



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:27 PM GMT

PDB ID : 4F0F
Title : Crystal Structure of the Roco4 Kinase Domain bound to AppCp from D. discoideum
Authors : Gilsbach, B.K.; Vetter, I.R.; Wittinghofer, A.; Kortholt, A.
Deposited on : 2012-05-04
Resolution : 1.80 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

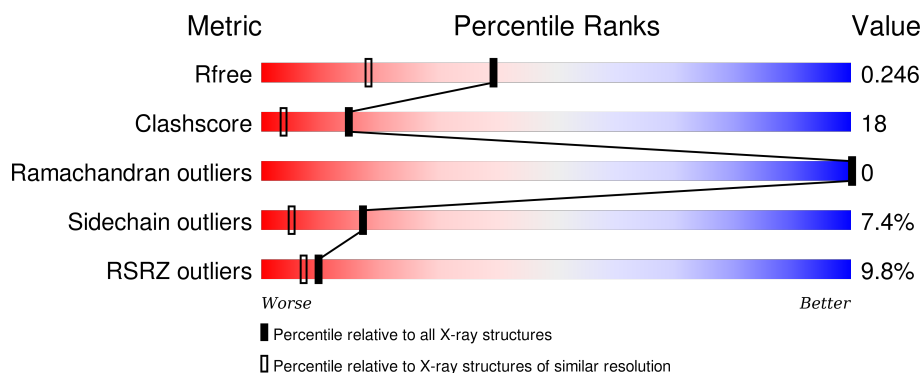
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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	287	<div> <div>9%</div> <div>66%</div> <div>24%</div> <div>5% . .</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2456 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.

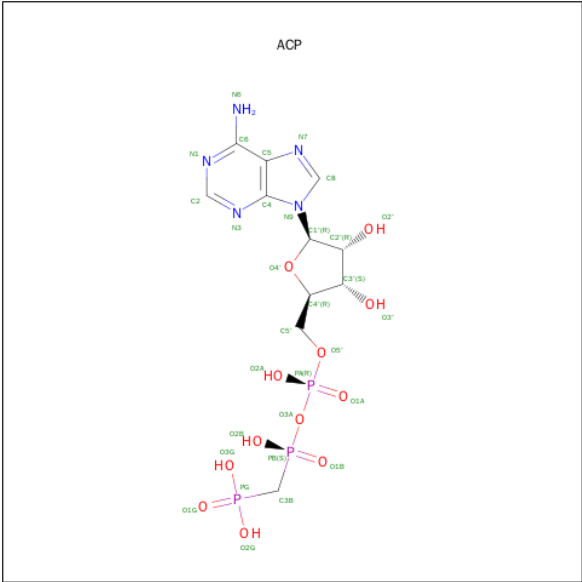
- Molecule 1 is a protein called Serine/threonine-protein kinase roco4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	1	0
			2214	1420	371	409	14			

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1006	GLY	-	EXPRESSION TAG	UNP Q6XHB2
A	1007	ALA	-	EXPRESSION TAG	UNP Q6XHB2
A	1008	MET	-	EXPRESSION TAG	UNP Q6XHB2
A	1009	GLY	-	EXPRESSION TAG	UNP Q6XHB2
A	1010	GLY	-	EXPRESSION TAG	UNP Q6XHB2
A	1011	SER	-	EXPRESSION TAG	UNP Q6XHB2
A	1012	GLU	-	EXPRESSION TAG	UNP Q6XHB2
A	1013	PHE	-	EXPRESSION TAG	UNP Q6XHB2
A	1014	PRO	-	EXPRESSION TAG	UNP Q6XHB2
A	1015	LYS	-	EXPRESSION TAG	UNP Q6XHB2
A	1016	SER	-	EXPRESSION TAG	UNP Q6XHB2
A	1017	ARG	-	EXPRESSION TAG	UNP Q6XHB2
A	1018	LEU	-	EXPRESSION TAG	UNP Q6XHB2

- Molecule 2 is PHOSPHOMETHYLPHOSPHONIC ACID ADENYLATE ESTER (three-letter code: ACP) (formula: $C_{11}H_{18}N_5O_{12}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	11	5	12	3		

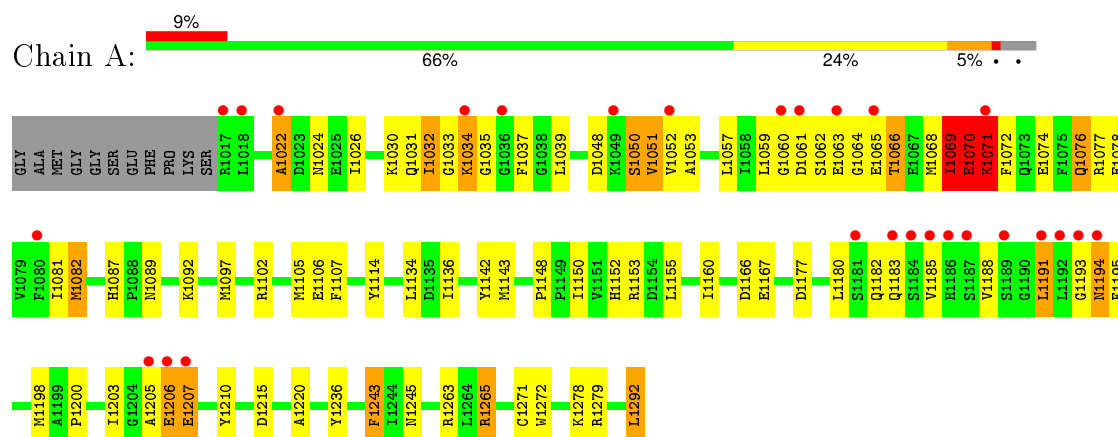
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	211	Total	O	0	0
			211	211		

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 ($\text{RSRZ} > 2$). 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 kinase roco4



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	42.72Å 42.72Å 339.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.10 – 1.80 19.10 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (19.10-1.80) 100.0 (19.10-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.03 (at 1.80Å)	Xtriage
Refinement program	REFMAC 5.6.0093	Depositor
R, R_{free}	0.195 , 0.245 0.198 , 0.246	Depositor DCC
R_{free} test set	1539 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	21.4	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 41.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 30792 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2456	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.52% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ACP

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	1.24	5/2270 (0.2%)	1.32	24/3067 (0.8%)

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	A	0	4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1236	TYR	CD1-CE1	7.20	1.50	1.39
1	A	1265	ARG	CD-NE	-6.17	1.35	1.46
1	A	1243	PHE	N-CA	5.62	1.57	1.46
1	A	1114	TYR	CD1-CE1	5.52	1.47	1.39
1	A	1142	TYR	CD1-CE1	5.25	1.47	1.39

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1265	ARG	NE-CZ-NH2	-19.78	110.41	120.30
1	A	1265	ARG	NE-CZ-NH1	14.35	127.47	120.30
1	A	1022	ALA	CB-CA-C	13.40	130.20	110.10
1	A	1195	PHE	N-CA-CB	11.42	131.16	110.60
1	A	1069	ILE	CG1-CB-CG2	-8.97	91.67	111.40
1	A	1024	ASN	N-CA-CB	-8.93	94.53	110.60
1	A	1193	GLY	N-CA-C	8.25	133.73	113.10
1	A	1071	LYS	CB-CA-C	-8.12	94.16	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1177	ASP	CB-CG-OD1	8.03	125.52	118.30
1	A	1053	ALA	N-CA-CB	-7.70	99.32	110.10
1	A	1053	ALA	N-CA-C	-7.27	91.37	111.00
1	A	1292	LEU	CB-CG-CD1	7.17	123.19	111.00
1	A	1050	SER	N-CA-CB	-6.91	100.13	110.50
1	A	1033	GLY	N-CA-C	6.48	129.29	113.10
1	A	1166	ASP	CB-CG-OD2	6.08	123.77	118.30
1	A	1215	ASP	CB-CG-OD1	5.71	123.44	118.30
1	A	1265	ARG	CG-CD-NE	-5.58	100.08	111.80
1	A	1195	PHE	N-CA-C	-5.40	96.42	111.00
1	A	1265	ARG	CD-NE-CZ	5.37	131.12	123.60
1	A	1134	LEU	CB-CG-CD1	-5.33	101.93	111.00
1	A	1105	MET	CA-CB-CG	5.30	122.32	113.30
1	A	1082	MET	CA-CB-CG	5.25	122.22	113.30
1	A	1279	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	A	1071	LYS	N-CA-C	5.08	124.73	111.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1022	ALA	Peptide
1	A	1051	VAL	Peptide
1	A	1070	GLU	Peptide
1	A	1265	ARG	Sidechain

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	2214	0	2208	77	4
2	A	31	0	14	4	0
3	A	211	0	0	27	1
All	All	2456	0	2222	79	4

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 18.

All (79) 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:1263:ARG:HD3	3:A:1603:HOH:O	1.53	1.07
1:A:1071:LYS:O	1:A:1071:LYS:HG3	1.26	1.03
1:A:1097:MET:CE	1:A:1102:ARG:HD2	1.88	1.03
1:A:1097:MET:HE3	1:A:1102:ARG:HD2	1.38	0.99
1:A:1263:ARG:CD	3:A:1603:HOH:O	2.11	0.98
1:A:1167:GLU:OE1	3:A:1610:HOH:O	1.85	0.95
1:A:1194:ASN:O	3:A:1541:HOH:O	1.87	0.93
1:A:1076:GLN:HG2	3:A:1585:HOH:O	1.68	0.93
1:A:1278:LYS:HE2	3:A:1562:HOH:O	1.73	0.89
1:A:1097:MET:HE3	1:A:1102:ARG:CD	2.03	0.87
1:A:1071:LYS:O	1:A:1071:LYS:CG	2.16	0.87
1:A:1032:ILE:HG22	3:A:1552:HOH:O	1.78	0.83
1:A:1153:ARG:HD2	1:A:1182:GLN:HG2	1.61	0.82
2:A:1301:ACP:H3B2	3:A:1542:HOH:O	1.81	0.81
2:A:1301:ACP:H3B1	3:A:1555:HOH:O	1.84	0.77
1:A:1087:HIS:HD2	1:A:1089:ASN:H	1.32	0.77
1:A:1198:MET:HG3	3:A:1541:HOH:O	1.88	0.74
1:A:1087:HIS:CD2	1:A:1089:ASN:H	2.05	0.73
1:A:1153:ARG:HD3	1:A:1180:LEU:O	1.88	0.72
1:A:1188:VAL:HB	3:A:1578:HOH:O	1.90	0.72
1:A:1245:ASN:HB2	3:A:1498:HOH:O	1.88	0.72
1:A:1210:TYR:CD2	3:A:1578:HOH:O	2.44	0.70
1:A:1245:ASN:CB	3:A:1498:HOH:O	2.39	0.70
1:A:1070:GLU:HA	1:A:1072:PHE:H	1.55	0.70
1:A:1082:MET:HA	1:A:1082:MET:HE3	1.74	0.69
1:A:1210:TYR:CE2	3:A:1578:HOH:O	2.45	0.69
1:A:1263:ARG:NE	3:A:1603:HOH:O	2.22	0.69
1:A:1057:LEU:HD21	1:A:1072:PHE:HD1	1.59	0.68
1:A:1048:ASP:HB3	1:A:1050:SER:HB2	1.77	0.65
1:A:1182:GLN:HG3	1:A:1188:VAL:CG1	2.31	0.60
1:A:1071:LYS:HA	3:A:1573:HOH:O	1.97	0.60
1:A:1051:VAL:HG12	1:A:1107:PHE:CB	2.30	0.60
1:A:1077:ARG:HD2	3:A:1508:HOH:O	2.02	0.60
1:A:1097:MET:HE3	1:A:1102:ARG:NE	2.17	0.59
1:A:1087:HIS:HE1	3:A:1425:HOH:O	1.87	0.57
1:A:1071:LYS:N	3:A:1573:HOH:O	1.62	0.57
1:A:1051:VAL:HG12	1:A:1107:PHE:HB3	1.84	0.57
1:A:1070:GLU:HA	1:A:1071:LYS:HB3	1.86	0.57
1:A:1065:GLU:HB3	1:A:1068:MET:SD	2.45	0.57
1:A:1148:PRO:HG2	1:A:1183:GLN:HG3	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1070:GLU:CA	1:A:1071:LYS:HB3	2.33	0.57
1:A:1048:ASP:CB	1:A:1050:SER:HB2	2.35	0.57
1:A:1191:LEU:HG	3:A:1489:HOH:O	2.05	0.56
1:A:1071:LYS:CA	3:A:1573:HOH:O	2.30	0.56
1:A:1070:GLU:HA	1:A:1072:PHE:N	2.19	0.55
1:A:1150:ILE:HD13	1:A:1183:GLN:HB2	1.87	0.55
1:A:1034:LYS:O	2:A:1301:ACP:H5'2	2.07	0.55
1:A:1031:GLN:HG2	1:A:1032:ILE:N	2.21	0.54
1:A:1026:ILE:HG13	1:A:1026:ILE:O	2.07	0.54
1:A:1191:LEU:CG	3:A:1489:HOH:O	2.56	0.53
1:A:1097:MET:HB2	1:A:1102:ARG:HB2	1.91	0.53
1:A:1057:LEU:HD21	1:A:1072:PHE:CD1	2.43	0.52
1:A:1037:PHE:HB3	1:A:1071:LYS:HD2	1.91	0.52
1:A:1060:GLY:O	1:A:1061:ASP:C	2.46	0.51
1:A:1182:GLN:HG3	1:A:1188:VAL:HG11	1.92	0.50
1:A:1092:LYS:H	1:A:1106:GLU:HG2	1.77	0.49
1:A:1097:MET:HE3	1:A:1102:ARG:CZ	2.43	0.49
1:A:1092:LYS:HD2	3:A:1597:HOH:O	2.14	0.48
1:A:1078:GLU:O	1:A:1082:MET:HG2	2.14	0.47
1:A:1077:ARG:O	1:A:1081:ILE:HG12	2.15	0.46
1:A:1185:VAL:HG23	3:A:1594:HOH:O	2.14	0.46
1:A:1035:GLY:HA2	2:A:1301:ACP:O2A	2.16	0.45
1:A:1220:ALA:HB2	1:A:1271:CYS:HB3	1.99	0.45
1:A:1153:ARG:HA	3:A:1476:HOH:O	2.16	0.45
1:A:1220:ALA:HB2	1:A:1271:CYS:CB	2.46	0.45
1:A:1062:SER:O	1:A:1063:GLU:C	2.54	0.45
1:A:1069:ILE:HG21	1:A:1069:ILE:HD13	1.24	0.44
1:A:1205:ALA:HB1	1:A:1207:GLU:HG3	1.48	0.43
1:A:1191:LEU:HD12	3:A:1489:HOH:O	2.17	0.43
1:A:1136:ILE:HD11	1:A:1160:ILE:HG21	1.99	0.43
1:A:1206:GLU:OE2	1:A:1206:GLU:O	2.36	0.43
1:A:1155:LEU:HA	1:A:1155:LEU:HD23	1.88	0.42
1:A:1148:PRO:HG2	1:A:1183:GLN:CG	2.49	0.42
1:A:1069:ILE:C	1:A:1071:LYS:HB3	2.40	0.42
1:A:1057:LEU:CD2	1:A:1072:PHE:CD1	3.02	0.41
1:A:1143:MET:HE1	1:A:1152:HIS:HB2	2.02	0.41
1:A:1037:PHE:HE1	1:A:1074:GLU:HG2	1.85	0.41
1:A:1057:LEU:CD2	1:A:1072:PHE:HD1	2.30	0.40
1:A:1200:PRO:HA	1:A:1203:ILE:HG12	2.04	0.40

All (4) 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)
1:A:1061:ASP:OD2	1:A:1064:GLY:O[7_465]	1.82	0.38
1:A:1066:THR:OG1	3:A:1522:HOH:O[7_465]	1.95	0.25
1:A:1072:PHE:CD2	1:A:1072:PHE:CE2[7_465]	2.12	0.08
1:A:1072:PHE:CD2	1:A:1072:PHE:CD2[7_465]	2.18	0.02

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	275/287 (96%)	265 (96%)	10 (4%)	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	245/253 (97%)	227 (93%)	18 (7%)	17	5

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1030	LYS
1	A	1032	ILE
1	A	1034	LYS
1	A	1039	LEU
1	A	1052	VAL

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Mol	Chain	Res	Type
1	A	1059	LEU
1	A	1066	THR
1	A	1069	ILE
1	A	1070	GLU
1	A	1071	LYS
1	A	1076	GLN
1	A	1191	LEU
1	A	1194	ASN
1	A	1206	GLU
1	A	1207	GLU
1	A	1243	PHE
1	A	1272	TRP
1	A	1292	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1087	HIS
1	A	1115	HIS
1	A	1194	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	$\# Z > 2$
2	ACP	A	1301	-	25,33,33	2.41	5 (20%)	31,52,52	2.20	6 (19%)

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
2	ACP	A	1301	-	-	0/15/38/38	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1301	ACP	PB-O2B	-3.10	1.48	1.56
2	A	1301	ACP	O4'-C1'	4.00	1.46	1.41
2	A	1301	ACP	PG-O3G	4.03	1.64	1.54
2	A	1301	ACP	PG-O2G	4.33	1.65	1.54
2	A	1301	ACP	PB-O3A	8.41	1.67	1.58

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1301	ACP	N3-C2-N1	-9.11	121.92	128.89
2	A	1301	ACP	C1'-N9-C4	-5.04	119.34	126.94
2	A	1301	ACP	C5'-C4'-C3'	-2.52	105.21	115.21
2	A	1301	ACP	PA-O3A-PB	-2.46	125.81	132.73
2	A	1301	ACP	C4-C5-N7	-2.15	107.50	109.48
2	A	1301	ACP	O2B-PB-O1B	3.33	120.60	110.12

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1301	ACP	4	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	276/287 (96%)	0.35	27 (9%) 10 7	11, 24, 64, 130	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1061	ASP	6.8
1	A	1186	HIS	6.6
1	A	1193	GLY	6.4
1	A	1022	ALA	6.1
1	A	1191	LEU	5.6
1	A	1187	SER	5.5
1	A	1206	GLU	4.7
1	A	1192	LEU	4.6
1	A	1194	ASN	4.5
1	A	1205	ALA	4.4
1	A	1034	LYS	4.2
1	A	1207	GLU	3.9
1	A	1185	VAL	3.8
1	A	1065	GLU	3.7
1	A	1052	VAL	3.6
1	A	1184	SER	3.5
1	A	1060	GLY	3.2
1	A	1063	GLU	2.9
1	A	1183	GLN	2.7
1	A	1017	ARG	2.7
1	A	1018	LEU	2.5
1	A	1189	SER	2.5
1	A	1071	LYS	2.4
1	A	1080	PHE	2.3
1	A	1181	SER	2.2
1	A	1036	GLY	2.1
1	A	1049	LYS	2.1

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(\AA^2)	Q<0.9
2	ACP	A	1301	31/31	0.82	0.19	0.61	22,46,97,109	0

6.5 Other polymers [i](#)

There are no such residues in this entry.