



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 06:58 AM UTC

PDB ID : 3OS1 / pdb_00003os1
Title : PFV target capture complex (TCC) at 2.97 Å resolution
Authors : Maertens, G.N.; Hare, S.; Cherepanov, P.
Deposited on : 2010-09-08
Resolution : 2.97 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

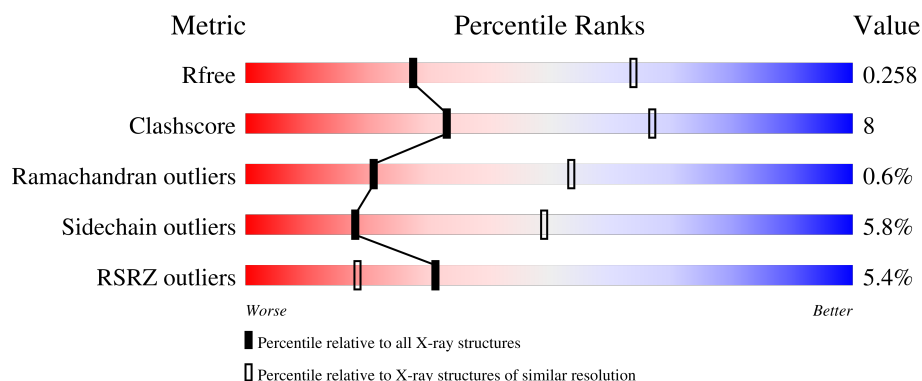
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.97 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	3580 (3.00-2.96)
Clashscore	190562	3904 (3.00-2.96)
Ramachandran outliers	187476	3761 (3.00-2.96)
Sidechain outliers	187428	3764 (3.00-2.96)
RSRZ outliers	180081	3579 (3.00-2.96)

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	3% 69% 21% 8%
1	B	395	4% 33% 8% 59%
2	C	19	5% 47% 47% 5%
3	D	17	82% 18%
4	T	30	7% 37% 13% 47%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 5229 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	365	Total	C	N	O	S	0	0	0
			2892	1855	509	524	4			
1	B	163	Total	C	N	O	S	0	0	0
			1270	833	204	232	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	SEE REMARK 999	UNP P14350
A	218	GLY	SER	SEE REMARK 999	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	SEE REMARK 999	UNP P14350
B	218	GLY	SER	SEE REMARK 999	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*(2DA))-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	17	Total	C	N	O	P	0	0	0
			344	166	65	97	16			

- Molecule 4 is a DNA chain called DNA (5'-D(*CP*CP*CP*GP*AP*GP*GP*CP*AP*CP*GP*TP*GP*CP*TP*AP*GP*CP*AP*CP*GP*TP*GP*CP*CP*TP*CP*GP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	T	16	Total	C	N	O	P	0	0	0
			328	155	61	96	16			

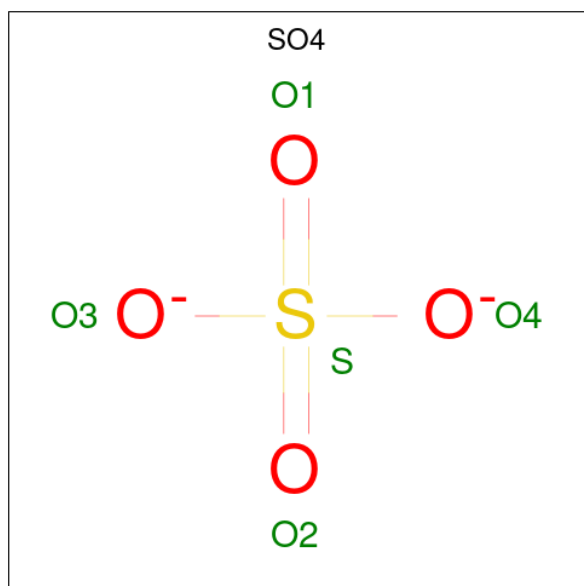
- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

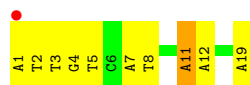
- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

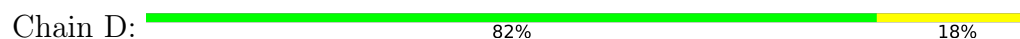
- Molecule 7 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



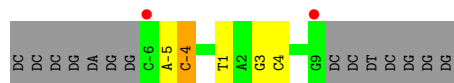
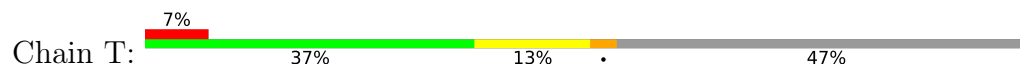
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	O	S	0	0
			5	4	1		



- Molecule 3: DNA (5'-D(*TP*GP*CP*GP*AP*AP*AP*TP*TP*CP*CP*AP*TP*GP*AP*CP*(2DA))-3')



- Molecule 4: DNA (5'-D(*CP*CP*CP*GP*AP*GP*GP*CP*AP*CP*GP*TP*GP*CP*TP*AP*GP*CP*AP*CP*GP*TP*GP*CP*CP*TP*CP*GP*GP*G)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	159.88Å 159.88Å 127.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.11 – 2.97 37.11 – 2.97	Depositor EDS
% Data completeness (in resolution range)	99.3 (37.11-2.97) 99.3 (37.11-2.97)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.95Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.231 , 0.264 0.227 , 0.258	Depositor DCC
R_{free} test set	1716 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	83.8	Xtriage
Anisotropy	0.436	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 44.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5229	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.95% 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: 2DA, MG, SO4, ZN

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	0.54	0/2971	0.85	1/4058 (0.0%)
1	B	0.54	0/1310	0.88	2/1796 (0.1%)
2	C	0.37	0/433	1.10	1/667 (0.1%)
3	D	0.33	0/363	1.02	0/558
4	T	0.35	0/367	0.96	1/564 (0.2%)
All	All	0.51	0/5444	0.90	5/7643 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	117	ARG	CA-C-N	5.67	126.92	119.84
1	B	117	ARG	C-N-CA	5.67	126.92	119.84
2	C	11	DA	P-O3'-C3'	5.51	128.47	120.20
4	T	-4	DC	P-O3'-C3'	5.32	128.18	120.20
1	A	256	THR	N-CA-C	5.31	117.18	108.52

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	238	ARG	Peptide

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	2892	0	2916	57	0
1	B	1270	0	1255	14	0
2	C	387	0	218	9	0
3	D	344	0	193	1	0
4	T	328	0	180	3	0
5	A	1	0	0	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	B	5	0	0	0	0
All	All	5229	0	4762	78	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 8.

All (78) 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:126:PHE:HB3	1:A:220:VAL:HG22	1.57	0.85
1:A:358:LEU:H	1:A:359:GLY:HA2	1.48	0.79
2:C:19:DA:H5''	2:C:19:DA:H8	1.48	0.77
1:A:358:LEU:N	1:A:359:GLY:HA2	2.01	0.75
1:B:258:SER:O	1:B:261:LEU:O	2.05	0.73
1:A:348:ASN:HB2	1:A:349:PRO:CD	2.23	0.68
1:A:356:ASP:HB3	1:A:360:ASN:H	1.62	0.64
1:A:73:LEU:HD13	1:A:86:ARG:HG3	1.80	0.63
2:C:19:DA:H5''	2:C:19:DA:C8	2.32	0.63
1:A:324:GLN:HG2	1:A:340:PRO:HA	1.82	0.62
1:A:126:PHE:HB3	1:A:220:VAL:CG2	2.30	0.62
2:C:7:DA:H1'	2:C:8:DT:H5'	1.83	0.60
2:C:1:DA:H2''	2:C:2:DT:H5'	1.81	0.60
1:A:275:ASN:HB3	1:B:178:ILE:HG22	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:T:3:DG:H2''	4:T:4:DC:H5'	1.84	0.60
1:A:356:ASP:O	1:A:359:GLY:HA2	2.03	0.57
1:A:14:GLN:HB3	1:A:19:HIS:HB2	1.86	0.57
1:A:109:SER:HB3	1:A:350:ARG:HH12	1.69	0.57
2:C:11:DA:H2''	2:C:12:DA:OP2	2.04	0.56
1:A:137:GLN:H	1:A:241:LYS:HE2	1.70	0.56
1:B:163:THR:HG21	1:B:194:THR:HB	1.87	0.56
1:A:12:LEU:HB3	1:A:21:ILE:HD13	1.87	0.56
1:A:315:TRP:CE2	1:A:371:PRO:HD3	2.41	0.55
1:A:209:SER:HB3	1:A:216:SER:HB2	1.88	0.54
1:A:139:TYR:CD1	1:A:157:PRO:HB2	2.42	0.54
1:A:129:TYR:CE2	1:A:143:LEU:HB2	2.42	0.53
1:A:126:PHE:CB	1:A:220:VAL:HG22	2.36	0.53
1:A:137:GLN:HE21	1:A:241:LYS:HD3	1.73	0.52
1:A:115:PRO:HG3	2:C:3:DT:C2	2.45	0.52
1:B:149:MET:C	1:B:149:MET:SD	2.94	0.51
2:C:1:DA:H3'	2:C:2:DT:H71	1.93	0.50
1:A:235:LEU:HD13	1:A:240:THR:HA	1.92	0.50
1:A:112:ILE:O	1:A:307:THR:HB	2.11	0.50
1:A:229:ARG:HG2	1:A:233:LYS:HE3	1.94	0.49
1:B:136:SER:HB2	1:B:242:TRP:NE1	2.28	0.49
1:A:348:ASN:HB2	1:A:349:PRO:HD2	1.94	0.48
1:A:118:PRO:HG2	1:A:148:GLY:HA3	1.93	0.48
1:A:127:ILE:HG22	1:A:145:VAL:HG22	1.94	0.48
1:A:222:ARG:NH1	1:A:226:ASP:OD2	2.47	0.48
1:A:82:TRP:HB2	1:A:83:PRO:HD2	1.96	0.48
1:A:274:SER:HB2	1:B:178:ILE:O	2.14	0.47
1:B:154:TRP:HB3	1:B:156:TYR:CE2	2.49	0.47
1:A:259:PRO:O	1:A:262:LYS:HD2	2.14	0.47
1:B:233:LYS:C	1:B:235:LEU:H	2.23	0.46
1:A:139:TYR:HD1	1:A:157:PRO:HB2	1.80	0.46
1:A:163:THR:HG21	1:A:194:THR:HB	1.96	0.46
1:A:66:HIS:HB3	1:A:99:CYS:SG	2.56	0.45
1:A:129:TYR:CZ	1:A:143:LEU:HD13	2.51	0.45
1:B:230:LEU:HG	1:B:249:VAL:HG13	1.98	0.45
1:A:358:LEU:N	1:A:359:GLY:CA	2.78	0.45
1:A:282:ASP:C	1:A:284:LEU:H	2.24	0.45
1:B:232:THR:O	1:B:235:LEU:HB3	2.17	0.45
1:A:149:MET:C	1:A:149:MET:SD	3.00	0.45
1:A:356:ASP:HB3	1:A:360:ASN:N	2.30	0.44
1:A:352:VAL:HG12	1:A:354:ILE:HG23	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:356:ASP:O	1:A:359:GLY:CA	2.66	0.43
1:A:163:THR:O	1:A:164:SER:C	2.61	0.43
1:B:220:VAL:HG12	1:B:224:ASN:HD22	1.84	0.43
4:T:-5:DA:H2''	4:T:-4:DC:OP2	2.18	0.43
1:A:174:THR:HA	1:A:177:ALA:O	2.19	0.43
1:A:221:GLU:O	1:A:224:ASN:HB2	2.19	0.43
1:A:185:ASP:C	1:A:187:GLY:H	2.27	0.42
1:B:161:PRO:O	1:B:189:ALA:HB2	2.19	0.42
1:A:132:PRO:HD2	4:T:1:DT:H5'	2.01	0.42
1:A:222:ARG:HD3	2:C:5:DT:O2	2.19	0.42
1:A:183:HIS:HA	1:A:207:GLU:O	2.20	0.42
1:A:350:ARG:HE	2:C:4:DG:P	2.43	0.41
1:B:167:VAL:O	1:B:168:LYS:C	2.63	0.41
1:A:54:ARG:O	1:A:55:GLN:C	2.64	0.41
1:A:230:LEU:HD23	1:A:249:VAL:HG13	2.02	0.41
1:A:348:ASN:HB2	1:A:349:PRO:HD3	2.03	0.41
1:A:356:ASP:HB2	1:A:362:ARG:HG3	2.02	0.41
3:D:11:DC:H2''	3:D:12:DA:N7	2.35	0.41
1:A:325:GLU:OE2	1:A:362:ARG:NH2	2.54	0.41
1:B:227:ILE:O	1:B:231:LEU:HB2	2.20	0.41
1:A:373:SER:C	1:A:374:HIS:CG	2.99	0.41
1:A:124:LYS:HG3	1:A:181:VAL:HB	2.03	0.40
1:A:269:LEU:HD23	1:A:270:PHE:CE1	2.57	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	363/395 (92%)	332 (92%)	30 (8%)	1 (0%)	36	67
1	B	161/395 (41%)	145 (90%)	14 (9%)	2 (1%)	10	38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	524/790 (66%)	477 (91%)	44 (8%)	3 (1%)	21 53

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	239	PRO
1	A	357	HIS
1	B	135	PRO

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	323/354 (91%)	303 (94%)	20 (6%)	16 46
1	B	140/354 (40%)	133 (95%)	7 (5%)	22 54
All	All	463/708 (65%)	436 (94%)	27 (6%)	18 49

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	LEU
1	A	102	THR
1	A	178	ILE
1	A	184	SER
1	A	194	THR
1	A	209	SER
1	A	210	THR
1	A	217	SER
1	A	220	VAL
1	A	226	ASP
1	A	256	THR
1	A	262	LYS
1	A	272	ILE
1	A	286	LEU
1	A	307	THR

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Mol	Chain	Res	Type
1	A	312	SER
1	A	316	SER
1	A	342	THR
1	A	351	THR
1	A	374	HIS
1	B	176	ILE
1	B	210	THR
1	B	216	SER
1	B	231	LEU
1	B	235	LEU
1	B	256	THR
1	B	272	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	137	GLN
1	A	255	ASN
1	A	338	HIS
1	A	357	HIS
1	A	374	HIS
1	B	183	HIS
1	B	205	HIS
1	B	224	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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
3	2DA	D	17	3	19,22,23	1.56	4 (21%)	25,31,34	3.16	12 (48%)

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
3	2DA	D	17	3	-	3/7/18/19	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	17	2DA	C1'-N9	2.80	1.51	1.46
3	D	17	2DA	O4'-C1'	2.76	1.48	1.42
3	D	17	2DA	O4'-C4'	2.73	1.50	1.44
3	D	17	2DA	C5-N7	-2.70	1.34	1.39

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	17	2DA	C4'-O4'-C1'	-8.69	101.61	109.81
3	D	17	2DA	O4'-C1'-N9	6.87	120.31	107.96
3	D	17	2DA	C5-C4-N3	-4.89	119.99	126.72
3	D	17	2DA	N3-C2-N1	-4.57	121.66	128.58
3	D	17	2DA	N3-C4-N9	3.65	133.38	127.17
3	D	17	2DA	O4'-C4'-C5'	3.49	115.64	109.34
3	D	17	2DA	C2-N3-C4	3.33	119.97	111.83
3	D	17	2DA	C3'-C2'-C1'	-2.85	99.60	102.87
3	D	17	2DA	N9-C8-N7	-2.66	110.16	113.94
3	D	17	2DA	O4'-C1'-C2'	-2.44	103.53	106.41
3	D	17	2DA	C5-N7-C8	2.20	106.91	103.45
3	D	17	2DA	C4-N9-C8	2.02	107.86	105.74

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	17	2DA	O4'-C1'-N9-C8
3	D	17	2DA	O4'-C1'-N9-C4
3	D	17	2DA	O4'-C4'-C5'-O5'

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 3 are monoatomic - leaving 1 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	$\# Z > 2$
7	SO4	B	394	-	4,4,4	0.23	0	6,6,6	0.17	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	365/395 (92%)	0.26	13 (3%) 46 29	59, 75, 117, 131	0
1	B	163/395 (41%)	0.64	15 (9%) 14 9	65, 85, 108, 115	0
2	C	19/19 (100%)	-0.07	1 (5%) 32 19	65, 78, 98, 112	0
3	D	16/17 (94%)	-0.18	0 100 100	69, 79, 102, 122	0
4	T	16/30 (53%)	0.68	2 (12%) 8 6	72, 85, 135, 144	0
All	All	579/856 (67%)	0.36	31 (5%) 31 18	59, 77, 115, 144	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	234	LEU	5.9
1	A	12	LEU	5.7
4	T	9	DG	4.0
1	B	233	LYS	3.7
1	A	10	ALA	3.5
1	B	236	VAL	3.4
1	B	257	TYR	3.2
1	B	212	TYR	3.2
1	A	22	LYS	3.0
1	B	239	PRO	3.0
1	A	28	TYR	2.9
1	A	297	GLU	2.8
1	B	237	GLY	2.7
1	B	278	PHE	2.6
1	B	213	HIS	2.5
1	B	256	THR	2.5
1	B	261	LEU	2.5
1	B	273	ASP	2.4
1	A	11	GLU	2.4
1	B	240	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	262	LYS	2.2
4	T	-6	DC	2.2
1	A	285	ASP	2.2
1	A	137	GLN	2.1
1	A	357	HIS	2.1
1	A	241	LYS	2.1
1	A	20	TYR	2.1
2	C	1	DA	2.1
1	A	138	GLY	2.0
1	A	16	LEU	2.0
1	B	260	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
3	2DA	D	17	20/21	0.94	0.17	69,80,81,81	0

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
7	SO4	B	394	5/5	0.89	0.16	138,138,138,138	0
6	MG	B	393	1/1	0.93	0.20	85,85,85,85	0
6	MG	A	394	1/1	0.97	0.10	64,64,64,64	0
5	ZN	A	393	1/1	1.00	0.03	70,70,70,70	0

6.5 Other polymers ⓘ

There are no such residues in this entry.