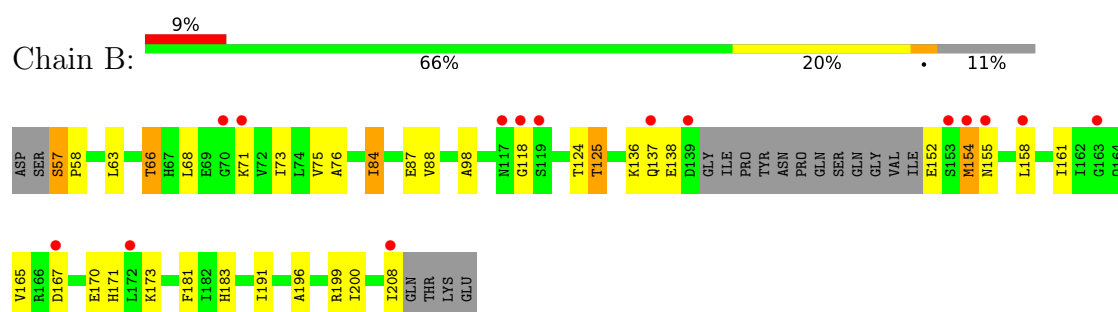
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: POL polypeptide



• Molecule 1: POL polypeptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	60.60Å 62.32Å 82.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.63 – 1.90 49.63 – 1.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.63-1.90) 100.0 (49.63-1.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.208 , 0.251 0.210 , 0.250	Depositor DCC
R_{free} test set	1267 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	28.9	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 43.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.024 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2306	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.44% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CD, SO4, ONX

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	1.38	4/1120 (0.4%)	1.19	2/1515 (0.1%)
1	B	1.35	4/1094 (0.4%)	1.24	8/1479 (0.5%)
All	All	1.37	8/2214 (0.4%)	1.22	10/2994 (0.3%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	124	THR	CB-OG1	-10.30	1.27	1.43
1	B	191	ILE	CA-CB	6.07	1.61	1.53
1	A	108	TRP	C-O	-5.95	1.18	1.24
1	A	72	VAL	CA-CB	5.81	1.61	1.54
1	A	141	ILE	CA-CB	5.73	1.61	1.54
1	B	191	ILE	CG1-CD1	5.48	1.73	1.51
1	A	182	ILE	CA-C	5.32	1.59	1.52
1	B	200	ILE	CA-CB	5.08	1.60	1.54

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	122	THR	N-CA-C	-6.51	105.33	113.28
1	B	84	ILE	N-CA-C	6.19	118.39	108.85
1	B	196	ALA	CA-C-N	5.95	126.72	119.99
1	B	196	ALA	C-N-CA	5.95	126.72	119.99
1	B	199	ARG	NE-CZ-NH2	5.91	124.52	119.20
1	A	84	ILE	N-CA-C	5.88	117.90	108.85
1	B	57	SER	CA-C-N	5.57	125.41	119.28
1	B	57	SER	C-N-CA	5.57	125.41	119.28
1	B	66	THR	CB-CA-C	5.15	119.54	109.35
1	B	199	ARG	NE-CZ-NH1	-5.05	116.44	121.50

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	1100	0	1101	21	0
1	B	1075	0	1075	26	0
2	A	2	0	0	0	1
2	B	2	0	0	0	0
3	A	5	0	0	3	0
4	A	26	0	18	1	0
4	B	13	0	9	1	0
5	A	40	0	0	3	0
5	B	43	0	0	2	1
All	All	2306	0	2203	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 11.

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:B:154:MET:HE2	1:B:183:HIS:HE1	1.27	0.97
3:A:303:SO4:O4	5:A:431:HOH:O	1.87	0.90
1:A:206:THR:HG23	4:A:304:ONX:H3	1.54	0.88
1:B:136:LYS:NZ	1:B:138:GLU:OE2	2.18	0.77
1:B:154:MET:HE2	1:B:183:HIS:CE1	2.18	0.72
1:B:208:ILE:HD11	4:B:303:ONX:N10	2.05	0.72
1:B:87:GLU:OE1	5:B:426:HOH:O	2.09	0.70
1:A:207:ASP:OD1	5:A:427:HOH:O	2.13	0.67
1:A:65:CYS:SG	5:A:434:HOH:O	2.54	0.66
1:A:66:THR:HG23	1:A:155:ASN:ND2	2.11	0.65
1:B:66:THR:HG23	1:B:73:ILE:HD12	1.79	0.64
1:A:139:ASP:O	1:A:141:ILE:HD12	1.98	0.63
1:A:84:ILE:HD11	1:A:180:VAL:HG22	1.80	0.62
1:B:170:GLU:HG2	1:B:171:HIS:CD2	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:ILE:HA	1:A:154:MET:HB2	1.85	0.59
1:B:57:SER:N	1:B:58:PRO:HD2	2.18	0.59
1:B:66:THR:HG22	1:B:73:ILE:HB	1.84	0.58
1:A:157:GLU:OE1	1:A:157:GLU:HA	2.04	0.58
1:A:77:VAL:HG22	1:A:84:ILE:HG22	1.86	0.57
1:B:118:GLY:HA3	5:B:430:HOH:O	2.05	0.56
1:B:84:ILE:CD1	1:B:158:LEU:HD13	2.36	0.55
1:A:72:VAL:HG11	1:A:92:GLU:HG3	1.88	0.54
1:B:154:MET:HG3	1:B:183:HIS:CE1	2.44	0.53
1:B:170:GLU:HG2	1:B:171:HIS:HD2	1.74	0.52
1:B:63:LEU:HD12	1:B:75:VAL:O	2.09	0.51
1:B:154:MET:CE	1:B:183:HIS:HE1	2.12	0.50
1:A:77:VAL:HG11	1:A:154:MET:HE3	1.93	0.50
1:A:188:LYS:NZ	1:A:199:ARG:HH12	2.12	0.48
1:B:63:LEU:HD13	1:B:76:ALA:HB2	1.96	0.48
1:B:170:GLU:HG2	1:B:171:HIS:N	2.29	0.47
1:A:158:LEU:O	1:A:162:ILE:HG13	2.14	0.47
1:B:98:ALA:CB	1:B:125:THR:HG22	2.45	0.47
1:B:170:GLU:CG	1:B:171:HIS:HD2	2.29	0.46
1:B:84:ILE:HD11	1:B:158:LEU:HD13	1.97	0.46
1:A:66:THR:HG23	1:A:155:ASN:CG	2.42	0.45
1:A:200:ILE:HG12	1:A:204:ILE:HD12	2.00	0.44
1:B:170:GLU:HG2	1:B:171:HIS:H	1.81	0.44
1:B:170:GLU:CG	1:B:171:HIS:CD2	2.99	0.43
1:A:188:LYS:HZ3	1:A:196:ALA:HB2	1.83	0.43
1:A:71:LYS:NZ	3:A:303:SO4:O4	2.52	0.42
1:A:188:LYS:HZ2	1:A:199:ARG:HH12	1.65	0.42
1:B:66:THR:CG2	1:B:73:ILE:HD12	2.48	0.42
1:A:105:ALA:HB2	1:A:135:ILE:HD11	2.02	0.42
1:B:88:VAL:HG21	1:B:173:LYS:HA	2.01	0.42
1:A:157:GLU:OE1	1:A:157:GLU:CA	2.66	0.41
1:B:181:PHE:CD2	1:B:181:PHE:C	2.99	0.41
1:A:71:LYS:NZ	3:A:303:SO4:S	2.94	0.40
1:B:161:ILE:O	1:B:165:VAL:HG22	2.21	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)
2:A:301:CD:CD	5:B:440:HOH:O[4_555]	1.22	0.98

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	140/158 (89%)	137 (98%)	3 (2%)	0	100	100
1	B	136/158 (86%)	131 (96%)	5 (4%)	0	100	100
All	All	276/316 (87%)	268 (97%)	8 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	116/129 (90%)	111 (96%)	5 (4%)	26	18
1	B	113/129 (88%)	105 (93%)	8 (7%)	13	7
All	All	229/258 (89%)	216 (94%)	13 (6%)	18	11

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	LEU
1	A	74	LEU
1	A	84	ILE
1	A	111	LYS
1	A	124	THR
1	B	68	LEU
1	B	71	LYS
1	B	125	THR

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Mol	Chain	Res	Type
1	B	137	GLN
1	B	152	GLU
1	B	154	MET
1	B	155	ASN
1	B	167	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	117	ASN
1	B	95	GLN
1	B	155	ASN
1	B	164	GLN
1	B	171	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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
4	ONX	B	303	-	14,14,14	0.91	1 (7%)	16,18,18	4.37	4 (25%)
3	SO4	A	303	-	4,4,4	0.35	0	6,6,6	0.75	0
4	ONX	A	304	-	14,14,14	1.88	3 (21%)	16,18,18	4.46	7 (43%)
4	ONX	A	305	-	14,14,14	0.77	1 (7%)	16,18,18	3.31	6 (37%)

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
4	ONX	B	303	-	-	0/6/6/6	0/2/2/2
4	ONX	A	304	-	-	0/6/6/6	0/2/2/2
4	ONX	A	305	-	-	0/6/6/6	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	304	ONX	C12-C9	4.56	1.53	1.50
4	A	304	ONX	C9-N10	2.49	1.37	1.29
4	A	304	ONX	C8-C7	2.47	1.39	1.35
4	A	305	ONX	C8-C7	2.11	1.38	1.35
4	B	303	ONX	C8-C7	2.07	1.38	1.35

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	304	ONX	O11-C7-C3	13.21	127.93	116.74
4	B	303	ONX	O11-C7-C3	11.92	126.84	116.74
4	A	305	ONX	O11-C7-C3	8.98	124.34	116.74
4	B	303	ONX	C7-O11-N10	8.42	112.96	108.73
4	B	303	ONX	O11-C7-C8	-8.07	104.22	109.23
4	A	304	ONX	O11-C7-C8	-7.34	104.68	109.23
4	A	305	ONX	O11-C7-C8	-6.30	105.32	109.23
4	A	304	ONX	C7-O11-N10	5.50	111.50	108.73
4	A	305	ONX	C7-O11-N10	5.34	111.42	108.73
4	A	304	ONX	O13-C12-C9	5.01	122.21	111.79
4	A	304	ONX	C8-C9-N10	-3.52	108.03	111.41
4	A	305	ONX	O13-C12-C9	-2.77	106.02	111.79
4	A	304	ONX	C3-C7-C8	-2.54	127.39	133.12
4	A	304	ONX	C4-C3-C2	2.43	121.66	118.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	305	0NX	C5-C4-C3	-2.24	118.16	120.36
4	B	303	0NX	C1-C2-C3	-2.09	118.31	120.36
4	A	305	0NX	C2-C3-C7	-2.07	117.48	120.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	303	0NX	1	0
3	A	303	SO4	3	0
4	A	304	0NX	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, 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	144/158 (91%)	0.69	15 (10%)	11 12	18, 29, 52, 75	0
1	B	140/158 (88%)	0.66	15 (10%)	11 11	18, 28, 58, 71	0
All	All	284/316 (89%)	0.68	30 (10%)	11 12	18, 29, 57, 75	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	151	ILE	5.6
1	B	139	ASP	4.6
1	B	117	ASN	4.4
1	A	189	GLY	4.1
1	A	207	ASP	3.9
1	A	194	TYR	3.8
1	A	142	PRO	3.7
1	A	152	GLU	3.7
1	B	154	MET	3.6
1	B	153	SER	3.3
1	A	153	SER	3.3
1	A	193	GLY	3.3
1	A	158	LEU	3.2
1	A	155	ASN	3.1
1	A	191	ILE	2.9
1	A	64	ASP	2.8
1	B	167	ASP	2.7
1	A	154	MET	2.6
1	A	57	SER	2.6
1	B	163	GLY	2.5
1	B	208	ILE	2.3
1	B	119	SER	2.3
1	B	118	GLY	2.2
1	B	172	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	190	GLY	2.2
1	B	71	LYS	2.2
1	B	155	ASN	2.2
1	B	70	GLY	2.2
1	B	158	LEU	2.1
1	B	137	GLN	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 oligosaccharides 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(Å ²)	Q<0.9
4	ONX	A	304	13/13	0.76	0.16	34,38,44,45	0
4	ONX	B	303	13/13	0.80	0.19	45,46,48,49	13
4	ONX	A	305	13/13	0.88	0.11	31,36,38,40	0
3	SO4	A	303	5/5	0.91	0.14	41,42,45,46	0
2	CD	B	302	1/1	0.95	0.06	52,52,52,52	0
2	CD	B	301	1/1	0.99	0.03	43,43,43,43	0
2	CD	A	301	1/1	0.99	0.02	24,24,24,24	1
2	CD	A	302	1/1	0.99	0.03	32,32,32,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.