



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:08 PM GMT

PDB ID : 1S95
Title : Structure of serine/threonine protein phosphatase 5
Authors : Swingle, M.R.; Honkanen, R.E.; Ciszak, E.M.
Deposited on : 2004-02-03
Resolution : 1.60 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

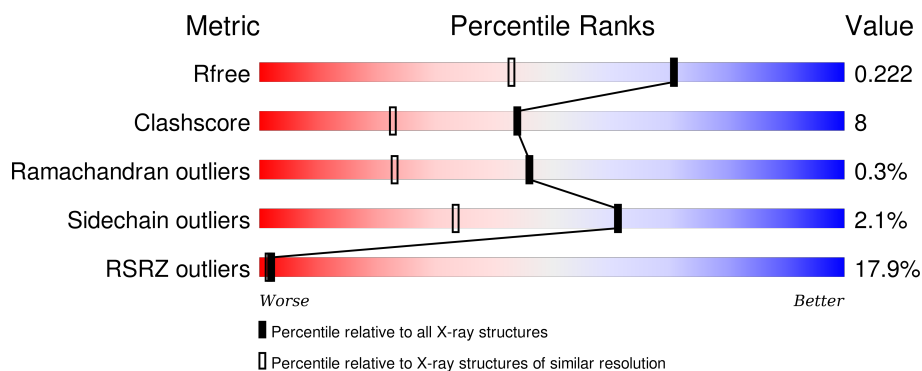
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	2475 (1.60-1.60)
Clashscore	102246	2732 (1.60-1.60)
Ramachandran outliers	100387	2654 (1.60-1.60)
Sidechain outliers	100360	2653 (1.60-1.60)
RSRZ outliers	91569	2479 (1.60-1.60)

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	
1	B	333	

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
4	MPD	A	3001	-	-	-	X
4	MPD	B	3002	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5748 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 Serine/threonine protein phosphatase 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	324	Total	C	N	O	S	0	7	0
			2624	1675	437	494	18			
1	B	325	Total	C	N	O	S	0	7	0
			2631	1681	439	491	20			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	167	GLY	-	CLONING ARTIFACT	UNP P53041
A	168	ALA	-	CLONING ARTIFACT	UNP P53041
B	167	GLY	-	CLONING ARTIFACT	UNP P53041
B	168	ALA	-	CLONING ARTIFACT	UNP P53041

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



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

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



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

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

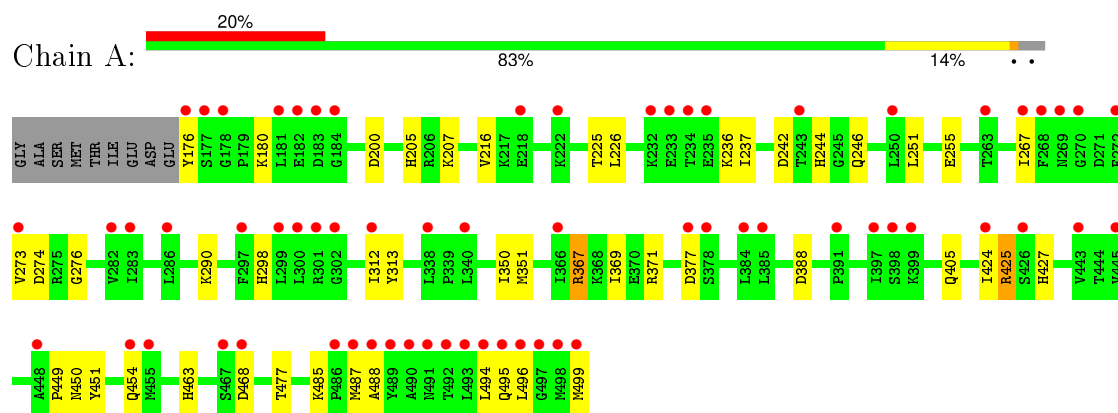
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	233	Total	O	0	0
			233	233		
5	B	230	Total	O	0	0
			230	230		

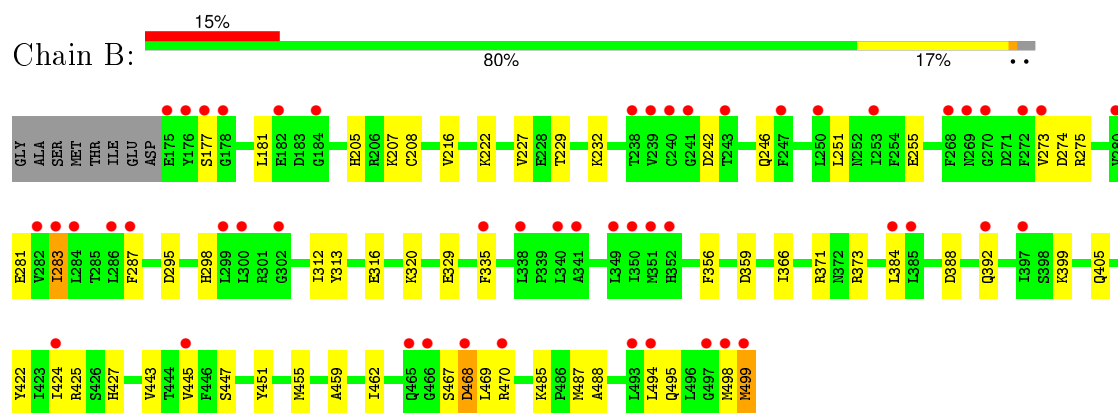
3 Residue-property plots [i](#)

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: Serine/threonine protein phosphatase 5



- Molecule 1: Serine/threonine protein phosphatase 5



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	40.81Å 80.31Å 92.19Å 90.00° 94.25° 90.00°	Depositor
Resolution (Å)	40.00 – 1.60 30.65 – 1.59	Depositor EDS
% Data completeness (in resolution range)	86.5 (40.00-1.60) 86.5 (30.65-1.59)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 1.59Å)	Xtriage
Refinement program	REFMAC 5.1.9999	Depositor
R, R_{free}	0.169 , 0.209 0.182 , 0.222	Depositor DCC
R_{free} test set	3913 reflections (6.01%)	DCC
Wilson B-factor (Å ²)	21.6	Xtriage
Anisotropy	0.170	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 68888 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5748	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 4.87% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: PO4, MN, 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	0.89	0/2716	0.91	5/3671 (0.1%)
1	B	0.95	1/2723 (0.0%)	0.95	5/3679 (0.1%)
All	All	0.92	1/5439 (0.0%)	0.93	10/7350 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	281	GLU	CD-OE1	5.61	1.31	1.25

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	295	ASP	CB-CG-OD2	7.94	125.45	118.30
1	B	373	ARG	NE-CZ-NH1	6.91	123.75	120.30
1	A	367	ARG	NE-CZ-NH2	-6.59	117.00	120.30
1	A	200	ASP	CB-CG-OD2	6.39	124.05	118.30
1	A	377	ASP	CB-CG-OD2	6.05	123.75	118.30
1	B	359	ASP	CB-CG-OD2	5.64	123.38	118.30
1	B	295	ASP	CB-CG-OD1	-5.49	113.36	118.30
1	A	468	ASP	CB-CG-OD2	5.25	123.03	118.30
1	B	371	ARG	NE-CZ-NH2	-5.21	117.69	120.30
1	A	371	ARG	NE-CZ-NH2	-5.06	117.77	120.30

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	2624	0	2554	35	0
1	B	2631	0	2573	55	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	8	0	14	0	0
4	B	8	0	14	0	0
5	A	233	0	0	5	0
5	B	230	0	0	8	0
All	All	5748	0	5155	84	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 (84) 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:424:ILE:HD13	1:B:443:VAL:HB	1.55	0.88
1:A:499:MET:HG2	1:B:399:LYS:HB3	1.67	0.76
1:B:366:ILE:HD12	1:B:384:LEU:HD11	1.65	0.76
1:B:485[B]:LYS:HG2	5:B:3079:HOH:O	1.85	0.74
1:B:312[B]:ILE:CD1	1:B:494:LEU:HD11	2.21	0.71
1:B:205:HIS:HD2	1:B:207:LYS:H	1.36	0.71
1:B:356:PHE:CZ	1:B:366:ILE:HD11	2.26	0.71
1:B:312[B]:ILE:HG13	1:B:494:LEU:HD11	1.74	0.70
1:B:283:ILE:HD11	1:B:335:PHE:HE1	1.56	0.69
1:B:356:PHE:HZ	1:B:366:ILE:HD11	1.57	0.69
1:A:477:THR:HG23	5:A:3232:HOH:O	1.94	0.67
1:B:468:ASP:O	1:B:470:ARG:N	2.28	0.66
1:B:485[B]:LYS:NZ	5:B:3157:HOH:O	2.28	0.65
5:A:3175:HOH:O	1:B:499:MET:HG3	1.97	0.64
1:B:366:ILE:CD1	1:B:384:LEU:HD11	2.27	0.64
1:A:388:ASP:O	1:A:405:GLN:HA	2.00	0.62
1:B:312[B]:ILE:CG1	1:B:494:LEU:HD11	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:267:ILE:HD12	1:A:298:HIS:HB2	1.82	0.61
1:A:454:GLN:HG3	1:B:455[B]:MET:CE	2.32	0.60
1:A:236:LYS:HZ1	1:A:463:HIS:CE1	2.20	0.59
1:B:485[B]:LYS:HG3	1:B:488:ALA:HB2	1.86	0.58
1:A:454:GLN:HG3	1:B:455[B]:MET:HE2	1.86	0.57
1:A:236:LYS:NZ	1:A:463:HIS:CE1	2.72	0.57
1:A:225:THR:HG21	1:A:369:ILE:HB	1.87	0.57
1:B:445:VAL:HG21	1:B:462:ILE:HD13	1.85	0.57
1:A:487:MET:CE	1:A:494:LEU:HB3	2.35	0.56
5:A:3190:HOH:O	1:B:499:MET:HA	2.06	0.55
1:A:487:MET:HE3	1:A:494:LEU:HB3	1.88	0.55
1:B:283:ILE:CD1	1:B:335:PHE:HE1	2.20	0.53
1:B:312[B]:ILE:HG13	1:B:494:LEU:CD1	2.37	0.53
1:A:273:VAL:O	1:A:274:ASP:HB2	2.09	0.53
1:B:275:ARG:HA	1:B:487[B]:MET:HG3	1.92	0.52
1:A:451:TYR:HE2	1:B:499:MET:CE	2.23	0.51
1:A:427:HIS:HA	5:A:3002:HOH:O	2.10	0.51
1:A:237:ILE:HD11	1:A:267:ILE:HD11	1.93	0.51
1:A:176:TYR:HA	1:A:180:LYS:HD3	1.93	0.51
1:A:205:HIS:HD2	1:A:207:LYS:H	1.59	0.51
1:A:267:ILE:CD1	1:A:298:HIS:ND1	2.75	0.50
1:B:298:HIS:HD2	5:B:3025:HOH:O	1.92	0.50
1:B:312[B]:ILE:HD12	1:B:494:LEU:HD11	1.94	0.50
1:A:451:TYR:CE2	1:B:499:MET:CE	2.95	0.50
1:B:388:ASP:O	1:B:405:GLN:HA	2.13	0.49
1:B:468:ASP:C	1:B:470:ARG:H	2.16	0.49
1:B:447:SER:HA	1:B:459:ALA:HB1	1.94	0.49
1:A:205:HIS:CD2	1:A:207:LYS:H	2.31	0.48
1:B:205:HIS:CD2	1:B:207:LYS:H	2.25	0.48
1:A:495:GLN:HG2	1:A:496:LEU:HD23	1.96	0.48
1:B:498:MET:HA	1:B:499:MET:HE2	1.96	0.48
1:A:244:HIS:O	1:A:276:GLY:HA3	2.14	0.47
1:A:451:TYR:CE2	1:B:499:MET:HE1	2.50	0.47
1:B:468:ASP:OD1	1:B:470:ARG:HG3	2.15	0.47
1:B:246:GLN:HE22	1:B:451:TYR:HA	1.79	0.47
1:A:351:MET:O	1:A:425:ARG:HA	2.15	0.46
1:A:485:LYS:O	1:A:488:ALA:HB2	2.16	0.46
1:B:316:GLU:HG2	1:B:320:LYS:HE2	1.96	0.46
1:B:462:ILE:N	1:B:462:ILE:HD12	2.29	0.46
1:A:246:GLN:HE22	1:A:451:TYR:HA	1.81	0.45
1:B:283:ILE:CD1	1:B:335:PHE:CE1	3.00	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:251:LEU:O	1:B:255:GLU:HG3	2.16	0.45
1:B:181:LEU:HG	1:B:208[B]:CYS:SG	2.55	0.45
1:A:298:HIS:HE1	5:A:3060:HOH:O	2.00	0.45
1:B:283:ILE:HD11	1:B:287:PHE:CE2	2.51	0.45
1:B:273:VAL:O	1:B:274:ASP:HB2	2.17	0.45
1:B:427:HIS:HA	5:B:3003:HOH:O	2.16	0.44
1:B:468:ASP:OD1	1:B:470:ARG:CG	2.65	0.44
1:B:298:HIS:HE1	5:B:3066:HOH:O	2.01	0.43
1:B:468:ASP:C	1:B:470:ARG:N	2.72	0.43
1:B:392:GLN:NE2	5:B:3137:HOH:O	2.51	0.43
1:A:226:LEU:HB3	1:A:367:ARG:HG2	2.01	0.43
1:A:267:ILE:HD12	1:A:298:HIS:CB	2.47	0.42
1:B:445:VAL:CG2	1:B:462:ILE:CD1	2.97	0.42
1:A:449:PRO:O	1:A:450:ASN:C	2.56	0.42
1:B:227:VAL:HG12	1:B:229:THR:HG23	2.00	0.42
1:A:312:ILE:HG13	1:A:313:TYR:CD2	2.54	0.42
1:B:316:GLU:OE2	1:B:320:LYS:HE2	2.20	0.42
1:A:350:ILE:HA	1:A:424:ILE:O	2.18	0.42
1:B:329:GLU:OE1	5:B:3103:HOH:O	2.21	0.42
1:B:422:TYR:HE2	1:B:424:ILE:HD11	1.84	0.42
1:B:312[A]:ILE:HG13	1:B:313:TYR:CD2	2.55	0.42
1:A:267:ILE:CD1	1:A:298:HIS:HB2	2.48	0.41
1:B:485[A]:LYS:CE	1:B:495:GLN:OE1	2.68	0.41
1:A:251:LEU:O	1:A:255:GLU:HG3	2.21	0.41
1:A:237:ILE:HD11	1:A:267:ILE:CD1	2.51	0.41
1:B:232:LYS:HE3	5:B:3140:HOH:O	2.21	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	329/333 (99%)	313 (95%)	16 (5%)	0	100	100
1	B	330/333 (99%)	318 (96%)	10 (3%)	2 (1%)	30	9
All	All	659/666 (99%)	631 (96%)	26 (4%)	2 (0%)	46	23

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	469	LEU
1	B	177	SER

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	294/294 (100%)	290 (99%)	4 (1%)	74	53
1	B	295/294 (100%)	287 (97%)	8 (3%)	52	23
All	All	589/588 (100%)	577 (98%)	12 (2%)	61	36

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	216	VAL
1	A	242	ASP
1	A	290	LYS
1	A	425	ARG
1	B	216	VAL
1	B	222	LYS
1	B	242	ASP
1	B	283	ILE
1	B	425	ARG
1	B	467	SER
1	B	468	ASP
1	B	499	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	205	HIS
1	A	246	GLN
1	A	264	ASN
1	A	326	GLN
1	A	405	GLN
1	A	491	ASN
1	B	205	HIS
1	B	246	GLN
1	B	298	HIS
1	B	392	GLN
1	B	393	ASN
1	B	405	GLN

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 carbohydrates 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	$\# Z > 2$
3	PO4	A	1001	2	4,4,4	1.09	0	6,6,6	0.30	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MPD	A	3001	-	6,7,7	0.39	0	7,10,10	0.70	0
3	PO4	B	2001	2	4,4,4	1.03	0	6,6,6	0.31	0
4	MPD	B	3002	-	6,7,7	0.34	0	7,10,10	0.41	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	PO4	A	1001	2	-	0/0/0/0	0/0/0/0
4	MPD	A	3001	-	-	0/5/5/5	0/0/0/0
3	PO4	B	2001	2	-	0/0/0/0	0/0/0/0
4	MPD	B	3002	-	-	0/5/5/5	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.

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

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	324/333 (97%)	1.34	65 (20%)  	18, 26, 36, 51	7 (2%)
1	B	325/333 (97%)	0.96	51 (15%)  	22, 26, 34, 57	0
All	All	649/666 (97%)	1.15	116 (17%)  	18, 26, 35, 57	7 (1%)

All (116) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	496	LEU	20.0
1	A	498	MET	17.6
1	A	499	MET	14.7
1	A	497	GLY	13.7
1	A	494	LEU	12.8
1	A	176	TYR	11.0
1	A	492	THR	9.8
1	B	177	SER	7.2
1	B	176	TYR	6.6
1	B	498	MET	6.1
1	A	397	ILE	6.1
1	B	470	ARG	6.0
1	A	183	ASP	5.4
1	A	467	SER	5.3
1	B	468	ASP	5.1
1	B	499	MET	5.1
1	A	234	THR	5.0
1	A	493	LEU	5.0
1	A	177	SER	4.9
1	B	300	LEU	4.9
1	A	490	ALA	4.7
1	A	495	GLN	4.5
1	A	300	LEU	4.5
1	B	268	PHE	4.3

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Mol	Chain	Res	Type	RSRZ
1	B	350	ILE	4.2
1	A	282	VAL	4.2
1	B	272	PHE	4.2
1	A	272	PHE	3.8
1	A	283	ILE	3.8
1	A	184	GLY	3.8
1	B	493	LEU	3.8
1	A	377	ASP	3.7
1	A	491	ASN	3.6
1	A	299	LEU	3.6
1	B	494	LEU	3.6
1	B	299	LEU	3.6
1	A	399	LYS	3.6
1	A	268	PHE	3.5
1	B	178	GLY	3.5
1	B	182	GLU	3.5
1	A	182	GLU	3.4
1	A	233[A]	GLU	3.4
1	B	175	GLU	3.4
1	B	286	LEU	3.4
1	B	273	VAL	3.3
1	B	385	LEU	3.3
1	A	286	LEU	3.3
1	B	338	LEU	3.3
1	A	218[A]	GLU	3.2
1	A	338	LEU	3.2
1	A	488	ALA	3.2
1	B	184	GLY	3.1
1	B	250	LEU	3.1
1	B	282	VAL	3.1
1	B	424	ILE	3.0
1	A	385	LEU	3.0
1	A	178	GLY	3.0
1	A	267	ILE	3.0
1	A	181	LEU	3.0
1	A	489	TYR	2.9
1	A	448	ALA	2.9
1	B	340	LEU	2.8
1	B	283	ILE	2.8
1	A	273	VAL	2.8
1	B	243	THR	2.8
1	A	270	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	397	ILE	2.7
1	B	240	CYS	2.6
1	A	454	GLN	2.6
1	B	497	GLY	2.6
1	B	466	GLY	2.6
1	A	487	MET	2.5
1	A	243	THR	2.5
1	A	384	LEU	2.5
1	B	284	LEU	2.5
1	B	384	LEU	2.5
1	A	486	PRO	2.5
1	B	270	GLY	2.5
1	B	253	ILE	2.4
1	A	340	LEU	2.4
1	B	280	VAL	2.4
1	B	238	THR	2.4
1	B	269	ASN	2.4
1	A	424	ILE	2.3
1	A	302	GLY	2.3
1	A	366	ILE	2.3
1	B	247	PHE	2.3
1	B	445	VAL	2.3
1	A	312	ILE	2.3
1	A	455	MET	2.3
1	B	349	LEU	2.2
1	A	398	SER	2.2
1	B	302	GLY	2.2
1	B	352	HIS	2.2
1	A	445	VAL	2.2
1	A	468	ASP	2.2
1	A	232	LYS	2.2
1	A	426	SER	2.2
1	A	297	PHE	2.2
1	A	391	PRO	2.2
1	A	263	THR	2.2
1	A	235	GLU	2.1
1	A	378	SER	2.1
1	A	443	VAL	2.1
1	B	241	GLY	2.1
1	B	287	PHE	2.1
1	A	250	LEU	2.1
1	A	269	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	351	MET	2.0
1	B	341	ALA	2.0
1	B	335	PHE	2.0
1	B	465	GLN	2.0
1	B	239	VAL	2.0
1	B	392	GLN	2.0
1	A	222	LYS	2.0
1	A	301	ARG	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 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
4	MPD	A	3001	8/8	0.83	0.27	6.35	35,37,39,39	0
4	MPD	B	3002	8/8	0.76	0.29	4.27	45,47,47,48	0
3	PO4	A	1001	5/5	0.97	0.23	1.17	27,27,28,29	0
2	MN	A	501	1/1	0.99	0.15	-0.30	24,24,24,24	0
3	PO4	B	2001	5/5	0.98	0.16	-0.50	19,21,21,21	0
2	MN	B	601	1/1	0.99	0.12	-1.77	19,19,19,19	0
2	MN	A	502	1/1	1.00	0.10	-2.81	21,21,21,21	0
2	MN	B	602	1/1	1.00	0.07	-3.88	14,14,14,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.