Full wwPDB X-ray Structure Validation Report (i)

Sep 20, 2023 – 03:56 AM EDT

PDB ID	:	5JQK
Title	:	The Xray Crystal Structure of P. falciparum Aminopeptidase P
Authors	:	Drinkwater, N.; McGowan, S.
Deposited on	:	2016-05-05
Resolution	:	2.35 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:

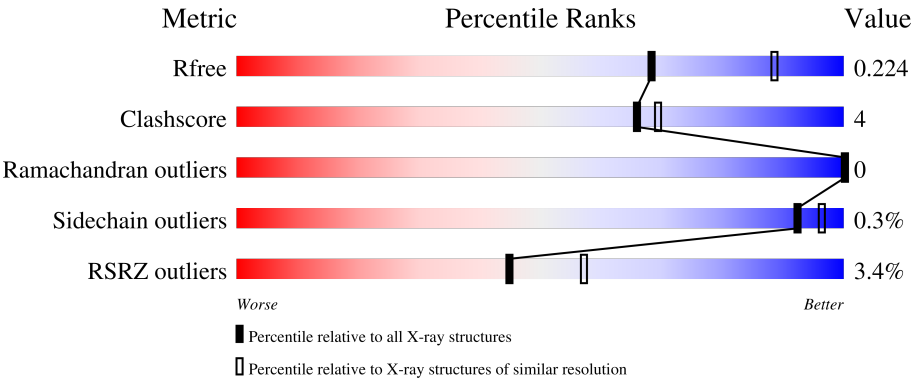
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.35 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	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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	664	2% 86% 9% 5%
1	B	664	5% 84% 9% 7%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 10122 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 Peptidase, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	631	Total	C	N	O	S	0	0	0
			4936	3180	812	928	16			
1	B	615	Total	C	N	O	S	0	0	0
			4661	2999	766	881	15			

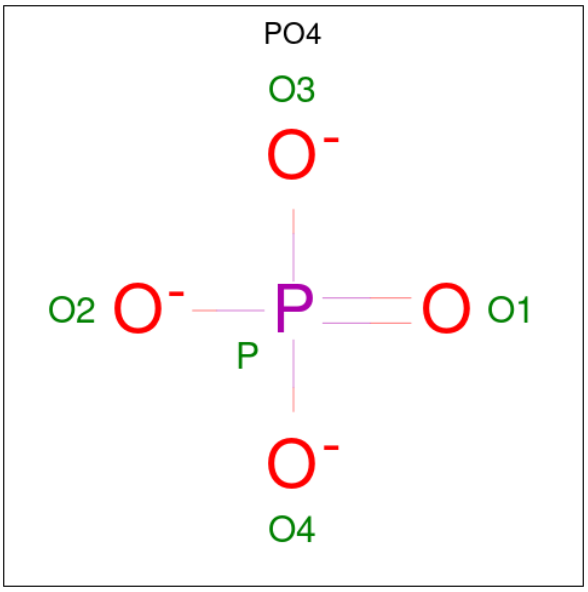
There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	120	MET	-	initiating methionine	UNP Q8IKT5
A	778	HIS	-	expression tag	UNP Q8IKT5
A	779	HIS	-	expression tag	UNP Q8IKT5
A	780	HIS	-	expression tag	UNP Q8IKT5
A	781	HIS	-	expression tag	UNP Q8IKT5
A	782	HIS	-	expression tag	UNP Q8IKT5
A	783	HIS	-	expression tag	UNP Q8IKT5
B	120	MET	-	initiating methionine	UNP Q8IKT5
B	778	HIS	-	expression tag	UNP Q8IKT5
B	779	HIS	-	expression tag	UNP Q8IKT5
B	780	HIS	-	expression tag	UNP Q8IKT5
B	781	HIS	-	expression tag	UNP Q8IKT5
B	782	HIS	-	expression tag	UNP Q8IKT5
B	783	HIS	-	expression tag	UNP Q8IKT5

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

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

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



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

- Molecule 4 is water.

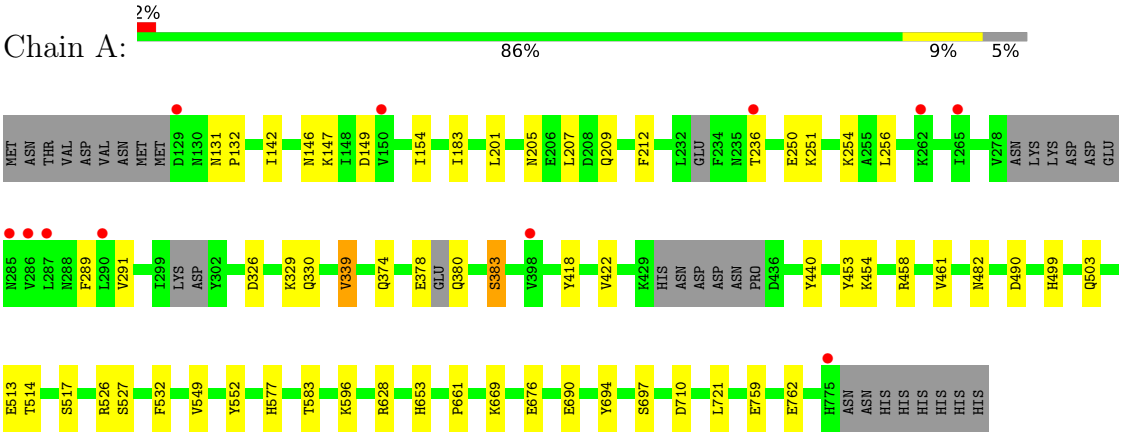
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	284	Total O 284 284	0	0
4	B	227	Total O 227 227	0	0

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.

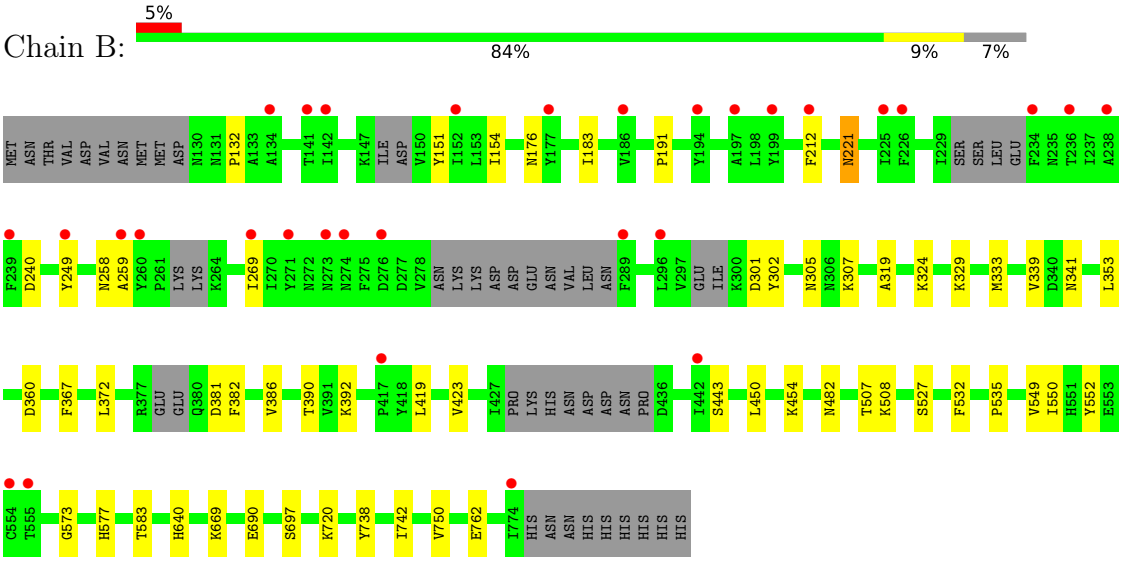
• Molecule 1: Peptidase, putative

Chain A:



• Molecule 1: Peptidase, putative

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	146.74Å 100.07Å 106.66Å 90.00° 105.37° 90.00°	Depositor
Resolution (Å)	44.99 – 2.35 70.75 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.4 (44.99-2.35) 99.4 (70.75-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.69 (at 2.34Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.189 , 0.226 0.190 , 0.224	Depositor DCC
R_{free} test set	3022 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å^2)	39.7	Xtriage
Anisotropy	0.509	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 58.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	10122	wwPDB-VP
Average B, all atoms (Å^2)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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: PO4, MN

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	0.24	0/5033	0.45	0/6831
1	B	0.24	0/4751	0.45	0/6459
All	All	0.24	0/9784	0.45	0/13290

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	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	258	ASN	Peptide
1	B	259	ALA	Peptide
1	B	301	ASP	Peptide
1	B	302	TYR	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	4936	0	4748	42	0
1	B	4661	0	4336	32	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	5	0	0	1	0
3	B	5	0	0	0	0
4	A	284	0	0	4	0
4	B	227	0	0	2	0
All	All	10122	0	9084	74	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 4.

All (74) 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:221:ASN:HD22	1:B:221:ASN:H	1.31	0.77
1:A:329:LYS:NZ	1:A:380:GLN:O	2.21	0.73
1:B:221:ASN:HD22	1:B:221:ASN:N	1.88	0.71
1:A:250:GLU:OE1	1:A:458:ARG:NH2	2.23	0.70
1:A:149:ASP:H	1:A:236:THR:HG23	1.57	0.70
1:A:378:GLU:O	1:A:380:GLN:N	2.25	0.69
1:A:549:VAL:HB	1:A:552:TYR:HB2	1.76	0.68
1:B:669:LYS:HD3	1:B:697:SER:HB2	1.80	0.63
1:B:549:VAL:HB	1:B:552:TYR:HB2	1.81	0.62
1:B:221:ASN:N	1:B:221:ASN:ND2	2.50	0.60
1:B:339:VAL:HG22	1:B:341:ASN:H	1.68	0.59
1:A:207:LEU:O	4:A:901:HOH:O	2.17	0.58
1:A:669:LYS:HD3	1:A:697:SER:HB2	1.85	0.58
1:A:251:LYS:NZ	4:A:910:HOH:O	2.38	0.57
1:A:418:TYR:O	1:A:422:VAL:HG12	2.05	0.56
1:A:453:TYR:CE1	1:A:461:VAL:HG21	2.40	0.56
1:A:513:GLU:OE1	1:A:517:SER:OG	2.19	0.56
1:A:374:GLN:O	1:A:383:SER:N	2.36	0.56
1:B:319:ALA:O	1:B:324:LYS:NZ	2.39	0.55
1:B:419:LEU:HA	1:B:423:VAL:HG22	1.89	0.54
1:B:191:PRO:HB2	1:B:212:PHE:HD1	1.72	0.54
1:A:147:LYS:O	1:A:236:THR:HG21	2.07	0.54
1:A:154:ILE:HG23	1:A:183:ILE:HG12	1.91	0.53
1:A:250:GLU:OE2	1:A:254:LYS:NZ	2.43	0.51
1:A:596:LYS:NZ	4:A:916:HOH:O	2.45	0.50
1:A:205:ASN:ND2	4:A:906:HOH:O	2.35	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:339:VAL:HG11	1:B:443:SER:HB2	1.95	0.49
1:A:453:TYR:HE1	1:A:461:VAL:HG21	1.77	0.48
1:B:381:ASP:OD1	1:B:382:PHE:N	2.42	0.48
1:B:535:PRO:HA	1:B:573:GLY:HA2	1.96	0.48
1:A:142:ILE:O	1:A:146:ASN:ND2	2.46	0.48
1:B:132:PRO:HG3	1:B:176:ASN:ND2	2.28	0.48
1:A:527:SER:HA	1:A:532:PHE:CD2	2.48	0.48
1:B:269:ILE:HD11	1:B:454:LYS:HB3	1.96	0.48
1:A:583:THR:H	1:A:690:GLU:HB3	1.79	0.48
1:A:499:HIS:O	1:A:503:GLN:HG2	2.14	0.47
1:B:151:TYR:OH	1:B:240:ASP:OD1	2.13	0.47
1:B:305:ASN:OD1	1:B:307:LYS:HB2	2.14	0.47
1:B:750:VAL:HG12	1:B:762:GLU:HG2	1.96	0.47
1:A:490:ASP:OD1	1:A:526:ARG:NE	2.44	0.46
1:B:372:LEU:HB3	1:B:386:VAL:HG13	1.97	0.46
1:B:454:LYS:NZ	4:B:909:HOH:O	2.40	0.46
1:A:759:GLU:HA	1:A:762:GLU:HG2	1.98	0.45
1:B:482:ASN:O	1:B:577:HIS:HB3	2.16	0.45
1:B:527:SER:HA	1:B:532:PHE:CD2	2.52	0.45
1:A:596:LYS:HG2	1:A:721:LEU:HA	1.98	0.45
1:B:329:LYS:NZ	1:B:333:MET:SD	2.89	0.45
1:A:256:LEU:HD12	1:A:289:PHE:CZ	2.52	0.44
1:B:550:ILE:HD13	1:B:640:HIS:HD2	1.81	0.44
1:B:507:THR:O	1:B:508:LYS:HB2	2.18	0.43
1:B:738:TYR:CE2	1:B:742:ILE:HD11	2.53	0.43
1:B:324:LYS:HB3	1:B:353:LEU:HD12	1.99	0.43
1:A:418:TYR:CE1	1:A:422:VAL:HG11	2.54	0.43
1:B:720:LYS:NZ	4:B:919:HOH:O	2.51	0.43
1:B:360:ASP:HB2	1:B:367:PHE:HA	2.01	0.43
1:A:628:ARG:HH21	1:A:661:PRO:HD3	1.84	0.43
1:A:514:THR:H	1:A:517:SER:HB3	1.83	0.42
1:B:583:THR:H	1:B:690:GLU:HB3	1.83	0.42
1:A:482:ASN:O	1:A:577:HIS:HB3	2.19	0.42
1:B:154:ILE:HG23	1:B:183:ILE:HG12	2.01	0.42
1:B:390:THR:HG22	1:B:392:LYS:HG2	2.01	0.42
1:A:458:ARG:O	1:A:461:VAL:HG22	2.19	0.42
1:A:291:VAL:HG21	1:A:454:LYS:HG3	2.03	0.41
1:A:256:LEU:HD12	1:A:289:PHE:HZ	1.85	0.41
1:B:249:TYR:CG	1:B:450:LEU:HD21	2.55	0.41
1:A:131:ASN:HA	1:A:132:PRO:HD3	1.77	0.41
1:A:339:VAL:CG1	1:A:440:TYR:HB2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:694:TYR:CE2	1:A:710:ASP:HB3	2.55	0.41
1:A:201:LEU:HD23	1:A:201:LEU:HA	1.92	0.40
1:A:209:GLN:HA	1:A:212:PHE:O	2.21	0.40
1:A:326:ASP:O	1:A:330:GLN:HG3	2.21	0.40
1:A:653:HIS:NE2	3:A:803:PO4:O3	2.44	0.40
1:A:583:THR:N	1:A:690:GLU:HB3	2.36	0.40
1:A:676:GLU:HB3	1:A:690:GLU:HB2	2.04	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	619/664 (93%)	612 (99%)	7 (1%)	0	100	100
1	B	599/664 (90%)	590 (98%)	9 (2%)	0	100	100
All	All	1218/1328 (92%)	1202 (99%)	16 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	516/613 (84%)	514 (100%)	2 (0%)	91	95

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	B	463/613 (76%)	462 (100%)	1 (0%)	93 97
All	All	979/1226 (80%)	976 (100%)	3 (0%)	92 96

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	339	VAL
1	A	383	SER
1	B	221	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	165	ASN
1	B	221	ASN

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 monosaccharides 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	PO4	A	803	2	4,4,4	0.93	0	6,6,6	0.43	0
3	PO4	B	803	2	4,4,4	0.91	0	6,6,6	0.54	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	803	PO4	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	631/664 (95%)	0.08	11 (1%)	70	78	26, 55, 104, 159	1 (0%)
1	B	615/664 (92%)	0.37	31 (5%)	28	41	27, 65, 132, 166	1 (0%)
All	All	1246/1328 (93%)	0.22	42 (3%)	45	57	26, 59, 117, 166	2 (0%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	775	HIS	7.8
1	A	286	VAL	4.9
1	B	186	VAL	4.6
1	B	141	THR	4.3
1	A	285	ASN	3.8
1	B	554	CYS	3.6
1	B	249	TYR	3.3
1	B	260	TYR	3.3
1	B	142	ILE	3.3
1	A	150	VAL	3.3
1	B	152	ILE	3.2
1	B	555	THR	3.2
1	B	274	ASN	3.2
1	B	236	THR	3.1
1	B	269	ILE	3.1
1	B	212	PHE	3.1
1	B	296	LEU	3.0
1	B	197	ALA	2.9
1	A	265	ILE	2.8
1	B	259	ALA	2.6
1	B	442	ILE	2.5
1	B	134	ALA	2.5
1	B	238	ALA	2.4
1	A	129	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	289	PHE	2.4
1	A	236	THR	2.4
1	A	398	VAL	2.4
1	B	417	PRO	2.4
1	B	194	TYR	2.4
1	B	774	ILE	2.3
1	B	276	ASP	2.3
1	A	290	LEU	2.3
1	B	226	PHE	2.3
1	B	271	TYR	2.2
1	B	177	TYR	2.1
1	B	273	ASN	2.1
1	A	287	LEU	2.1
1	B	199	TYR	2.1
1	B	225	ILE	2.1
1	A	262	LYS	2.0
1	B	234	PHE	2.0
1	B	239	PHE	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 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	PO4	A	803	5/5	0.97	0.15	31,39,43,47	0
3	PO4	B	803	5/5	0.98	0.14	49,51,58,59	0
2	MN	B	802	1/1	0.99	0.17	37,37,37,37	0
2	MN	A	801	1/1	0.99	0.17	30,30,30,30	0
2	MN	A	802	1/1	0.99	0.17	28,28,28,28	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	MN	B	801	1/1	1.00	0.13	40,40,40,40	0

6.5 Other polymers (i)

There are no such residues in this entry.