



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:08 AM UTC

PDB ID : 5WCP / pdb_00005wcp
Title : Phosphotriesterase variant S7
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Deposited on : 2017-07-01
Resolution : 1.50 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

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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

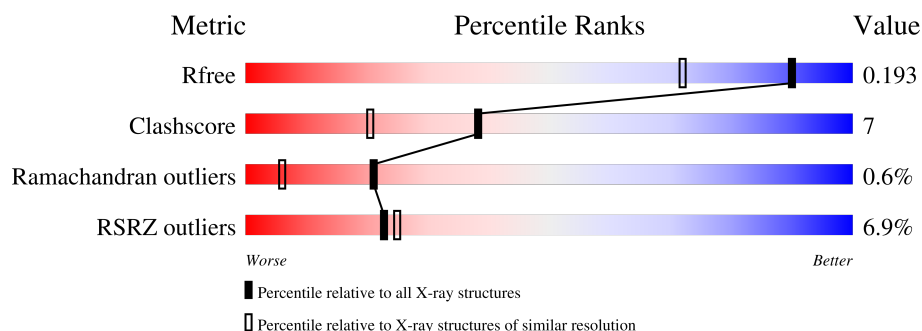
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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 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	333	6% 81% 14% 5%
1	G	333	7% 80% 13% 6%

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	CAC	G	2403	-	X	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5653 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 Phosphotriesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	316	Total	C	N	O	S	0	24	0
			2609	1633	483	486	7			
1	G	312	Total	C	N	O	S	0	25	0
			2589	1625	471	486	7			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	216	LEU	PHE	conflict	UNP A0A060GZX0
A	233	ALA	ASP	conflict	UNP A0A060GZX0
A	234	ALA	THR	conflict	UNP A0A060GZX0
A	254	SER	ARG	conflict	UNP A0A060GZX0
A	263	GLY	GLU	conflict	UNP A0A060GZX0
A	271	HIS	LEU	conflict	UNP A0A060GZX0
A	293	THR	MET	conflict	UNP A0A060GZX0
A	306	ILE	PHE	conflict	UNP A0A060GZX0
A	320	GLY	VAL	conflict	UNP A0A060GZX0
G	216	LEU	PHE	conflict	UNP A0A060GZX0
G	233	ALA	ASP	conflict	UNP A0A060GZX0
G	234	ALA	THR	conflict	UNP A0A060GZX0
G	254	SER	ARG	conflict	UNP A0A060GZX0
G	263	GLY	GLU	conflict	UNP A0A060GZX0
G	271	HIS	LEU	conflict	UNP A0A060GZX0
G	293	THR	MET	conflict	UNP A0A060GZX0
G	306	ILE	PHE	conflict	UNP A0A060GZX0
G	320	GLY	VAL	conflict	UNP A0A060GZX0

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

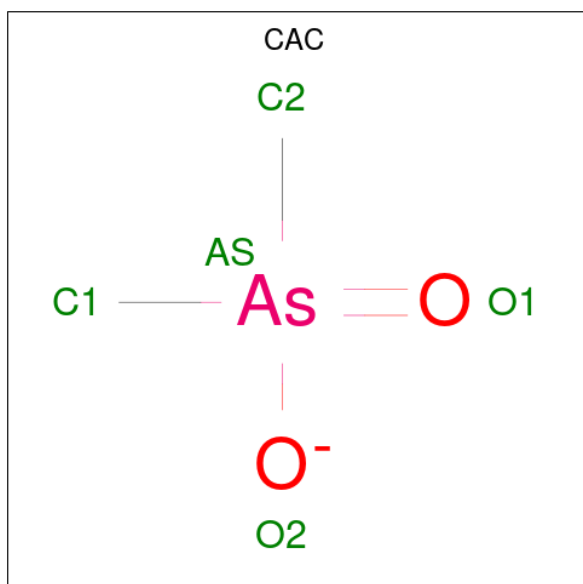
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Zn	0	0
			2	2		

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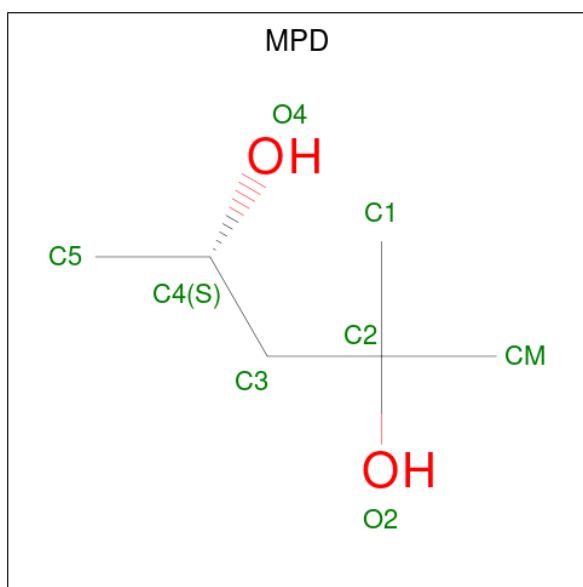
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	2	Total	Zn	0	0
			2	2		

- Molecule 3 is CACODYLATE ION (CCD ID: CAC') (formula: $\text{C}_2\text{H}_6\text{AsO}_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	As	C	O	0	0
			5	1	2	2		
3	G	1	Total	As	C	O	0	0
			5	1	2	2		

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: $\text{C}_6\text{H}_{14}\text{O}_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 8 6 2	0	0
4	A	1	Total C O 8 6 2	0	0
4	A	1	Total C O 8 6 2	0	0
4	A	1	Total C O 8 6 2	0	0
4	G	1	Total C O 8 6 2	0	0
4	G	1	Total C O 8 6 2	0	0

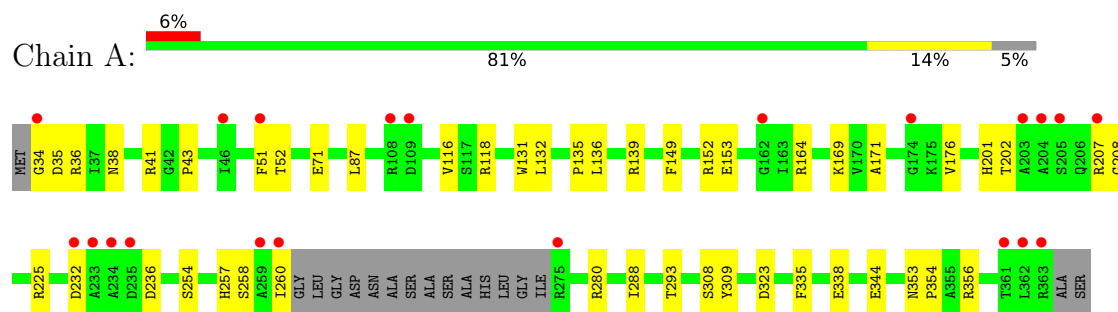
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	212	Total O 212 212	0	0
5	G	181	Total O 181 181	0	0

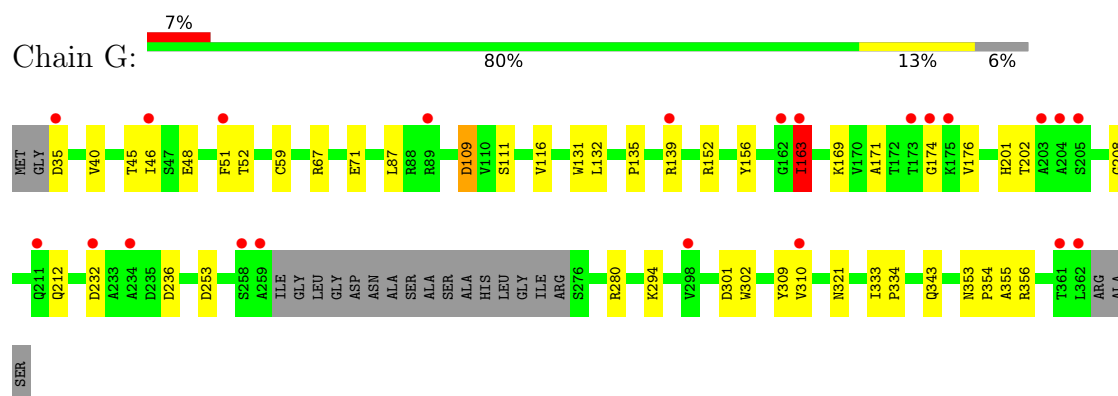
3 Residue-property plots

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: Phosphotriesterase



• Molecule 1: Phosphotriesterase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	85.60Å 85.95Å 88.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.48 – 1.50 29.48 – 1.50	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.48-1.50) 99.8 (29.48-1.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.45 (at 1.50Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.151 , 0.175 (Not available) , 0.193	Depositor DCC
R_{free} test set	5319 reflections (3.33%)	wwPDB-VP
Wilson B-factor (Å ²)	18.6	Xtriage
Anisotropy	0.661	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.008 for -h,l,k 0.009 for -l,-k,-h 0.010 for k,h,-l 0.000 for k,l,h 0.000 for l,h,k	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5653	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.71% 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: KCX, ZN, MPD, CAC

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	0.82	0/2639	0.85	2/3577 (0.1%)
1	G	0.74	0/2619	0.83	2/3555 (0.1%)
All	All	0.78	0/5258	0.84	4/7132 (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	G	0	4

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	G	163	ILE	O-C-N	-10.49	111.24	122.68
1	G	109	ASP	O-C-N	-7.26	113.38	122.82
1	A	34	GLY	CA-C-N	5.31	131.68	121.54
1	A	34	GLY	C-N-CA	5.31	131.68	121.54

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	G	109	ASP	Mainchain
1	G	163	ILE	Mainchain

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	2609	0	2632	37	0
1	G	2589	0	2605	36	0
2	A	2	0	0	0	0
2	G	2	0	0	0	0
3	A	5	0	0	0	0
3	G	5	0	0	0	0
4	A	32	0	56	7	0
4	G	16	0	28	1	0
5	A	212	0	0	7	0
5	G	181	0	0	3	0
All	All	5653	0	5321	73	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 7.

All (73) 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:323:ASP:OD1	5:A:2501:HOH:O	2.08	0.70
1:G:236:ASP:OD2	5:G:2501:HOH:O	2.10	0.70
1:G:202[B]:THR:HG23	1:G:212:GLN:NE2	2.10	0.66
1:A:71[B]:GLU:HG2	4:A:2407:MPD:H12	1.79	0.64
1:A:135:PRO:O	1:A:139[A]:ARG:HG3	1.99	0.62
1:G:71[B]:GLU:H	1:G:71[B]:GLU:CD	2.06	0.62
1:G:87:LEU:HD12	1:G:116[B]:VAL:HG12	1.81	0.61
1:A:51:PHE:HD1	1:A:52:THR:N	2.00	0.60
1:A:257:HIS:HA	1:A:260:ILE:HD12	1.83	0.60
1:A:309:TYR:HB2	4:A:2406:MPD:C1	2.33	0.59
1:A:87:LEU:HD12	1:A:116[B]:VAL:HG12	1.83	0.58
1:G:309:TYR:CD2	1:G:310:VAL:HG22	2.38	0.58
1:A:38[A]:ASN:OD1	5:A:2502:HOH:O	2.16	0.58
1:G:294:LYS:O	1:G:356[C]:ARG:NH2	2.36	0.58
1:G:353:ASN:HB2	1:G:354:PRO:HD3	1.85	0.57
1:A:308[A]:SER:O	4:A:2406:MPD:H13	2.05	0.56
1:A:149:PHE:O	1:A:152:ARG:HG3	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:ILE:HA	1:A:293[B]:THR:HG22	1.87	0.55
1:A:344:GLU:CD	1:A:344:GLU:H	2.15	0.55
1:G:202[B]:THR:HG22	5:G:2542:HOH:O	2.06	0.55
1:A:71[B]:GLU:CD	1:G:152:ARG:HE	2.15	0.54
1:A:51:PHE:C	1:A:51:PHE:CD1	2.87	0.52
1:G:51:PHE:HD1	1:G:52:THR:N	2.07	0.52
1:G:131:TRP:CG	1:G:132:LEU:H	2.27	0.52
1:A:71[A]:GLU:HG2	4:A:2407:MPD:H12	1.91	0.52
1:G:87:LEU:HD12	1:G:116[B]:VAL:CG1	2.39	0.52
1:A:171:ALA:HA	1:A:201:HIS:HB3	1.92	0.51
1:A:207[A]:ARG:NH1	1:A:236:ASP:OD2	2.43	0.51
1:A:51:PHE:HD1	1:A:51:PHE:C	2.20	0.50
1:G:135:PRO:O	1:G:139:ARG:HG3	2.10	0.50
1:A:131:TRP:CG	1:A:132:LEU:H	2.31	0.49
1:G:67[A]:ARG:NH2	1:G:111[A]:SER:OG	2.46	0.49
1:G:232:ASP:OD2	1:G:280:ARG:NH1	2.46	0.48
1:G:35[A]:ASP:OD1	1:G:35[A]:ASP:N	2.47	0.48
1:A:225:ARG:CD	5:A:2514:HOH:O	2.61	0.47
1:A:225:ARG:HD3	5:A:2514:HOH:O	2.15	0.47
1:A:356[B]:ARG:HH11	1:A:356[B]:ARG:HG3	1.79	0.47
1:G:333:ILE:HB	1:G:334:PRO:HD3	1.96	0.47
1:A:335:PHE:O	1:A:338:GLU:HG2	2.14	0.47
1:A:87:LEU:HD12	1:A:116[B]:VAL:CG1	2.44	0.47
1:A:232:ASP:OD2	1:A:254[B]:SER:OG	2.26	0.46
1:A:136:LEU:HG	1:A:139[A]:ARG:NH2	2.30	0.46
1:A:152:ARG:HG3	1:A:153:GLU:N	2.30	0.46
1:A:41[B]:ARG:HB2	5:A:2574:HOH:O	2.14	0.46
4:A:2404:MPD:H52	4:A:2404:MPD:O2	2.16	0.45
1:A:41[A]:ARG:NH2	1:A:118:ARG:HD3	2.31	0.45
1:G:51:PHE:C	1:G:51:PHE:CD1	2.95	0.45
1:G:46:ILE:HA	1:G:46:ILE:HD12	1.78	0.45
1:G:171:ALA:HA	1:G:201:HIS:HB3	1.99	0.44
1:G:46:ILE:HD11	1:G:355:ALA:HB1	2.00	0.44
1:A:71[B]:GLU:OE2	1:G:152:ARG:HG3	2.16	0.44
1:G:302:TRP:CH2	1:G:321:ASN:HB3	2.53	0.44
1:A:353:ASN:HB2	1:A:354:PRO:HD3	1.98	0.44
1:G:202[B]:THR:HG22	1:G:208:GLY:HA3	1.98	0.44
4:A:2406:MPD:H13	5:A:2531:HOH:O	2.17	0.44
1:G:51:PHE:HD1	1:G:51:PHE:C	2.25	0.43
1:A:202[A]:THR:OG1	1:A:208:GLY:HA3	2.18	0.43
1:A:356[B]:ARG:HG3	1:A:356[B]:ARG:NH1	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:202[B]:THR:HG21	1:G:208:GLY:C	2.43	0.43
1:G:59[A]:CYS:O	1:G:59[A]:CYS:SG	2.77	0.43
4:A:2407:MPD:H31	1:G:156:TYR:CE1	2.53	0.42
1:G:353:ASN:OD1	1:G:356[A]:ARG:NH2	2.51	0.42
1:G:171:ALA:CB	1:G:201:HIS:HB3	2.50	0.42
1:G:253:ASP:O	1:G:301:ASP:HB2	2.19	0.42
1:A:258:SER:OG	1:A:280:ARG:NE	2.52	0.41
1:G:40:VAL:HA	1:G:163:ILE:HG23	2.00	0.41
1:G:59[A]:CYS:O	5:G:2502:HOH:O	2.21	0.41
1:A:232:ASP:OD2	1:A:254[A]:SER:HB2	2.20	0.41
1:G:343:GLN:NE2	4:G:2405:MPD:H4	2.35	0.41
1:G:309:TYR:CE2	1:G:310:VAL:HG22	2.56	0.41
1:A:36[B]:ARG:HD3	1:A:43:PRO:HB2	2.03	0.41
1:G:45:THR:OG1	1:G:48[A]:GLU:HG3	2.20	0.41
1:A:164:ARG:NH1	5:A:2512:HOH:O	2.53	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	335/333 (101%)	323 (96%)	10 (3%)	2 (1%)	21	6
1	G	332/333 (100%)	320 (96%)	10 (3%)	2 (1%)	21	6
All	All	667/666 (100%)	643 (96%)	20 (3%)	4 (1%)	21	6

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	174	GLY
1	A	35	ASP
1	A	176	VAL

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Mol	Chain	Res	Type
1	G	176	VAL

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are 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$
1	KCX	A	169	1,2	10,11,12	2.44	3 (30%)	6,12,14	0.79	0
1	KCX	G	169	1,2	10,11,12	1.55	2 (20%)	6,12,14	2.04	1 (16%)

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
1	KCX	A	169	1,2	-	0/9/10/12	-
1	KCX	G	169	1,2	-	0/9/10/12	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	169	KCX	CE-NZ	5.04	1.57	1.46
1	A	169	KCX	CX-NZ	-4.54	1.27	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	169	KCX	CE-NZ	3.54	1.54	1.46
1	A	169	KCX	OQ1-CX	3.12	1.27	1.21
1	G	169	KCX	CB-CA	2.11	1.56	1.53

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	169	KCX	OQ1-CX-NZ	-4.72	117.74	124.92

There are no chirality outliers.

There are no torsion outliers.

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 12 ligands modelled in this entry, 4 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	$\# Z > 2$
4	MPD	A	2404	-	7,7,7	0.55	0	9,10,10	0.99	0
3	CAC	A	2403	2	2,4,4	1.72	0	4,6,6	2.40	1 (25%)
4	MPD	G	2404	-	7,7,7	0.69	0	9,10,10	0.98	0
4	MPD	G	2405	-	7,7,7	0.65	0	9,10,10	0.47	0
4	MPD	A	2407	-	7,7,7	0.68	0	9,10,10	0.34	0
3	CAC	G	2403	2	2,4,4	1.87	1 (50%)	4,6,6	2.78	3 (75%)
4	MPD	A	2406	-	7,7,7	0.62	0	9,10,10	1.11	0
4	MPD	A	2405	-	7,7,7	0.72	0	9,10,10	0.61	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
4	MPD	A	2404	-	-	4/5/5/5	-
4	MPD	G	2404	-	-	1/5/5/5	-
4	MPD	G	2405	-	-	3/5/5/5	-
4	MPD	A	2407	-	-	4/5/5/5	-
4	MPD	A	2406	-	-	0/5/5/5	-
4	MPD	A	2405	-	-	2/5/5/5	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	2403	CAC	AS-C1	2.28	1.95	1.90

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2403	CAC	O1-AS-C2	-4.57	105.81	111.50
3	G	2403	CAC	O1-AS-C1	-3.92	106.62	111.50
3	G	2403	CAC	O1-AS-C2	3.21	115.49	111.50
3	G	2403	CAC	O2-AS-C1	2.30	111.40	105.84

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	2404	MPD	C2-C3-C4-O4
4	A	2404	MPD	C2-C3-C4-C5
4	A	2407	MPD	C1-C2-C3-C4
4	A	2407	MPD	C2-C3-C4-O4
4	A	2407	MPD	C2-C3-C4-C5
4	G	2404	MPD	O2-C2-C3-C4
4	A	2405	MPD	C2-C3-C4-O4
4	A	2404	MPD	C1-C2-C3-C4
4	G	2405	MPD	C1-C2-C3-C4
4	G	2405	MPD	CM-C2-C3-C4
4	A	2405	MPD	C2-C3-C4-C5
4	G	2405	MPD	O2-C2-C3-C4
4	A	2404	MPD	CM-C2-C3-C4

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Mol	Chain	Res	Type	Atoms
4	A	2407	MPD	CM-C2-C3-C4

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	2404	MPD	1	0
4	G	2405	MPD	1	0
4	A	2407	MPD	3	0
4	A	2406	MPD	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	315/333 (94%)	0.08	21 (6%) 24 26	8, 21, 41, 72	24 (7%)
1	G	311/333 (93%)	0.20	22 (7%) 22 24	9, 25, 44, 62	25 (8%)
All	All	626/666 (93%)	0.14	43 (6%) 23 25	8, 24, 42, 72	49 (7%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	162	GLY	8.6
1	G	362	LEU	4.6
1	G	162	GLY	4.2
1	A	362	LEU	4.1
1	A	34	GLY	4.1
1	A	235[A]	ASP	3.6
1	G	35[A]	ASP	3.5
1	G	203	ALA	3.4
1	A	259	ALA	3.3
1	G	204	ALA	3.3
1	A	363	ARG	3.3
1	A	260	ILE	3.2
1	A	46	ILE	3.1
1	A	234	ALA	3.1
1	A	51	PHE	3.0
1	G	51	PHE	2.9
1	A	232	ASP	2.9
1	G	205	SER	2.9
1	G	259	ALA	2.8
1	G	46	ILE	2.8
1	A	108	ARG	2.7
1	A	204	ALA	2.6
1	A	109	ASP	2.6
1	G	361	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	205	SER	2.6
1	A	233	ALA	2.4
1	G	258	SER	2.4
1	G	163	ILE	2.4
1	A	203	ALA	2.3
1	A	275	ARG	2.3
1	G	89	ARG	2.3
1	G	139	ARG	2.3
1	G	232	ASP	2.2
1	G	298[A]	VAL	2.2
1	G	174	GLY	2.2
1	A	361	THR	2.2
1	G	310	VAL	2.2
1	A	174	GLY	2.1
1	A	207[A]	ARG	2.1
1	G	211[A]	GLN	2.1
1	G	175	LYS	2.1
1	G	173	THR	2.0
1	G	234	ALA	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
1	KCX	G	169	12/13	0.94	0.08	18,21,23,23	0
1	KCX	A	169	12/13	0.97	0.05	13,15,17,19	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($\text{\AA}^2$)	Q<0.9
4	MPD	A	2405	8/8	0.58	0.31	34,39,42,44	8
4	MPD	A	2407	8/8	0.76	0.21	42,45,48,50	8
4	MPD	G	2404	8/8	0.80	0.17	33,41,47,48	0
4	MPD	G	2405	8/8	0.81	0.13	47,50,53,54	0
4	MPD	A	2404	8/8	0.85	0.14	35,42,48,48	0
4	MPD	A	2406	8/8	0.88	0.16	28,31,35,40	8
2	ZN	G	2401	1/1	0.97	0.04	19,19,19,19	1
2	ZN	G	2402	1/1	0.99	0.04	19,19,19,19	1
3	CAC	A	2403	5/5	0.99	0.08	14,14,18,18	5
3	CAC	G	2403	5/5	0.99	0.08	17,18,21,23	5
2	ZN	A	2401	1/1	1.00	0.03	16,16,16,16	0
2	ZN	A	2402	1/1	1.00	0.02	16,16,16,16	1

6.5 Other polymers [i](#)

There are no such residues in this entry.