Full wwPDB X-ray Structure Validation Report (i)

Nov 6, 2023 – 10:51 AM EST

PDB ID	:	5WJ0
Title	:	Phosphotriesterase variant S5+254R
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Deposited on	:	2017-07-21
Resolution	:	1.65 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 (i) 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 (1)) were used in the production of this report:

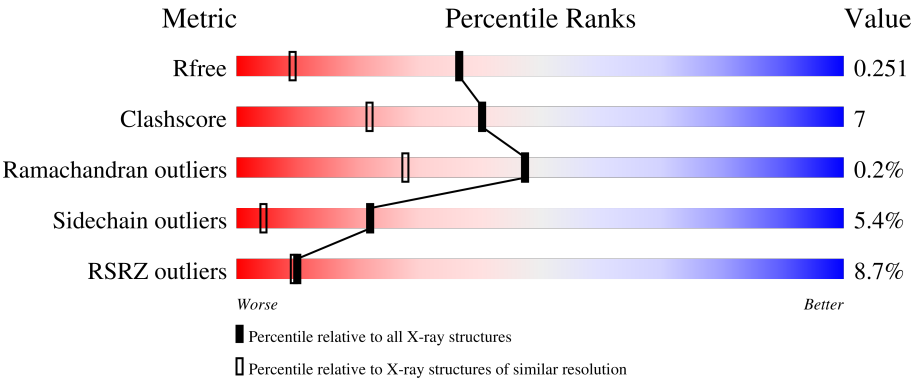
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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}	130704	1827 (1.66-1.66)
Clashscore	141614	1931 (1.66-1.66)
Ramachandran outliers	138981	1891 (1.66-1.66)
Sidechain outliers	138945	1891 (1.66-1.66)
RSRZ outliers	127900	1791 (1.66-1.66)

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 <=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	4% 80% 13% • 5%
1	G	333	13% 78% 14% • 5%

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	2404	-	-	-	X
3	MPD	A	2405	-	-	-	X
3	MPD	G	401	-	-	-	X
3	MPD	G	402	-	-	-	X

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5200 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	4	0
			2444	1544	442	452	6			
1	G	315	Total	C	N	O	S	0	9	0
			2489	1572	453	458	6			

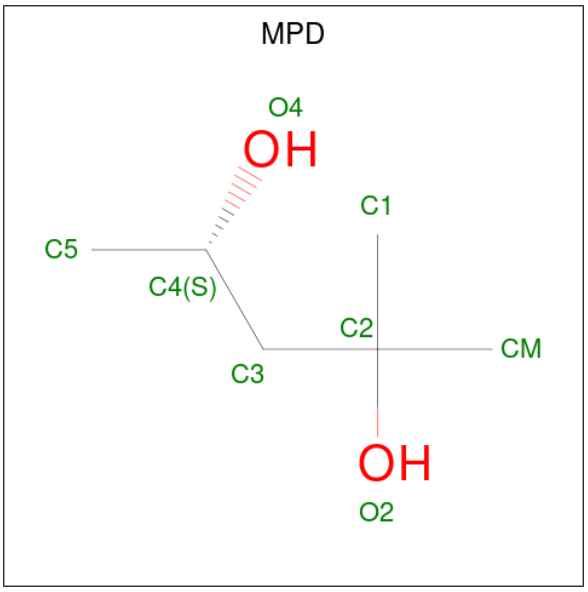
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	216	LEU	PHE	conflict	UNP A0A060GZX0
A	233	ALA	ASP	conflict	UNP A0A060GZX0
A	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	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 (three-letter code: ZN) (formula: Zn).

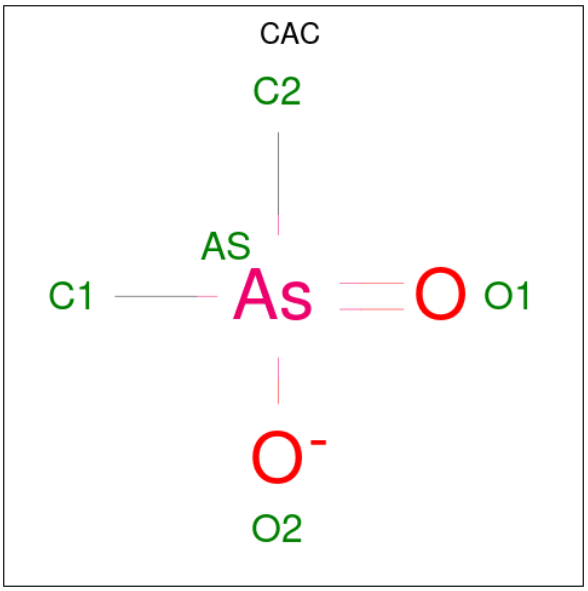
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Zn	0	0
			2	2		
2	G	2	Total	Zn	0	0
			2	2		

- Molecule 3 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



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

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



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

- Molecule 5 is water.

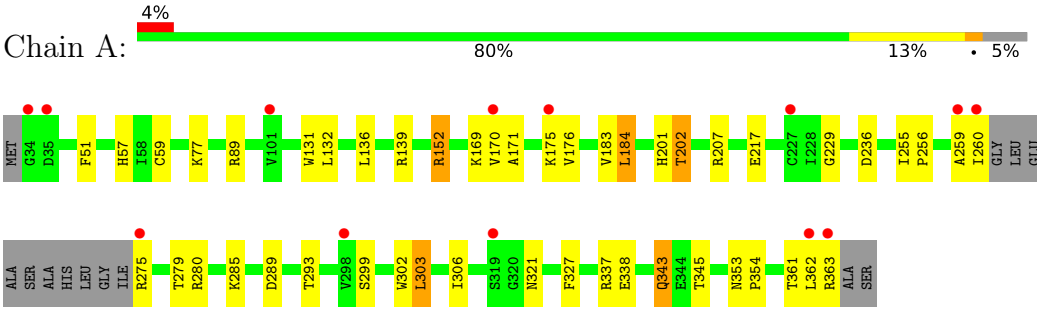
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	113	Total	O	0	0
			113	113		
5	G	92	Total	O	0	0
			92	92		

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.

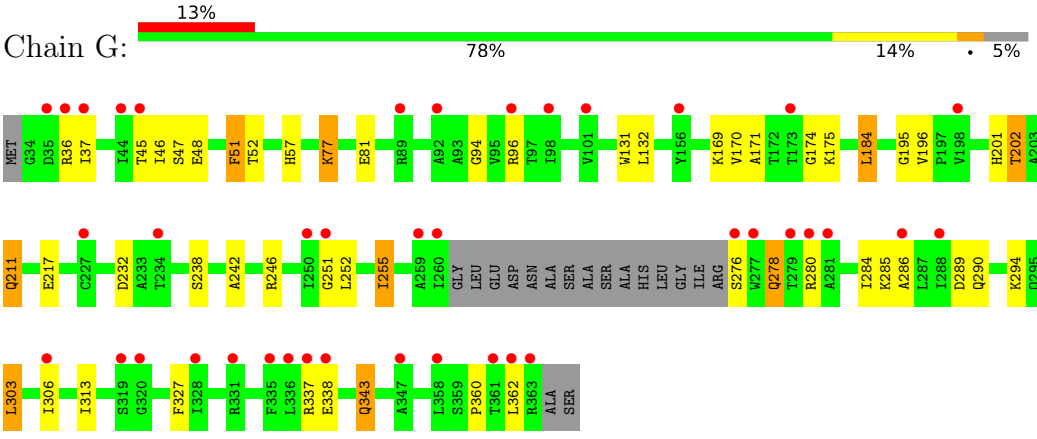
• Molecule 1: Phosphotriesterase

Chain A:



• Molecule 1: Phosphotriesterase

Chain G:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	85.33Å 85.78Å 88.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.43 – 1.65 29.43 – 1.65	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.43-1.65) 99.9 (29.43-1.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.48 (at 1.65Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.215 , 0.248 0.217 , 0.251	Depositor DCC
R_{free} test set	4051 reflections (5.16%)	wwPDB-VP
Wilson B-factor (Å^2)	18.7	Xtriage
Anisotropy	0.455	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 46.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.009 for -h,l,k 0.012 for -l,-k,-h 0.013 for k,h,-l 0.000 for k,l,h 0.000 for l,h,k	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5200	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.78% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: ZN, CAC, KCX, MPD

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	0.47	0/2474	0.67	0/3358
1	G	0.42	0/2520	0.64	0/3418
All	All	0.44	0/4994	0.65	0/6776

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	2444	0	2484	34	0
1	G	2489	0	2523	37	0
2	A	2	0	0	0	0
2	G	2	0	0	0	0
3	A	24	0	42	2	0
3	G	24	0	42	2	0
4	A	5	0	0	0	0
4	G	5	0	0	0	0
5	A	113	0	0	5	0
5	G	92	0	0	4	0
All	All	5200	0	5091	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:170[B]:VAL:HG11	1:A:184:LEU:HD13	1.55	0.89
1:G:170[B]:VAL:HG11	1:G:184:LEU:HD13	1.61	0.82
1:A:170[B]:VAL:HG11	1:A:184:LEU:CD1	2.12	0.80
1:G:170[B]:VAL:HG11	1:G:184:LEU:CD1	2.14	0.78
1:G:202:THR:HG21	5:G:522:HOH:O	1.94	0.67
1:G:251:GLY:HA3	5:G:547:HOH:O	1.96	0.64
1:G:276:SER:HB2	1:G:278[A]:GLN:HE21	1.63	0.62
1:G:36:ARG:HH11	1:G:45:THR:HG23	1.64	0.61
1:A:202:THR:HG21	5:A:2515:HOH:O	2.00	0.61
1:G:217:GLU:OE2	1:G:246[B]:ARG:NH1	2.35	0.60
1:G:57:HIS:HB2	1:G:303:LEU:HB3	1.84	0.59
1:G:284:ILE:HG23	1:G:296:ILE:HG21	1.85	0.58
1:G:303:LEU:HD13	1:G:306[A]:ILE:HG12	1.86	0.57
1:A:57:HIS:HB2	1:A:303:LEU:HB3	1.88	0.56
1:G:51:PHE:HD1	1:G:52:THR:N	2.04	0.56
3:A:2403:MPD:O4	3:A:2403:MPD:O2	2.24	0.54
1:A:361:THR:HG22	1:A:363:ARG:H	1.73	0.54
1:A:362:LEU:HD23	5:A:2605:HOH:O	2.08	0.53
1:G:337:ARG:HH21	1:G:343:GLN:HG2	1.74	0.52
1:G:174:GLY:O	1:G:211:GLN:NE2	2.43	0.52
1:A:170[A]:VAL:HG21	1:A:184:LEU:CD1	2.40	0.51
1:A:337:ARG:HH12	1:A:343:GLN:HG2	1.75	0.51
1:A:170[B]:VAL:CG1	1:A:184:LEU:HD13	2.34	0.51
1:A:59:CYS:HB3	5:A:2593:HOH:O	2.11	0.50
1:G:57:HIS:O	1:G:303:LEU:HA	2.11	0.50
1:A:255:ILE:HD13	1:A:280:ARG:HB3	1.94	0.50
1:A:337:ARG:HH22	1:A:343:GLN:HG2	1.76	0.50
3:G:401:MPD:O4	3:G:401:MPD:O2	2.08	0.50
1:A:259:ALA:O	1:A:260:ILE:HB	2.13	0.48
1:G:285:LYS:NZ	1:G:289:ASP:OD2	2.46	0.48
1:A:275:ARG:HH21	1:A:279:THR:HG22	1.79	0.48
1:A:255:ILE:HG12	5:A:2502:HOH:O	2.13	0.47
1:A:57:HIS:O	1:A:303:LEU:HA	2.14	0.47
1:A:337:ARG:HH22	1:A:343:GLN:CG	2.26	0.47
1:A:256:PRO:HG3	1:A:327:PHE:CD1	2.50	0.47
1:A:89:ARG:HG3	5:A:2538:HOH:O	2.14	0.46
1:G:37:ILE:HG13	1:G:46:ILE:HG22	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:174:GLY:HA3	1:G:206:GLN:OE1	2.15	0.46
1:G:51:PHE:C	1:G:51:PHE:CD1	2.89	0.46
1:G:303:LEU:HD22	1:G:306[B]:ILE:HG23	1.97	0.46
1:G:232:ASP:OD2	1:G:280:ARG:NH1	2.48	0.45
1:A:302:TRP:CH2	1:A:321:ASN:HB3	2.52	0.45
1:A:217:GLU:HG2	3:A:2405:MPD:H53	1.98	0.45
1:G:306[B]:ILE:CD1	1:G:313:ILE:HG23	2.47	0.44
1:G:303:LEU:HD22	1:G:306[B]:ILE:CG2	2.48	0.44
1:G:51:PHE:HD1	1:G:51:PHE:C	2.21	0.44
1:G:242:ALA:O	1:G:246[B]:ARG:HG3	2.18	0.44
1:G:286:ALA:O	1:G:290:GLN:HG2	2.18	0.44
1:A:207:ARG:NE	1:A:236:ASP:OD2	2.50	0.44
1:A:255:ILE:HB	1:A:256:PRO:HD3	1.99	0.44
1:G:252:LEU:HD12	1:G:284:ILE:HG12	2.01	0.43
1:G:232:ASP:HB3	1:G:252:LEU:HA	2.01	0.43
1:A:285:LYS:NZ	1:A:289:ASP:OD2	2.41	0.42
1:G:77:LYS:HE2	1:G:77:LYS:HB3	1.58	0.42
1:G:171:ALA:HA	1:G:201:HIS:HB3	2.01	0.42
1:A:170[B]:VAL:HG13	1:A:183:VAL:HG12	2.02	0.42
1:A:152:ARG:O	1:A:152:ARG:HG2	2.20	0.42
1:A:136:LEU:HG	1:A:139:ARG:NH2	2.35	0.41
1:A:171:ALA:HA	1:A:201:HIS:HB3	2.02	0.41
1:A:184:LEU:HA	1:A:184:LEU:HD12	1.78	0.41
1:G:94:GLY:O	1:G:96[B]:ARG:NE	2.49	0.41
1:A:131:TRP:CG	1:A:132:LEU:N	2.89	0.41
1:A:303:LEU:HD11	1:A:306[A]:ILE:HD11	2.02	0.41
1:A:353:ASN:HB2	1:A:354:PRO:HD3	2.01	0.41
1:G:81:GLU:HG2	3:G:401:MPD:O2	2.20	0.41
1:G:51:PHE:HB2	1:G:96[B]:ARG:HH21	1.85	0.41
1:G:251:GLY:CA	5:G:547:HOH:O	2.60	0.41
1:G:48:GLU:HG2	5:G:588:HOH:O	2.21	0.41
1:G:255:ILE:HB	1:G:327[B]:PHE:HE1	1.85	0.41
1:G:195:GLY:O	1:G:360:PRO:HA	2.21	0.41
1:A:202:THR:HG22	1:A:229:GLY:O	2.21	0.40
1:A:293:THR:HG21	1:A:345:THR:HG23	2.04	0.40
1:G:131:TRP:CG	1:G:132:LEU:N	2.90	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	315/333 (95%)	303 (96%)	11 (4%)	1 (0%)	41 22
1	G	319/333 (96%)	309 (97%)	10 (3%)	0	100 100
All	All	634/666 (95%)	612 (96%)	21 (3%)	1 (0%)	47 28

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	176	VAL

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	256/264 (97%)	246 (96%)	10 (4%)	32 9
1	G	260/264 (98%)	242 (93%)	18 (7%)	15 2
All	All	516/528 (98%)	488 (95%)	28 (5%)	22 4

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	PHE
1	A	77	LYS
1	A	152	ARG
1	A	175	LYS
1	A	184	LEU

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Mol	Chain	Res	Type
1	A	202	THR
1	A	299	SER
1	A	303	LEU
1	A	338	GLU
1	A	343	GLN
1	G	47	SER
1	G	51	PHE
1	G	77	LYS
1	G	175	LYS
1	G	184	LEU
1	G	196	VAL
1	G	202	THR
1	G	211	GLN
1	G	238	SER
1	G	255	ILE
1	G	278[A]	GLN
1	G	278[B]	GLN
1	G	294	LYS
1	G	299	SER
1	G	303	LEU
1	G	338	GLU
1	G	343	GLN
1	G	362	LEU

Sometimes 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	G	169	2,1	9,11,12	2.48	2 (22%)	5,12,14	0.67	0
1	KCX	A	169	2,1	9,11,12	1.72	2 (22%)	5,12,14	2.19	2 (40%)

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

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	169	KCX	CX-NZ	5.95	1.45	1.35
1	A	169	KCX	CX-NZ	4.24	1.42	1.35
1	G	169	KCX	OQ1-CX	3.99	1.29	1.21
1	A	169	KCX	OQ1-CX	2.13	1.25	1.21

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	169	KCX	OQ1-CX-NZ	-3.51	119.52	124.96
1	A	169	KCX	CE-NZ-CX	-2.32	118.16	121.89

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 monosaccharides 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	$\#\|Z\| > 2$
3	MPD	G	402	-	7,7,7	0.72	0	9,10,10	0.26	0
3	MPD	A	2404	-	7,7,7	0.77	0	9,10,10	0.66	0
4	CAC	A	2406	2	0,4,4	-	-	0,6,6	-	-
3	MPD	A	2403	-	7,7,7	0.60	0	9,10,10	0.55	0
4	CAC	G	406	2	0,4,4	-	-	0,6,6	-	-
3	MPD	G	405	-	7,7,7	0.78	0	9,10,10	0.65	0
3	MPD	G	401	-	7,7,7	0.67	0	9,10,10	0.44	0
3	MPD	A	2405	-	7,7,7	0.73	0	9,10,10	0.66	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
3	MPD	G	402	-	-	2/5/5/5	-
3	MPD	A	2404	-	-	1/5/5/5	-
3	MPD	A	2403	-	-	4/5/5/5	-
3	MPD	G	405	-	-	1/5/5/5	-
3	MPD	G	401	-	-	2/5/5/5	-
3	MPD	A	2405	-	-	1/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	2403	MPD	O2-C2-C3-C4
3	A	2405	MPD	C2-C3-C4-C5
3	G	401	MPD	C2-C3-C4-C5
3	G	402	MPD	C2-C3-C4-C5

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Mol	Chain	Res	Type	Atoms
3	G	405	MPD	C2-C3-C4-C5
3	G	402	MPD	C2-C3-C4-O4
3	A	2403	MPD	C1-C2-C3-C4
3	A	2403	MPD	CM-C2-C3-C4
3	A	2404	MPD	C2-C3-C4-C5
3	A	2403	MPD	C2-C3-C4-O4
3	G	401	MPD	C2-C3-C4-O4

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2403	MPD	1	0
3	G	401	MPD	2	0
3	A	2405	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, 95^{th} 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%)	0.32	13 (4%) 37 37	12, 20, 39, 64	0
1	G	314/333 (94%)	0.87	42 (13%) 3 2	15, 30, 50, 74	0
All	All	629/666 (94%)	0.59	55 (8%) 10 9	12, 26, 46, 74	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	363	ARG	8.0
1	G	362	LEU	7.6
1	A	34	GLY	7.4
1	A	275	ARG	6.1
1	G	260	ILE	5.9
1	A	260	ILE	5.5
1	A	35	ASP	5.4
1	A	362	LEU	4.9
1	G	37	ILE	4.6
1	G	45	THR	4.5
1	G	259	ALA	4.4
1	G	204	ALA	3.8
1	A	363	ARG	3.8
1	G	156	TYR	3.6
1	G	36	ARG	3.4
1	G	101	VAL	3.4
1	A	259	ALA	3.4
1	G	328	ILE	3.3
1	G	35	ASP	3.2
1	G	335	PHE	3.1
1	A	227	CYS	3.0
1	G	337	ARG	3.0
1	A	175	LYS	2.9
1	G	205	SER	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	319	SER	2.7
1	G	277	TRP	2.7
1	G	336	LEU	2.7
1	G	234	THR	2.7
1	G	358	LEU	2.6
1	G	361	THR	2.6
1	G	96[A]	ARG	2.6
1	A	170[A]	VAL	2.6
1	G	173	THR	2.5
1	G	276	SER	2.5
1	G	320	GLY	2.5
1	G	251	GLY	2.4
1	G	331	ARG	2.4
1	G	338	GLU	2.4
1	G	227	CYS	2.4
1	G	92	ALA	2.4
1	G	89	ARG	2.3
1	G	281	ALA	2.2
1	G	280	ARG	2.2
1	A	101	VAL	2.2
1	G	98	ILE	2.2
1	G	288	ILE	2.2
1	G	286	ALA	2.1
1	G	279	THR	2.1
1	G	198	VAL	2.1
1	G	347	ALA	2.1
1	A	298	VAL	2.1
1	G	44	ILE	2.1
1	G	250	ILE	2.1
1	G	319	SER	2.1
1	G	306[A]	ILE	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, 95^{th} 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	0.87	0.17	20,23,28,30	0
1	KCX	A	169	12/13	0.95	0.14	9,13,15,16	0

6.3 Carbohydrates (i)

There are no monosaccharides 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, 95^{th} 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	MPD	G	402	8/8	0.16	0.59	57,61,62,62	0
3	MPD	A	2404	8/8	0.23	0.62	54,60,63,69	0
3	MPD	A	2405	8/8	0.27	0.52	27,48,52,56	0
3	MPD	G	401	8/8	0.44	0.41	35,43,49,51	0
3	MPD	G	405	8/8	0.69	0.26	54,55,56,58	0
3	MPD	A	2403	8/8	0.77	0.20	40,44,50,54	0
4	CAC	A	2406	5/5	0.85	0.26	11,37,40,43	5
4	CAC	G	406	5/5	0.86	0.26	19,32,41,43	5
2	ZN	G	403	1/1	0.94	0.04	21,21,21,21	0
2	ZN	G	404	1/1	0.96	0.04	27,27,27,27	0
2	ZN	A	2402	1/1	0.98	0.06	19,19,19,19	0
2	ZN	A	2401	1/1	0.99	0.10	14,14,14,14	0

6.5 Other polymers (i)

There are no such residues in this entry.