



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 10:15 PM GMT

PDB ID : 5CJV
Title : Isobutyryl-CoA mutase fused with bound adenosylcobalamin, GDP, Mg (holo-IcmF/GDP), and substrate isovaleryl-coenzyme A
Authors : Jost, M.; Drennan, C.L.
Deposited on : 2015-07-15
Resolution : 3.45 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-20026982
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-20026982

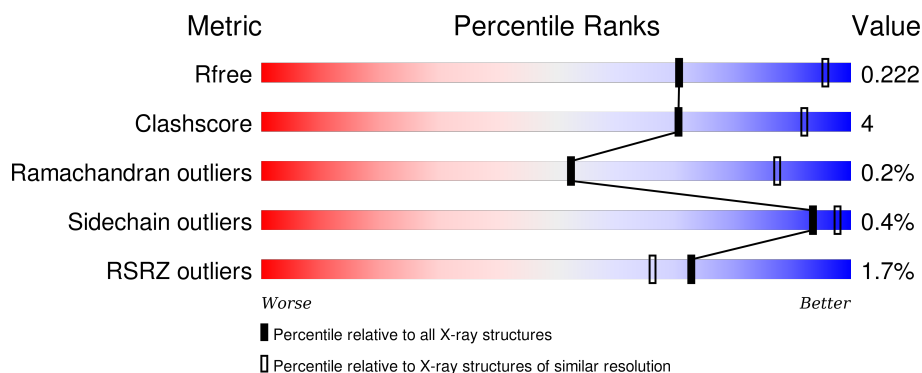
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.45 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	1000 (3.56-3.36)
Clashscore	102246	1090 (3.56-3.36)
Ramachandran outliers	100387	1057 (3.56-3.36)
Sidechain outliers	100360	1058 (3.56-3.36)
RSRZ outliers	91569	1005 (3.56-3.36)

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 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	1113	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -10px; left: 0; width: 100%; height: 1px; background-color: red;"></div> <div style="position: absolute; bottom: 0; left: 0; width: 100%; height: 10px; background-color: green;"></div> <div style="position: absolute; bottom: 0; left: 84%; width: 11%; height: 10px; background-color: yellow;"></div> <div style="position: absolute; bottom: 0; left: 95%; width: 5%; height: 10px; background-color: grey;"></div> </div> <div>84% 11% 5%</div> </div>
1	B	1113	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -10px; left: 0; width: 100%; height: 1px; background-color: red;"></div> <div style="position: absolute; bottom: 0; left: 0; width: 100%; height: 10px; background-color: green;"></div> <div style="position: absolute; bottom: 0; left: 85%; width: 10%; height: 10px; background-color: yellow;"></div> <div style="position: absolute; bottom: 0; left: 95%; width: 5%; height: 10px; background-color: grey;"></div> </div> <div>85% 10% 5%</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
6	MG	A	1106	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 16519 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 Isobutyryl-CoA mutase fused.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1061	Total	C	N	O	S	0	0	0
			8093	5058	1449	1549	37			
1	B	1059	Total	C	N	O	S	0	0	0
			8067	5034	1453	1543	37			

There are 40 discrepancies between the modelled and reference sequences:

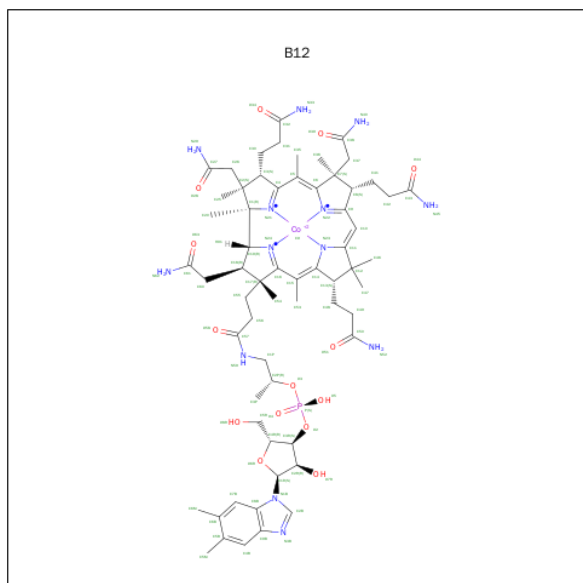
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP Q1LRY0
A	-18	GLY	-	expression tag	UNP Q1LRY0
A	-17	SER	-	expression tag	UNP Q1LRY0
A	-16	SER	-	expression tag	UNP Q1LRY0
A	-15	HIS	-	expression tag	UNP Q1LRY0
A	-14	HIS	-	expression tag	UNP Q1LRY0
A	-13	HIS	-	expression tag	UNP Q1LRY0
A	-12	HIS	-	expression tag	UNP Q1LRY0
A	-11	HIS	-	expression tag	UNP Q1LRY0
A	-10	HIS	-	expression tag	UNP Q1LRY0
A	-9	SER	-	expression tag	UNP Q1LRY0
A	-8	SER	-	expression tag	UNP Q1LRY0
A	-7	GLY	-	expression tag	UNP Q1LRY0
A	-6	LEU	-	expression tag	UNP Q1LRY0
A	-5	VAL	-	expression tag	UNP Q1LRY0
A	-4	PRO	-	expression tag	UNP Q1LRY0
A	-3	ARG	-	expression tag	UNP Q1LRY0
A	-2	GLY	-	expression tag	UNP Q1LRY0
A	-1	SER	-	expression tag	UNP Q1LRY0
A	0	HIS	-	expression tag	UNP Q1LRY0
B	-19	MET	-	initiating methionine	UNP Q1LRY0
B	-18	GLY	-	expression tag	UNP Q1LRY0
B	-17	SER	-	expression tag	UNP Q1LRY0
B	-16	SER	-	expression tag	UNP Q1LRY0
B	-15	HIS	-	expression tag	UNP Q1LRY0

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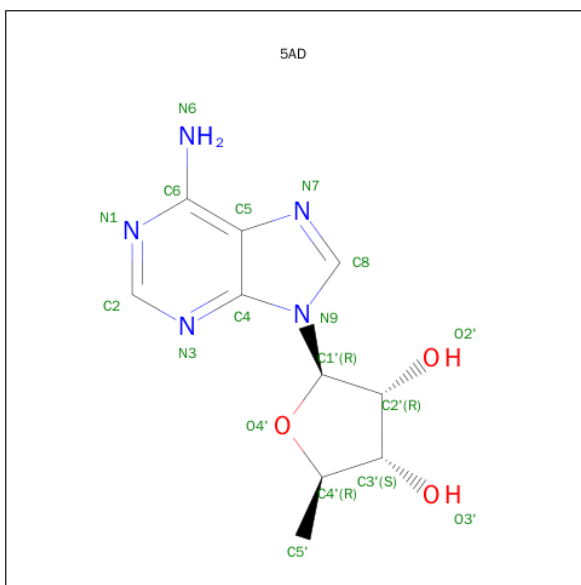
Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP Q1LRY0
B	-13	HIS	-	expression tag	UNP Q1LRY0
B	-12	HIS	-	expression tag	UNP Q1LRY0
B	-11	HIS	-	expression tag	UNP Q1LRY0
B	-10	HIS	-	expression tag	UNP Q1LRY0
B	-9	SER	-	expression tag	UNP Q1LRY0
B	-8	SER	-	expression tag	UNP Q1LRY0
B	-7	GLY	-	expression tag	UNP Q1LRY0
B	-6	LEU	-	expression tag	UNP Q1LRY0
B	-5	VAL	-	expression tag	UNP Q1LRY0
B	-4	PRO	-	expression tag	UNP Q1LRY0
B	-3	ARG	-	expression tag	UNP Q1LRY0
B	-2	GLY	-	expression tag	UNP Q1LRY0
B	-1	SER	-	expression tag	UNP Q1LRY0
B	0	HIS	-	expression tag	UNP Q1LRY0

- Molecule 2 is COBALAMIN (three-letter code: B12) (formula: $C_{62}H_{89}CoN_{13}O_{14}P$).



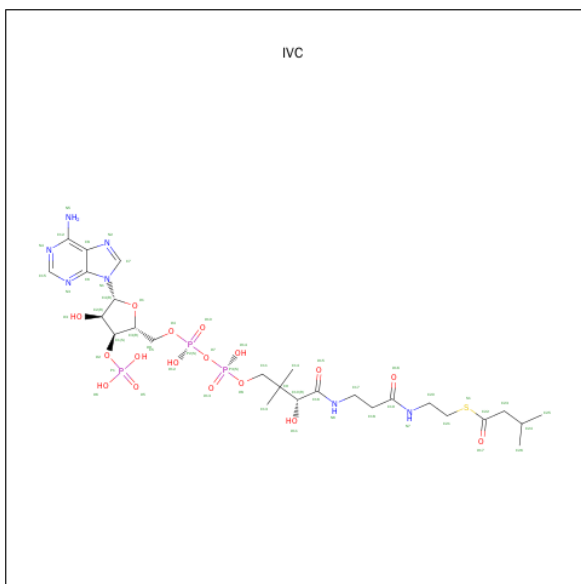
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	Co	N	O	P	0	0
			91	62	1	13	14	1		
2	B	1	Total	C	Co	N	O	P	0	0
			91	62	1	13	14	1		

- Molecule 3 is 5'-DEOXYADENOSINE (three-letter code: 5AD) (formula: $C_{10}H_{13}N_5O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	1
			36	20	10	6		

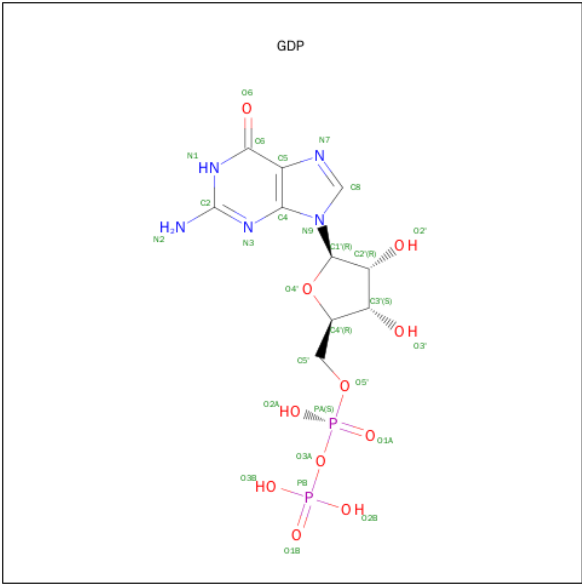
- Molecule 4 is Isovaleryl-coenzyme A (three-letter code: IVC) (formula: C₂₆H₄₄N₇O₁₇P₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	S	0	0
			54	26	7	17	3	1		
4	B	1	Total	C	N	O	P		0	0
			27	10	5	10	2			

- Molecule 5 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula:

C₁₀H₁₅N₅O₁₁P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			28	10	5	11	2		
5	B	1	Total	C	N	O	P	0	0
			28	10	5	11	2		

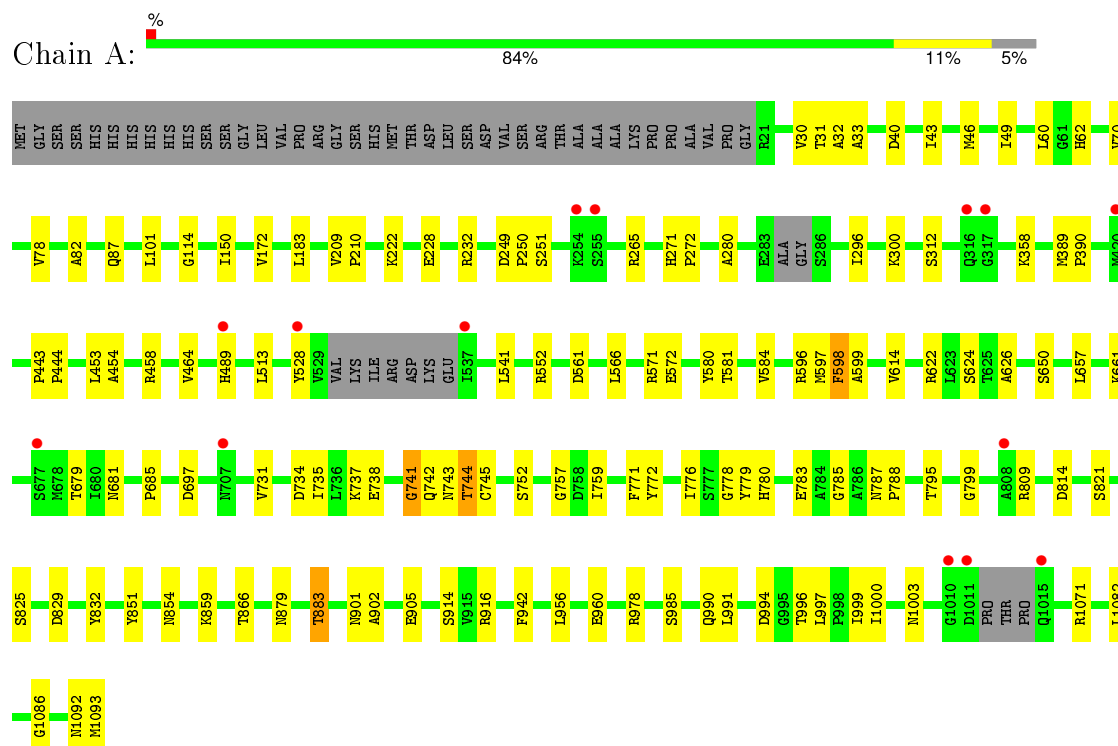
- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	2	Total	Mg	0	0
			2	2		
6	A	2	Total	Mg	0	0
			2	2		

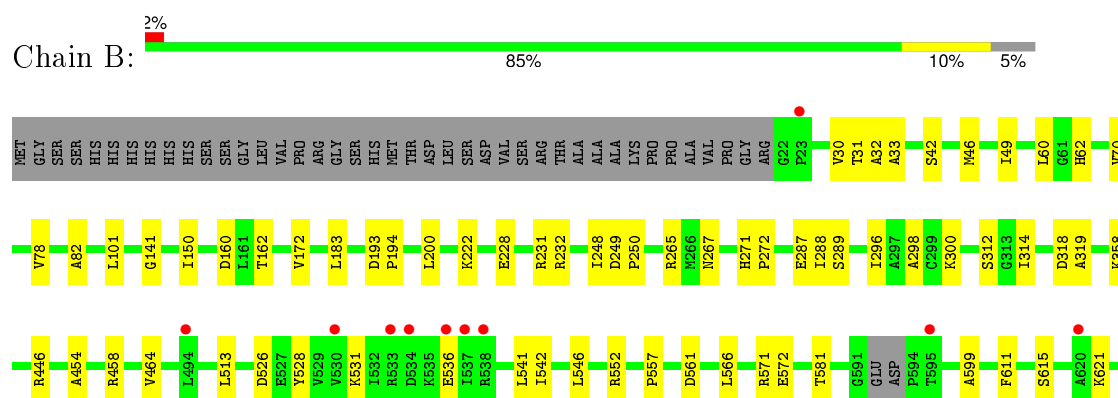
3 Residue-property plots

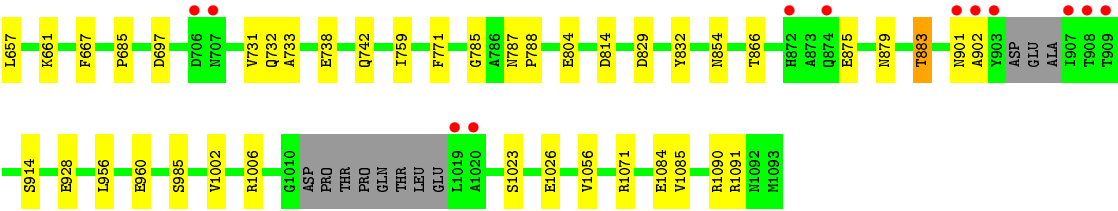
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Isobutyryl-CoA mutase fused



• Molecule 1: Isobutyryl-CoA mutase fused





4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	317.56Å 317.56Å 343.52Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.91 – 3.45 34.91 – 3.45	Depositor EDS
% Data completeness (in resolution range)	99.0 (34.91-3.45) 99.1 (34.91-3.45)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 3.47Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.193 , 0.214 0.201 , 0.222	Depositor DCC
R_{free} test set	4357 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	89.2	Xtriage
Anisotropy	0.274	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 54.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 86230 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	16519	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.08% 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: GDP, MG, B12, 5AD, IVC

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	0.21	0/8234	0.40	3/11154 (0.0%)
1	B	0.21	0/8205	0.39	3/11107 (0.0%)
All	All	0.21	0/16439	0.39	6/22261 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	552	ARG	NE-CZ-NH1	10.07	125.33	120.30
1	A	552	ARG	NE-CZ-NH2	-9.65	115.48	120.30
1	B	552	ARG	NE-CZ-NH2	9.57	125.08	120.30
1	B	552	ARG	NE-CZ-NH1	-9.56	115.52	120.30
1	B	552	ARG	CD-NE-CZ	5.34	131.07	123.60
1	A	552	ARG	CD-NE-CZ	5.21	130.89	123.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	8093	0	7834	81	0
1	B	8067	0	7806	59	0
2	A	91	0	88	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	91	0	88	6	0
3	A	36	0	26	3	0
4	A	54	0	42	7	0
4	B	27	0	13	0	0
5	A	28	0	12	3	0
5	B	28	0	12	3	0
6	A	2	0	0	0	0
6	B	2	0	0	0	0
All	All	16519	0	15921	144	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 (144) 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:87:GLN:HE22	2:A:1101:B12:H421	1.44	0.82
1:A:783:GLU:HB3	2:A:1101:B12:H532	1.64	0.78
1:B:615:SER:O	1:B:621:LYS:NZ	2.25	0.69
1:A:779:TYR:OH	3:A:1102[B]:5AD:O3'	2.13	0.66
1:A:821:SER:HG	4:A:1103:IVC:H25	1.39	0.66
1:B:314:ILE:HG23	1:B:318:ASP:HB2	1.81	0.62
1:A:734:ASP:HB3	1:A:737:LYS:HZ2	1.65	0.62
1:A:681:ASN:ND2	1:A:742:GLN:OE1	2.32	0.62
1:A:31:THR:HG22	1:A:82:ALA:HB3	1.82	0.61
1:A:622:ARG:NH1	4:A:1103:IVC:O13	2.24	0.60
1:A:991:LEU:HD22	1:A:996:THR:HB	1.84	0.59
1:B:454:ALA:HA	1:B:956:LEU:HD22	1.84	0.59
1:B:31:THR:HG22	1:B:82:ALA:HB3	1.83	0.59
1:A:735:ILE:HG13	1:A:752:SER:HB3	1.85	0.58
1:A:990:GLN:O	1:A:994:ASP:HB2	2.02	0.58
1:A:780:HIS:HA	2:A:1101:B12:H461	1.86	0.57
1:A:1082:LEU:O	1:A:1086:GLY:N	2.37	0.57
1:A:33:ALA:HB3	1:A:62:HIS:HA	1.87	0.56
1:A:454:ALA:HA	1:A:956:LEU:HD22	1.88	0.55
1:B:33:ALA:HB3	1:B:62:HIS:HA	1.88	0.55
1:B:1002:VAL:O	1:B:1006:ARG:NH1	2.36	0.55
1:A:561:ASP:HB2	1:B:561:ASP:HB2	1.89	0.54
1:A:785:GLY:HA3	1:A:985:SER:HB2	1.90	0.53
1:A:942:PHE:CZ	1:B:557:PRO:HB3	2.43	0.53
1:A:596:ARG:HG2	1:A:624:SER:HB2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:742:GLN:HG3	2:A:1101:B12:H472	1.91	0.53
1:A:679:THR:HG21	4:A:1103:IVC:H40	1.91	0.52
1:A:251:SER:HB2	1:A:280:ALA:HB1	1.92	0.52
1:B:738:GLU:OE2	1:B:742:GLN:NE2	2.41	0.52
1:A:228:GLU:OE1	1:A:232:ARG:NH1	2.43	0.52
1:B:572:GLU:HG2	1:B:581:THR:HG21	1.92	0.52
1:A:458:ARG:NH1	1:A:960:GLU:OE2	2.41	0.52
1:A:879:ASN:O	1:A:883:THR:OG1	2.26	0.51
1:A:265:ARG:NH1	5:A:1104:GDP:O1A	2.44	0.51
3:A:1102[A]:5AD:H5'1	4:A:1103:IVC:H43	1.92	0.51
1:B:785:GLY:HA3	1:B:985:SER:HB2	1.92	0.51
1:A:572:GLU:HG2	1:A:581:THR:HG21	1.93	0.51
1:A:916:ARG:NH1	1:B:875:GLU:OE2	2.44	0.51
2:A:1101:B12:H252	2:A:1101:B12:H601	1.94	0.50
1:B:879:ASN:O	1:B:883:THR:OG1	2.25	0.50
1:B:228:GLU:OE1	1:B:232:ARG:NH1	2.44	0.50
1:B:265:ARG:NH1	5:B:1103:GDP:O2A	2.44	0.50
1:B:30:VAL:HG23	1:B:78:VAL:HG11	1.93	0.50
1:A:30:VAL:HG23	1:A:78:VAL:HG11	1.93	0.49
1:A:697:ASP:OD1	1:A:1071:ARG:NH1	2.45	0.49
1:A:787:ASN:HB2	1:A:788:PRO:HD2	1.93	0.49
1:B:697:ASP:OD1	1:B:1071:ARG:NH1	2.45	0.49
1:A:513:LEU:HD21	1:A:566:LEU:HB2	1.94	0.49
1:B:787:ASN:HB2	1:B:788:PRO:HD2	1.94	0.49
1:B:296:ILE:O	1:B:300:LYS:HG3	2.13	0.49
1:A:296:ILE:O	1:A:300:LYS:HG3	2.13	0.49
1:B:731:VAL:O	1:B:733:ALA:N	2.47	0.48
1:A:734:ASP:O	1:A:737:LYS:HG2	2.13	0.48
1:B:513:LEU:HD21	1:B:566:LEU:HB2	1.94	0.48
1:B:267:ASN:HD21	1:B:1090:ARG:HA	1.78	0.48
1:A:250:PRO:HD3	1:A:312:SER:HB2	1.96	0.48
1:B:458:ARG:NH1	1:B:960:GLU:OE2	2.43	0.48
1:A:814:ASP:CG	1:A:854:ASN:HD22	2.17	0.48
1:B:866:THR:O	1:B:901:ASN:ND2	2.48	0.47
1:A:599:ALA:O	1:A:626:ALA:N	2.38	0.47
1:B:814:ASP:CG	1:B:854:ASN:HD22	2.17	0.47
1:B:685:PRO:HB3	1:B:759:ILE:HD11	1.97	0.47
1:A:780:HIS:CD2	4:A:1103:IVC:H44	2.50	0.47
1:A:49:ILE:HB	1:A:150:ILE:HG13	1.97	0.46
1:B:49:ILE:HB	1:B:150:ILE:HG13	1.97	0.46
1:B:141:GLY:HA3	2:B:1101:B12:H5R2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:250:PRO:HD3	1:B:312:SER:HB2	1.97	0.46
1:A:779:TYR:CD2	1:A:825:SER:HB2	2.51	0.46
1:A:528:TYR:HB2	1:A:541:LEU:HD21	1.98	0.46
1:B:446:ARG:NH2	1:B:804:GLU:OE2	2.48	0.46
1:B:528:TYR:HB2	1:B:541:LEU:HD21	1.98	0.46
1:A:172:VAL:HG21	1:A:183:LEU:HD22	1.97	0.46
1:A:46:MET:SD	1:A:150:ILE:HD11	2.56	0.46
1:B:531:LYS:HA	1:B:536:GLU:HA	1.96	0.46
1:B:231:ARG:NH1	1:B:1084:GLU:HA	2.31	0.46
1:A:905:GLU:O	2:A:1101:B12:N29	2.38	0.45
1:A:738:GLU:HB3	1:A:745:CYS:HB2	1.96	0.45
1:A:776:ILE:HD13	1:A:799:GLY:HA2	1.98	0.45
1:B:731:VAL:HG23	1:B:771:PHE:HE1	1.80	0.45
1:A:902:ALA:HB1	1:A:914:SER:OG	2.16	0.45
1:B:902:ALA:HB1	1:B:914:SER:OG	2.17	0.45
2:A:1101:B12:H261	3:A:1102[B]:5AD:H2'	1.98	0.45
1:A:598:PHE:HA	1:A:624:SER:HB3	1.99	0.45
2:A:1101:B12:H463	4:A:1103:IVC:H43	1.99	0.45
1:B:599:ALA:HB3	1:B:611:PHE:CE1	2.52	0.45
2:A:1101:B12:H552	2:A:1101:B12:H531	1.99	0.45
1:B:731:VAL:HG23	1:B:771:PHE:CE1	2.52	0.45
1:B:46:MET:SD	1:B:150:ILE:HD11	2.57	0.45
1:B:172:VAL:HG21	1:B:183:LEU:HD22	1.98	0.45
1:A:657:LEU:HG	1:A:661:LYS:HE3	1.99	0.45
1:B:287:GLU:HG2	1:B:288:ILE:HG23	1.98	0.44
2:B:1101:B12:H363	2:B:1101:B12:H411	1.80	0.44
1:B:32:ALA:HB2	1:B:60:LEU:HB2	1.99	0.44
1:A:32:ALA:HB2	1:A:60:LEU:HB2	1.98	0.44
1:B:657:LEU:HG	1:B:661:LYS:HE3	1.99	0.44
1:A:685:PRO:HB3	1:A:759:ILE:HD11	1.99	0.44
1:A:735:ILE:HG21	1:A:752:SER:HB2	1.99	0.44
1:A:62:HIS:CE1	1:A:743:ASN:HB3	2.52	0.44
1:A:597:MET:HB2	1:A:614:VAL:HG11	1.98	0.44
1:A:866:THR:O	1:A:901:ASN:ND2	2.51	0.44
1:A:978:ARG:NE	1:A:1093:MET:OXT	2.45	0.43
1:A:453:LEU:HD21	1:A:851:TYR:CE2	2.53	0.43
1:A:584:VAL:HB	1:A:859:LYS:HB3	1.99	0.43
1:A:997:LEU:HD22	1:B:546:LEU:HD11	2.01	0.43
1:A:735:ILE:HA	1:A:738:GLU:HG2	2.01	0.43
2:B:1101:B12:H531	2:B:1101:B12:H552	2.00	0.43
1:A:681:ASN:HB2	1:A:744:THR:HB	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1101:B12:H253	2:B:1101:B12:H301	1.70	0.42
1:A:731:VAL:HG12	1:A:771:PHE:HE1	1.85	0.42
1:B:193:ASP:HA	1:B:194:PRO:HD3	1.94	0.42
1:A:757:GLY:HA3	1:A:809:ARG:NH2	2.34	0.42
1:A:999:ILE:N	1:A:1003:ASN:OD1	2.47	0.42
1:B:70:VAL:HG13	1:B:101:LEU:HD23	2.02	0.42
1:B:1056:VAL:HB	1:B:1085:VAL:HG21	2.01	0.42
1:A:70:VAL:HG13	1:A:101:LEU:HD23	2.02	0.42
1:A:114:GLY:HA3	2:A:1101:B12:C9B	2.49	0.42
1:A:87:GLN:O	1:A:650:SER:OG	2.38	0.42
1:B:611:PHE:HB3	1:B:667:PHE:CZ	2.55	0.42
1:B:200:LEU:HD11	1:B:298:ALA:HA	2.01	0.42
1:B:464:VAL:HG11	1:B:571:ARG:HG3	2.02	0.42
1:B:160:ASP:OD2	1:B:162:THR:OG1	2.26	0.41
1:A:271:HIS:CG	1:A:272:PRO:HD2	2.55	0.41
1:B:222:LYS:NZ	5:B:1103:GDP:O1B	2.42	0.41
1:B:271:HIS:CG	1:B:272:PRO:HD2	2.55	0.41
1:A:778:GLY:HA3	1:A:795:THR:OG1	2.21	0.41
1:A:209:VAL:HA	1:A:210:PRO:HD3	1.88	0.41
1:A:772:TYR:CZ	4:A:1103:IVC:H20	2.55	0.41
1:A:222:LYS:NZ	5:A:1104:GDP:O1B	2.41	0.41
2:B:1101:B12:H481	2:B:1101:B12:H473	1.81	0.41
1:B:42:SER:OG	2:B:1101:B12:N3B	2.50	0.41
1:A:40:ASP:HA	1:A:43:ILE:HG22	2.03	0.41
1:A:1000:ILE:N	1:B:928:GLU:OE2	2.42	0.41
1:A:580:TYR:CE1	1:A:859:LYS:HE2	2.56	0.41
1:B:526:ASP:HA	1:B:542:ILE:HG12	2.03	0.41
1:A:737:LYS:O	1:A:741:GLY:N	2.47	0.40
1:B:358:LYS:HG2	5:B:1103:GDP:C6	2.56	0.40
1:A:443:PRO:HA	1:A:444:PRO:HD3	1.89	0.40
1:A:464:VAL:HG11	1:A:571:ARG:HG3	2.02	0.40
1:A:389:MET:HA	1:A:390:PRO:HD3	1.95	0.40
1:B:1023:SER:HB3	1:B:1026:GLU:HG3	2.03	0.40
1:A:829:ASP:HB2	1:A:832:TYR:CD2	2.55	0.40
1:B:248:ILE:HD13	1:B:289:SER:H	1.87	0.40
1:B:829:ASP:HB2	1:B:832:TYR:CD2	2.57	0.40
1:A:358:LYS:HG2	5:A:1104:GDP:C6	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	1053/1113 (95%)	1009 (96%)	42 (4%)	2 (0%)	52	87
1	B	1051/1113 (94%)	1014 (96%)	34 (3%)	3 (0%)	46	84
All	All	2104/2226 (94%)	2023 (96%)	76 (4%)	5 (0%)	52	87

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	732	GLN
1	B	1091	ARG
1	A	489	HIS
1	B	319	ALA
1	A	741	GLY

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	823/906 (91%)	818 (99%)	5 (1%)	90	97
1	B	817/906 (90%)	815 (100%)	2 (0%)	95	99
All	All	1640/1812 (90%)	1633 (100%)	7 (0%)	93	98

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	249	ASP

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Mol	Chain	Res	Type
1	A	598	PHE
1	A	744	THR
1	A	883	THR
1	A	1092	ASN
1	B	249	ASP
1	B	883	THR

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

Mol	Chain	Res	Type
1	A	865	GLN
1	B	899	HIS

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 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	$\# Z > 2$
2	B12	A	1101	1,3	74,101,101	0.59	1 (1%)	114,166,166	1.36	13 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	5AD	A	1102[A]	-	17,20,20	0.58	0	15,30,30	0.74	1 (6%)
3	5AD	A	1102[B]	2	17,20,20	0.57	0	15,30,30	0.34	0
4	IVC	A	1103	-	46,56,56	0.55	0	58,83,83	0.61	1 (1%)
5	GDP	A	1104	6	24,30,30	2.51	9 (37%)	26,47,47	1.67	6 (23%)
2	B12	B	1101	1	74,101,101	0.56	1 (1%)	114,166,166	1.40	14 (12%)
4	IVC	B	1102	-	25,29,56	0.60	0	30,45,83	0.72	1 (3%)
5	GDP	B	1103	6	24,30,30	2.54	9 (37%)	26,47,47	1.68	6 (23%)

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	B12	A	1101	1,3	-	0/51/223/223	0/3/11/11
3	5AD	A	1102[A]	-	-	0/0/20/20	0/3/3/3
3	5AD	A	1102[B]	2	-	0/0/20/20	0/3/3/3
4	IVC	A	1103	-	-	0/50/71/71	0/3/3/3
5	GDP	A	1104	6	-	0/12/32/32	0/3/3/3
2	B12	B	1101	1	-	0/51/223/223	0/3/11/11
4	IVC	B	1102	-	-	0/11/31/71	0/3/3/3
5	GDP	B	1103	6	-	0/12/32/32	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1104	GDP	O5'-C5'	-3.78	1.29	1.44
5	B	1103	GDP	O5'-C5'	-3.77	1.29	1.44
5	A	1104	GDP	C2-N1	-3.76	1.28	1.35
5	B	1103	GDP	C2-N1	-3.73	1.28	1.35
5	A	1104	GDP	C5-C4	-3.29	1.33	1.40
5	B	1103	GDP	C5-C4	-3.29	1.33	1.40
5	B	1103	GDP	C2'-C1'	-3.10	1.48	1.53
5	A	1104	GDP	C2'-C1'	-2.91	1.49	1.53
5	B	1103	GDP	PA-O2A	-2.89	1.42	1.55
5	B	1103	GDP	C2'-C3'	-2.78	1.45	1.53
5	A	1104	GDP	C2'-C3'	-2.73	1.46	1.53
5	B	1103	GDP	PB-O3B	-2.40	1.46	1.54
5	A	1104	GDP	PB-O3B	-2.36	1.46	1.54
5	A	1104	GDP	PA-O1A	-2.31	1.42	1.51
2	A	1101	B12	C1P-C2P	2.26	1.57	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1101	B12	C1P-C2P	2.30	1.57	1.51
5	A	1104	GDP	O6-C6	4.90	1.37	1.24
5	B	1103	GDP	O6-C6	4.93	1.37	1.24
5	A	1104	GDP	C2-N2	6.67	1.48	1.34
5	B	1103	GDP	C2-N2	6.69	1.48	1.34

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1104	GDP	N3-C2-N1	-4.28	121.74	127.56
5	B	1103	GDP	N3-C2-N1	-4.23	121.80	127.56
2	B	1101	B12	C35-C5-C4	-3.95	112.38	117.81
2	A	1101	B12	C2-C1-C19	-3.83	115.00	118.62
2	A	1101	B12	C35-C5-C4	-3.56	112.91	117.81
2	A	1101	B12	C30-C3-C2	-3.56	112.00	119.01
2	B	1101	B12	C9-C10-C11	-3.41	124.06	132.31
2	B	1101	B12	C30-C3-C2	-3.25	112.61	119.01
2	A	1101	B12	C9-C10-C11	-3.19	124.59	132.31
2	A	1101	B12	C13-C14-C15	-3.18	122.07	132.12
5	A	1104	GDP	C4'-O4'-C1'	-3.07	106.39	109.64
5	B	1103	GDP	C4'-O4'-C1'	-3.06	106.39	109.64
2	B	1101	B12	C13-C14-C15	-2.99	122.67	132.12
2	B	1101	B12	C3-C4-C5	-2.97	122.72	132.12
2	B	1101	B12	C2-C1-C19	-2.97	115.81	118.62
5	B	1103	GDP	C5-C6-N1	-2.94	119.68	123.52
2	A	1101	B12	C3-C4-C5	-2.94	122.84	132.12
5	A	1104	GDP	C5-C6-N1	-2.87	119.77	123.52
3	A	1102[A]	5AD	C4'-O4'-C1'	-2.73	106.75	109.64
5	A	1104	GDP	C1'-N9-C4	-2.54	123.97	126.81
2	B	1101	B12	C25-C2-C3	-2.50	112.87	115.90
5	B	1103	GDP	C1'-N9-C4	-2.49	124.03	126.81
2	B	1101	B12	C55-C17-C18	-2.49	106.57	111.03
5	A	1104	GDP	C6-C5-C4	-2.33	118.20	120.86
5	B	1103	GDP	C6-C5-C4	-2.26	118.27	120.86
2	A	1101	B12	C20-C1-C19	-2.22	107.16	109.36
2	B	1101	B12	C16-C15-C14	-2.16	120.42	124.08
2	B	1101	B12	C20-C1-C19	-2.15	107.23	109.36
2	A	1101	B12	C16-C15-C14	-2.10	120.53	124.08
2	A	1101	B12	C53-C15-C16	2.01	120.58	117.81
2	B	1101	B12	O3-P-O4	2.21	116.58	108.83
2	B	1101	B12	P-O2-C3R	2.24	126.73	120.10
4	A	1103	IVC	P1-O2-C1	2.31	127.47	121.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1101	B12	C8-C7-C6	2.41	104.21	101.00
4	B	1102	IVC	O4-P2-O10	2.70	113.87	107.08
2	A	1101	B12	C19-C1-N21	3.34	105.69	102.16
5	B	1103	GDP	C6-N1-C2	3.61	120.11	115.88
5	A	1104	GDP	C6-N1-C2	3.63	120.14	115.88
2	A	1101	B12	C35-C5-C6	4.97	124.64	117.81
2	A	1101	B12	P-O3-C2P	5.15	128.21	120.94
2	B	1101	B12	P-O3-C2P	5.36	128.50	120.94
2	B	1101	B12	C35-C5-C6	5.88	125.89	117.81

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 29 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1101	B12	10	0
3	A	1102[A]	5AD	1	0
3	A	1102[B]	5AD	2	0
4	A	1103	IVC	7	0
5	A	1104	GDP	3	0
2	B	1101	B12	6	0
5	B	1103	GDP	3	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	1061/1113 (95%)	-0.08	14 (1%) 79 72	58, 88, 139, 180	0
1	B	1059/1113 (95%)	-0.11	22 (2%) 67 61	55, 82, 137, 212	0
All	All	2120/2226 (95%)	-0.09	36 (1%) 73 66	55, 85, 139, 212	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	909	THR	5.5
1	B	908	THR	4.1
1	A	1011	ASP	3.7
1	B	907	ILE	3.6
1	A	489	HIS	3.4
1	A	255	SER	3.3
1	A	420	MET	3.2
1	A	528	TYR	3.2
1	A	537	ILE	3.2
1	A	1015	GLN	2.9
1	A	316	GLN	2.7
1	A	317	GLY	2.7
1	B	1020	ALA	2.7
1	B	874	GLN	2.6
1	A	808	ALA	2.6
1	B	536	GLU	2.6
1	B	534	ASP	2.6
1	B	23	PRO	2.5
1	B	902	ALA	2.5
1	B	494	LEU	2.4
1	B	1019	LEU	2.4
1	A	677	SER	2.4
1	B	707	ASN	2.3
1	B	537	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	254	LYS	2.3
1	B	872	HIS	2.3
1	B	706	ASP	2.2
1	B	903	TYR	2.2
1	B	595	THR	2.2
1	B	530	VAL	2.2
1	A	1010	GLY	2.1
1	B	533	ARG	2.1
1	A	707	ASN	2.0
1	B	901	ASN	2.0
1	B	620	ALA	2.0
1	B	538	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
6	MG	A	1106	1/1	0.94	0.28	2.16	78,78,78,78	0
6	MG	B	1105	1/1	0.98	0.31	1.23	43,43,43,43	0
4	IVC	B	1102	27/54	0.83	0.27	0.68	130,147,183,185	0
6	MG	B	1104	1/1	0.99	0.28	0.51	42,42,42,42	0
3	5AD	A	1102[A]	18/18	0.93	0.23	0.42	107,112,115,115	18
3	5AD	A	1102[B]	18/18	0.93	0.23	0.40	108,113,115,115	18
2	B12	B	1101	91/91	0.93	0.25	0.39	91,113,136,139	0
5	GDP	B	1103	28/28	0.95	0.20	-0.06	55,65,72,123	0
2	B12	A	1101	91/91	0.97	0.19	-0.33	51,65,80,83	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	IVC	A	1103	54/54	0.91	0.22	-0.51	76,114,161,164	0
5	GDP	A	1104	28/28	0.96	0.16	-0.77	70,89,96,98	0
6	MG	A	1105	1/1	0.98	0.29	-	64,64,64,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.