



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 19, 2016 – 02:26 PM EDT

PDB ID : 4XAY
Title : Cycles of destabilization and repair underlie evolutionary transitions in enzymes
Authors : Jackson, C.J.; Campbell, E.; Kaltenbach, M.; Tokuriki, N.
Deposited on : 2014-12-16
Resolution : 1.84 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027939
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : rb-20027939

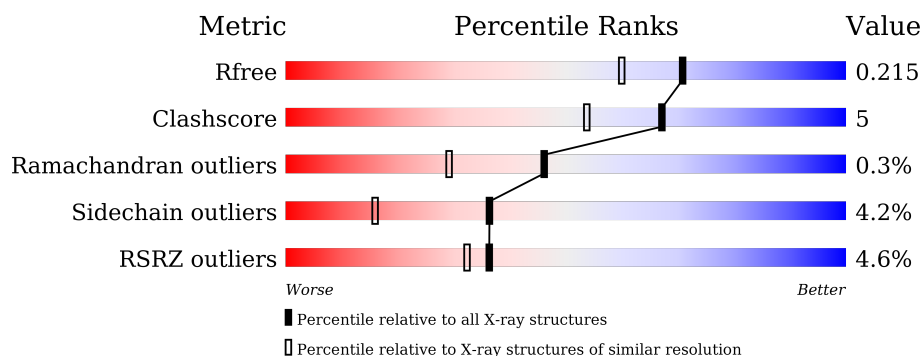
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.84 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}	91344	2634 (1.86-1.82)
Clashscore	102246	2862 (1.86-1.82)
Ramachandran outliers	100387	2831 (1.86-1.82)
Sidechain outliers	100360	2832 (1.86-1.82)
RSRZ outliers	91569	2639 (1.86-1.82)

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	333	<div> <div>7%</div> <div>86%</div> <div>12%</div> <div>..</div> </div>
1	G	333	<div> <div>2%</div> <div>85%</div> <div>13%</div> <div>..</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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	MPD	A	2403	-	-	-	X
4	CAC	G	401	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5573 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.

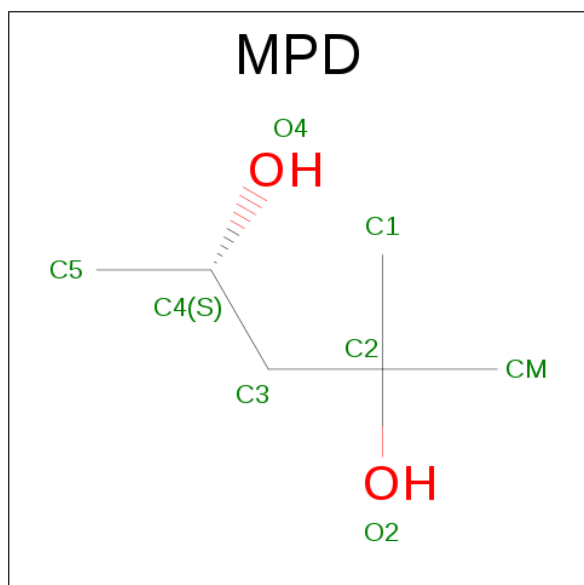
- Molecule 1 is a protein called Phosphotriesterase variant PTE-R8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	8	0
			2578	1626	462	484	6			
1	G	329	Total	C	N	O	S	0	11	0
			2607	1644	469	488	6			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

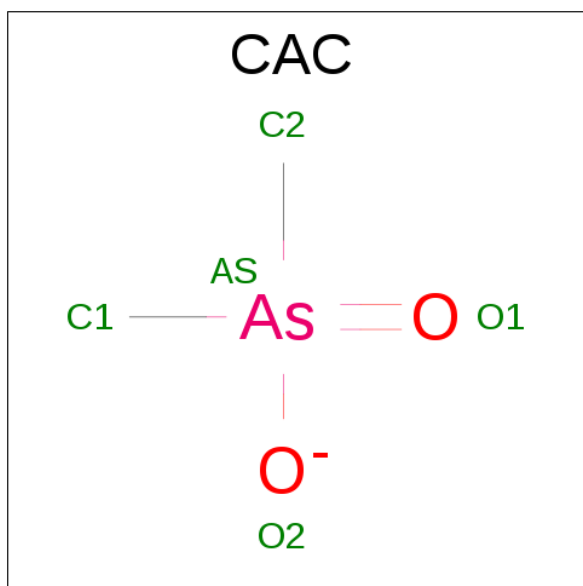
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	2	Total	Zn	0	0
			2	2		
2	A	2	Total	Zn	0	0
			2	2		

- Molecule 3 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: C₆H₁₄O₂).



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

- Molecule 4 is CACODYLATE ION (three-letter code: CAC) (formula: $C_2H_6AsO_2$).



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

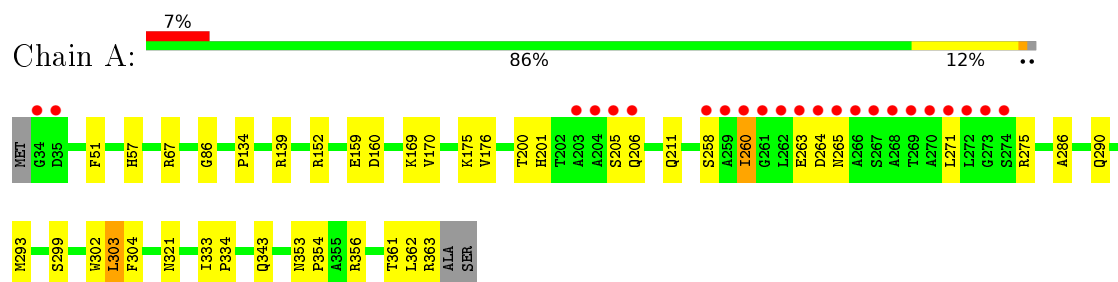
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	189	Total	O	0	0
			189	189		
5	G	182	Total	O	0	0
			182	182		

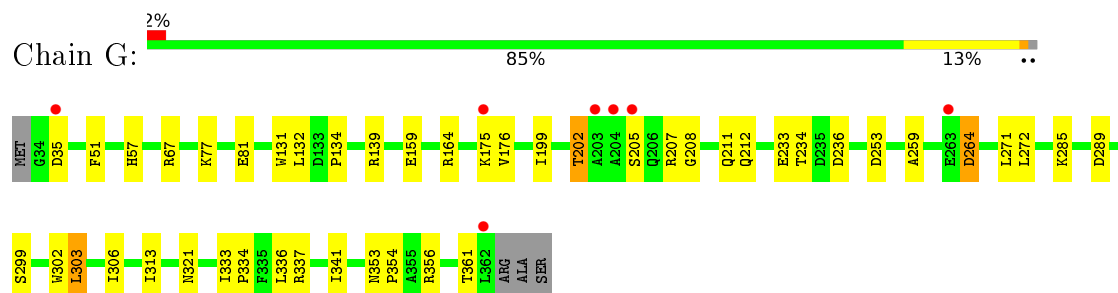
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 variant PTE-R8



- Molecule 1: Phosphotriesterase variant PTE-R8



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	85.76Å 85.92Å 88.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.41 – 1.84 44.34 – 1.84	Depositor EDS
% Data completeness (in resolution range)	99.9 (38.41-1.84) 92.0 (44.34-1.84)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.86 (at 1.84Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.176 , 0.219 0.173 , 0.215	Depositor DCC
R_{free} test set	2651 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	19.6	Xtriage
Anisotropy	0.596	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 53.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.016 for -h,l,k 0.015 for -l,-k,-h 0.016 for k,h,-l 0.003 for k,l,h 0.003 for l,h,k	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5573	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 4.46% 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: CAC, ZN, MPD, KCX

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	0.37	0/2610	0.55	0/3544
1	G	0.37	0/2639	0.54	0/3583
All	All	0.37	0/5249	0.54	0/7127

There are no bond length outliers.

There are no bond angle outliers.

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	2578	0	2609	20	0
1	G	2607	0	2641	33	0
2	A	2	0	0	0	0
2	G	2	0	0	0	0
3	A	8	0	14	1	0
4	G	5	0	0	0	0
5	A	189	0	0	1	2
5	G	182	0	0	6	0
All	All	5573	0	5264	50	2

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 (50) 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:343:GLN:HE22	3:A:2403:MPD:HM2	1.39	0.87
1:A:170:VAL:HG23	1:A:200[A]:THR:HG23	1.68	0.76
1:A:160:ASP:OD1	1:G:67[B]:ARG:NH1	2.26	0.68
1:G:253:ASP:OD2	5:G:646:HOH:O	2.13	0.66
1:G:57:HIS:HB2	1:G:303[B]:LEU:HB3	1.80	0.63
1:G:164:ARG:NH1	5:G:502:HOH:O	2.33	0.61
1:A:159:GLU:HG2	1:G:67[B]:ARG:HG2	1.82	0.60
1:G:264:ASP:OD1	1:G:264:ASP:N	2.28	0.57
1:A:361:THR:HG22	1:A:363:ARG:H	1.69	0.56
1:A:363:ARG:NH1	5:A:2501:HOH:O	2.39	0.55
1:G:175:LYS:HA	1:G:211[C]:GLN:HE22	1.70	0.55
1:G:353:ASN:OD1	1:G:356:ARG:NH2	2.39	0.55
1:G:77:LYS:NZ	1:G:81:GLU:OE1	2.38	0.54
1:G:205:SER:HA	1:G:233:GLU:HB3	1.90	0.53
1:A:57:HIS:HB2	1:A:303:LEU:HB3	1.90	0.53
1:G:302:TRP:CH2	1:G:321:ASN:HB3	2.45	0.51
1:G:207:ARG:HD3	1:G:234[A]:THR:HG21	1.93	0.51
1:G:57:HIS:HB2	1:G:303[A]:LEU:HB3	1.92	0.51
1:G:207:ARG:O	5:G:653:HOH:O	2.19	0.50
1:A:169:KCX:HE2	1:A:201:HIS:HB2	1.93	0.50
1:G:57:HIS:O	1:G:303[A]:LEU:HA	2.13	0.48
1:A:258:SER:HB2	1:A:275:ARG:O	2.13	0.48
1:A:333:ILE:HB	1:A:334:PRO:HD3	1.96	0.48
1:G:259:ALA:HB2	1:G:272:LEU:HB2	1.96	0.48
1:G:306:ILE:HD12	1:G:313:ILE:HG23	1.96	0.47
1:A:67[B]:ARG:HG2	1:G:159:GLU:HG2	1.98	0.46
1:A:134:PRO:HG2	1:A:139:ARG:HG2	1.97	0.46
1:G:356:ARG:NE	5:G:632:HOH:O	2.36	0.46
1:G:202:THR:OG1	5:G:655:HOH:O	2.07	0.45
1:A:302:TRP:CH2	1:A:321:ASN:HB3	2.52	0.45
1:G:131:TRP:CG	1:G:132:LEU:N	2.85	0.44
1:A:169:KCX:OQ1	1:A:201:HIS:HB2	2.17	0.44
1:G:57:HIS:O	1:G:303[B]:LEU:HA	2.17	0.43
1:G:199:ILE:HD11	5:G:646:HOH:O	2.18	0.43
1:A:57:HIS:O	1:A:303:LEU:HA	2.18	0.43
1:G:208:GLY:O	1:G:212:GLN:HG3	2.19	0.42
1:G:207:ARG:HG3	1:G:234[B]:THR:HG21	2.00	0.42
1:G:234[A]:THR:HG22	1:G:236:ASP:H	1.84	0.42
1:G:134:PRO:HG2	1:G:139:ARG:HG2	2.01	0.41
1:A:175:LYS:O	1:A:211:GLN:HG2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:176:VAL:N	1:G:211[C]:GLN:OE1	2.52	0.41
1:G:285:LYS:NZ	1:G:289:ASP:OD2	2.50	0.41
1:G:303[A]:LEU:HD13	1:G:306:ILE:HG12	2.03	0.41
1:A:353:ASN:HB2	1:A:354:PRO:HD3	2.02	0.41
1:A:362:LEU:HA	1:A:362:LEU:HD23	1.83	0.41
1:G:333:ILE:HB	1:G:334:PRO:HD3	2.01	0.41
1:G:336:LEU:HB3	1:G:341:ILE:HD13	2.02	0.41
1:G:353:ASN:HB2	1:G:354:PRO:HD3	2.02	0.40
1:A:286:ALA:O	1:A:290:GLN:HG2	2.21	0.40
1:A:86:GLY:HA3	1:A:304:PHE:CE1	2.57	0.40

All (2) 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)
5:A:2506:HOH:O	5:A:2531:HOH:O[3_445]	2.11	0.09
5:A:2528:HOH:O	5:A:2533:HOH:O[3_445]	2.17	0.03

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%)	317 (95%)	16 (5%)	2 (1%)	30	14
1	G	338/333 (102%)	324 (96%)	14 (4%)	0	100	100
All	All	673/666 (101%)	641 (95%)	30 (4%)	2 (0%)	46	29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	260	ILE
1	A	176	VAL

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	272/266 (102%)	259 (95%)	13 (5%)	31	12
1	G	275/266 (103%)	264 (96%)	11 (4%)	38	18
All	All	547/532 (103%)	523 (96%)	24 (4%)	36	15

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	PHE
1	A	152	ARG
1	A	205	SER
1	A	206	GLN
1	A	260	ILE
1	A	263	GLU
1	A	264	ASP
1	A	265	ASN
1	A	271	LEU
1	A	293	MET
1	A	299	SER
1	A	303	LEU
1	A	356	ARG
1	G	35	ASP
1	G	51	PHE
1	G	202	THR
1	G	264	ASP
1	G	271	LEU
1	G	299	SER
1	G	303[A]	LEU
1	G	303[B]	LEU
1	G	337[A]	ARG
1	G	337[B]	ARG
1	G	361	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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	6,11,12	0.70	0	7,12,14	2.26	1 (14%)
1	KCX	G	169	1,2	6,11,12	0.73	0	7,12,14	1.87	2 (28%)

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	-	0/6/10/12	0/0/0/0
1	KCX	G	169	1,2	-	0/6/10/12	0/0/0/0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	169	KCX	CE-NZ-CX	-5.37	117.89	123.53
1	G	169	KCX	CE-NZ-CX	-4.21	119.10	123.53
1	G	169	KCX	O-C-CA	-2.14	119.97	125.72

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	169	KCX	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	$\# Z > 2$
3	MPD	A	2403	-	6,7,7	0.28	0	6,10,10	0.21	0
4	CAC	G	401	2	0,4,4	0.00	-	0,6,6	0.00	-

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	MPD	A	2403	-	-	0/5/5/5	0/0/0/0
4	CAC	G	401	2	-	0/0/0/0	0/0/0/0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2403	MPD	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	329/333 (98%)	0.21	23 (6%) 19 17	12, 21, 74, 138	0
1	G	328/333 (98%)	-0.26	7 (2%) 67 64	14, 22, 53, 82	0
All	All	657/666 (98%)	-0.03	30 (4%) 36 32	12, 22, 57, 138	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	266	ALA	17.0
1	A	262	LEU	14.5
1	A	265	ASN	13.9
1	A	268	ALA	13.4
1	A	260	ILE	11.2
1	A	271	LEU	10.9
1	A	261	GLY	10.6
1	A	267	SER	9.6
1	A	270	ALA	9.1
1	A	264	ASP	8.2
1	A	274	SER	7.5
1	A	272	LEU	6.8
1	A	258	SER	4.6
1	A	205	SER	4.5
1	G	204	ALA	4.4
1	A	273	GLY	4.4
1	A	35	ASP	4.3
1	A	206	GLN	4.1
1	A	269	THR	3.7
1	A	34	GLY	3.4
1	A	204	ALA	3.1
1	G	35	ASP	3.1
1	G	362	LEU	2.9
1	A	263	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	G	203	ALA	2.5
1	G	175	LYS	2.4
1	A	203	ALA	2.4
1	A	259	ALA	2.1
1	G	205	SER	2.1
1	G	263	GLU	2.1

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. 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
1	KCX	A	169	12/13	0.93	0.10	-	10,19,31,32	0
1	KCX	G	169	12/13	0.93	0.10	-	13,19,30,34	0

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
3	MPD	A	2403	8/8	0.88	0.23	6.00	42,48,55,59	0
4	CAC	G	401	5/5	0.93	0.18	4.52	12,21,32,33	5
2	ZN	A	2402	1/1	0.99	0.06	-	21,21,21,21	0
2	ZN	G	402	1/1	1.00	0.07	-	19,19,19,19	0
2	ZN	A	2401	1/1	1.00	0.09	-	16,16,16,16	0
2	ZN	G	403	1/1	0.98	0.06	-	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.