



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 5, 2018 – 11:12 PM EST

PDB ID : 1K98
Title : AdoMet complex of MetH C-terminal fragment
Authors : Bandarian, V.; Pattridge, K.A.; Lennon, B.W.; Huddler, D.P.; Matthews, R.G.; Ludwig, M.L.
Deposited on : 2001-10-27
Resolution : 3.75 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030736

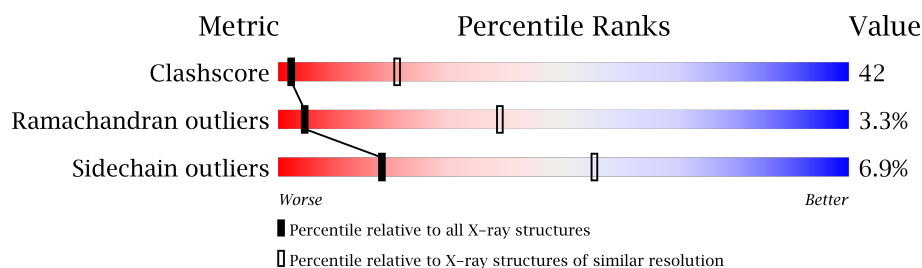
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.75 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))
Clashscore	112137	1087 (4.00-3.52)
Ramachandran outliers	110173	1047 (4.00-3.52)
Sidechain outliers	110143	1041 (4.00-3.52)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	577	

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	B12	A	1248	X	-	-	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4662 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 Methionine synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	577	4566	2888	787	874	17	0	0	0

There is a discrepancy between the modelled and reference sequences:

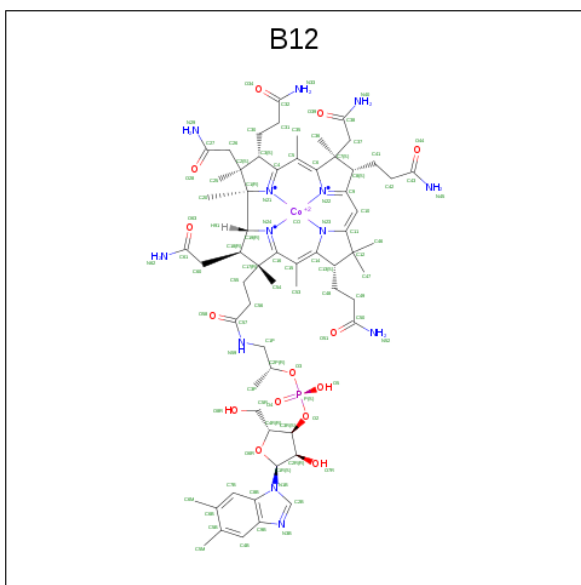
Chain	Residue	Modelled	Actual	Comment	Reference
A	759	GLY	HIS	ENGINEERED	UNP P13009

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0

- Molecule 3 is COBALAMIN (three-letter code: B12) (formula: C₆₂H₈₉CoN₁₃O₁₄P).



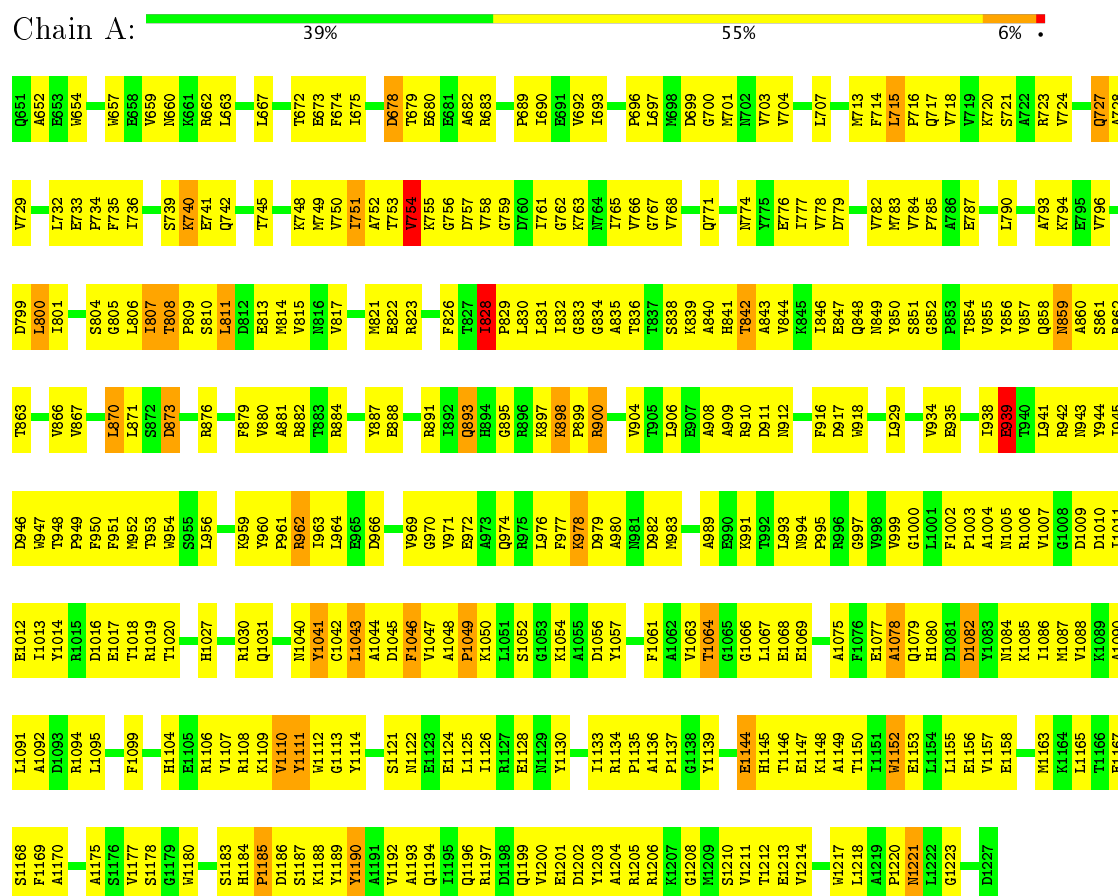
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	Co	N	O	P	0	0
			91	62	1	13	14	1		

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 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.

Note EDS was not executed.

• Molecule 1: Methionine synthase



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	108.50 Å 108.50 Å 146.47 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 3.75	Depositor
% Data completeness (in resolution range)	5.3 (15.00-3.75)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.308 , 0.363	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4662	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: B12, SO4

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	0.56	1/4664 (0.0%)	0.77	3/6326 (0.0%)

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	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	740	LYS	CB-CG	5.01	1.66	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	740	LYS	CD-CE-NZ	9.21	132.88	111.70
1	A	740	LYS	CB-CG-CD	8.88	134.68	111.60
1	A	811	LEU	CA-CB-CG	5.33	127.56	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1190	TYR	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	4566	0	4484	381	0
2	A	5	0	0	0	0
3	A	91	0	77	20	0
All	All	4662	0	4561	391	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 42.

All (391) 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)
3:A:1248:B12:C56	3:A:1248:B12:C55	1.75	1.57
3:A:1248:B12:C26	3:A:1248:B12:C27	1.80	1.55
3:A:1248:B12:C18	3:A:1248:B12:C60	1.80	1.52
3:A:1248:B12:C53	3:A:1248:B12:C15	1.88	1.51
3:A:1248:B12:C57	3:A:1248:B12:C55	2.38	1.00
1:A:1152:TRP:HA	1:A:1157:VAL:HG23	1.45	0.97
1:A:1178:SER:H	3:A:1248:B12:H402	1.09	0.96
1:A:898:LYS:O	1:A:900:ARG:N	1.97	0.96
1:A:748:LYS:HE2	1:A:776:GLU:OE2	1.67	0.92
1:A:1165:LEU:HD23	1:A:1170:ALA:O	1.70	0.91
1:A:813:GLU:O	1:A:817:VAL:HG23	1.73	0.89
1:A:672:THR:HG23	1:A:724:VAL:HG22	1.55	0.88
1:A:752:ALA:HB1	1:A:784:VAL:HG11	1.59	0.85
1:A:840:ALA:O	1:A:844:VAL:HG23	1.78	0.83
1:A:866:VAL:HG12	1:A:870:LEU:HD22	1.59	0.82
1:A:1075:ALA:O	1:A:1078:ALA:HB3	1.80	0.81
1:A:1078:ALA:O	1:A:1080:HIS:N	2.12	0.81
1:A:809:PRO:HB2	1:A:1168:SER:OG	1.83	0.79
1:A:948:THR:HB	1:A:949:PRO:HD3	1.65	0.79
1:A:768:VAL:HA	1:A:771:GLN:HE21	1.48	0.79
1:A:1152:TRP:HA	1:A:1157:VAL:CG2	2.12	0.78
1:A:793:ALA:HB2	1:A:801:ILE:HD11	1.65	0.78
1:A:753:THR:HB	1:A:782:VAL:HA	1.66	0.78
1:A:822:GLU:OE2	1:A:822:GLU:HA	1.84	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:966:ASP:OD2	1:A:969:VAL:HG12	1.82	0.78
1:A:1112:TRP:CZ2	1:A:1183:SER:HB3	2.20	0.77
1:A:997:GLY:HA3	1:A:1063:VAL:HG12	1.67	0.76
1:A:762:GLY:O	1:A:766:VAL:HG23	1.85	0.76
1:A:748:LYS:HE2	1:A:776:GLU:CD	2.06	0.76
1:A:897:LYS:HA	1:A:900:ARG:HH21	1.50	0.75
1:A:964:LEU:HD22	1:A:974:GLN:HG2	1.69	0.75
1:A:954:TRP:O	1:A:956:LEU:HD13	1.86	0.74
1:A:754:VAL:HG13	1:A:755:LYS:N	2.03	0.74
1:A:1090:ALA:O	1:A:1094:ARG:HG2	1.89	0.73
1:A:663:LEU:HD11	1:A:679:THR:HA	1.70	0.72
1:A:753:THR:CB	1:A:782:VAL:HG12	2.20	0.71
1:A:1201:GLU:O	1:A:1204:ALA:HB3	1.91	0.71
1:A:1104:HIS:ND1	1:A:1134:ARG:NH1	2.37	0.71
1:A:736:ILE:O	1:A:742:GLN:HG2	1.90	0.71
1:A:1108:ARG:HG3	1:A:1114:TYR:OH	1.90	0.71
1:A:956:LEU:HG	1:A:969:VAL:HG21	1.73	0.70
1:A:1056:ASP:OD1	1:A:1057:TYR:N	2.24	0.70
1:A:1048:ALA:HB3	1:A:1056:ASP:HB2	1.73	0.69
1:A:1133:ILE:HG13	1:A:1135:PRO:HD3	1.75	0.69
1:A:822:GLU:OE1	1:A:851:SER:HB3	1.92	0.69
1:A:1006:ARG:HB3	1:A:1049:PRO:HA	1.74	0.69
1:A:790:LEU:HB3	1:A:826:PHE:CE2	2.27	0.69
1:A:946:ASP:HB2	1:A:1126:ILE:CD1	2.21	0.69
1:A:749:MET:HE2	1:A:800:LEU:HG	1.74	0.69
1:A:970:GLY:O	1:A:974:GLN:HG3	1.92	0.69
1:A:759:GLY:HA3	1:A:806:LEU:HD22	1.74	0.68
1:A:1137:PRO:HB2	1:A:1148:LYS:HG3	1.73	0.68
1:A:1202:ASP:O	1:A:1206:ARG:HG3	1.93	0.68
3:A:1248:B12:C17	3:A:1248:B12:C60	2.71	0.68
1:A:1188:LYS:HB2	1:A:1190:TYR:CE1	2.29	0.67
1:A:962:ARG:CZ	1:A:962:ARG:HB2	2.25	0.67
1:A:971:VAL:O	1:A:974:GLN:HB2	1.95	0.66
1:A:1078:ALA:C	1:A:1080:HIS:H	1.99	0.66
1:A:947:TRP:O	1:A:950:PHE:HB3	1.95	0.66
1:A:1046:PHE:CZ	1:A:1187:SER:HA	2.30	0.66
1:A:1178:SER:N	3:A:1248:B12:H402	1.88	0.66
1:A:876:ARG:HH12	1:A:880:VAL:HG11	1.61	0.65
1:A:980:ALA:HB1	1:A:1091:LEU:HD21	1.77	0.65
1:A:793:ALA:CB	1:A:801:ILE:HD11	2.26	0.65
1:A:1014:TYR:OH	1:A:1050:LYS:HE2	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:753:THR:O	1:A:754:VAL:O	2.15	0.65
1:A:859:ASN:ND2	1:A:861:SER:HB3	2.11	0.65
1:A:1205:ARG:O	1:A:1208:GLY:N	2.30	0.64
1:A:906:LEU:HD22	1:A:1199:GLN:HA	1.80	0.64
1:A:1134:ARG:HA	1:A:1180:TRP:O	1.98	0.64
1:A:693:ILE:HD12	1:A:729:VAL:HG22	1.81	0.63
1:A:859:ASN:HD21	1:A:861:SER:HB3	1.63	0.63
1:A:1011:ILE:HD11	1:A:1044:ALA:HA	1.78	0.63
1:A:811:LEU:O	1:A:815:VAL:HG23	1.99	0.63
1:A:768:VAL:CG2	1:A:1082:ASP:OD2	2.46	0.63
1:A:749:MET:CE	1:A:831:LEU:HD22	2.28	0.63
3:A:1248:B12:C60	3:A:1248:B12:H18	2.18	0.63
1:A:946:ASP:HB2	1:A:1126:ILE:HD12	1.81	0.63
1:A:1107:VAL:HA	1:A:1111:TYR:HB2	1.80	0.63
1:A:1077:GLU:HG2	1:A:1085:LYS:HD2	1.80	0.62
1:A:950:PHE:HE2	1:A:976:LEU:HD22	1.64	0.62
1:A:748:LYS:NZ	1:A:796:VAL:HG22	2.14	0.62
1:A:830:LEU:HD23	1:A:854:THR:HG23	1.81	0.62
1:A:809:PRO:CB	1:A:1168:SER:OG	2.47	0.62
1:A:768:VAL:HA	1:A:771:GLN:NE2	2.15	0.61
1:A:751:ILE:HG23	1:A:763:LYS:HD2	1.83	0.61
1:A:959:LYS:HG2	1:A:960:TYR:N	2.13	0.61
1:A:999:VAL:HG22	1:A:1000:GLY:H	1.65	0.61
1:A:749:MET:HB2	1:A:800:LEU:HD21	1.83	0.60
1:A:999:VAL:HG22	1:A:1000:GLY:N	2.16	0.60
1:A:929:LEU:HD22	1:A:1003:PRO:HD3	1.84	0.60
1:A:954:TRP:CG	1:A:976:LEU:HD11	2.36	0.60
1:A:876:ARG:NH1	1:A:880:VAL:HG11	2.16	0.60
1:A:1177:VAL:HA	3:A:1248:B12:N40	2.16	0.60
1:A:749:MET:CE	1:A:800:LEU:HG	2.32	0.60
1:A:918:TRP:O	1:A:1054:LYS:HE3	2.03	0.59
1:A:1188:LYS:HB2	1:A:1190:TYR:CZ	2.37	0.59
3:A:1248:B12:C57	3:A:1248:B12:H552	2.33	0.59
1:A:839:LYS:HA	1:A:856:TYR:CD2	2.38	0.59
1:A:749:MET:HE3	1:A:831:LEU:HD22	1.83	0.58
1:A:993:LEU:HD11	1:A:1066:GLY:HA3	1.86	0.58
1:A:753:THR:OG1	1:A:782:VAL:HG12	2.03	0.58
1:A:842:THR:HG23	1:A:846:ILE:HD12	1.85	0.58
1:A:1067:LEU:C	1:A:1069:GLU:H	2.08	0.58
1:A:880:VAL:O	1:A:884:ARG:HG3	2.04	0.58
1:A:1004:ALA:HB1	1:A:1011:ILE:HG22	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:756:GLY:HA2	1:A:1167:GLU:OE1	2.04	0.57
1:A:654:TRP:HA	1:A:657:TRP:CE2	2.39	0.57
1:A:1145:HIS:HD2	1:A:1148:LYS:NZ	2.02	0.57
1:A:785:PRO:C	1:A:787:GLU:N	2.55	0.57
1:A:790:LEU:HD23	1:A:801:ILE:HD13	1.86	0.57
1:A:1077:GLU:HG2	1:A:1085:LYS:CD	2.34	0.57
1:A:1152:TRP:CD2	1:A:1157:VAL:HG21	2.39	0.57
1:A:750:VAL:HA	1:A:778:VAL:HG13	1.86	0.57
1:A:1144:GLU:OE1	1:A:1146:THR:HG23	2.04	0.56
1:A:754:VAL:HG13	1:A:755:LYS:H	1.68	0.56
1:A:714:PHE:HD2	1:A:972:GLU:HG3	1.70	0.56
1:A:753:THR:O	1:A:784:VAL:HG12	2.04	0.56
1:A:938:ILE:O	1:A:939:GLU:C	2.42	0.56
1:A:950:PHE:CE2	1:A:976:LEU:HD22	2.40	0.56
1:A:826:PHE:CD1	1:A:828:ILE:HD11	2.41	0.56
1:A:951:PHE:HZ	1:A:977:PHE:HB2	1.71	0.56
1:A:1196:GLN:O	1:A:1200:VAL:HG23	2.06	0.56
1:A:1121:SER:OG	1:A:1124:GLU:HG3	2.05	0.56
1:A:866:VAL:CG1	1:A:870:LEU:HD22	2.30	0.56
1:A:1067:LEU:O	1:A:1069:GLU:N	2.39	0.55
1:A:939:GLU:HA	1:A:939:GLU:OE1	2.07	0.55
1:A:832:ILE:HD11	1:A:842:THR:OG1	2.06	0.55
1:A:946:ASP:O	1:A:949:PRO:HD2	2.06	0.55
1:A:859:ASN:ND2	1:A:862:ARG:H	2.04	0.55
1:A:1135:PRO:HG2	1:A:1180:TRP:HB2	1.88	0.55
1:A:1004:ALA:HB1	1:A:1011:ILE:CG2	2.36	0.55
1:A:1007:VAL:HG21	1:A:1012:GLU:OE2	2.07	0.55
1:A:790:LEU:HD23	1:A:801:ILE:CD1	2.36	0.55
1:A:1109:LYS:O	1:A:1113:GLY:HA2	2.07	0.55
1:A:859:ASN:HD22	1:A:861:SER:H	1.55	0.55
1:A:753:THR:HB	1:A:782:VAL:HG12	1.89	0.54
1:A:1133:ILE:HB	1:A:1189:TYR:CZ	2.43	0.54
1:A:997:GLY:CA	1:A:1063:VAL:HG12	2.36	0.54
1:A:828:ILE:H	1:A:828:ILE:HD13	1.71	0.54
1:A:959:LYS:CG	1:A:960:TYR:N	2.70	0.54
1:A:704:VAL:HG13	1:A:713:MET:CE	2.37	0.54
1:A:790:LEU:HB3	1:A:826:PHE:HE2	1.70	0.54
1:A:822:GLU:HB2	1:A:849:ASN:O	2.07	0.54
1:A:906:LEU:O	1:A:906:LEU:HD12	2.07	0.54
1:A:690:ILE:HG13	1:A:690:ILE:O	2.06	0.53
1:A:1014:TYR:HH	1:A:1050:LYS:HE2	1.72	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:693:ILE:CD1	1:A:729:VAL:HG22	2.39	0.53
1:A:697:LEU:HD21	1:A:728:ALA:CB	2.38	0.53
1:A:715:LEU:O	1:A:715:LEU:HD13	2.08	0.53
1:A:916:PHE:CZ	1:A:1184:HIS:HE1	2.26	0.53
1:A:960:TYR:CD2	1:A:961:PRO:HA	2.44	0.53
1:A:1046:PHE:CE1	1:A:1187:SER:HA	2.44	0.53
1:A:778:VAL:HG13	1:A:778:VAL:O	2.08	0.53
1:A:1145:HIS:CD2	1:A:1148:LYS:NZ	2.77	0.52
1:A:1188:LYS:HD2	1:A:1190:TYR:OH	2.08	0.52
1:A:758:VAL:HG23	1:A:782:VAL:HG21	1.90	0.52
1:A:785:PRO:C	1:A:787:GLU:H	2.12	0.52
1:A:810:SER:O	1:A:813:GLU:HB2	2.09	0.52
1:A:1007:VAL:HG23	1:A:1007:VAL:O	2.10	0.52
1:A:1221:ASN:N	1:A:1221:ASN:OD1	2.42	0.52
1:A:736:ILE:O	1:A:739:SER:OG	2.25	0.52
1:A:989:ALA:O	1:A:991:LYS:HG2	2.10	0.52
1:A:804:SER:HA	1:A:833:GLY:O	2.10	0.52
1:A:909:ALA:O	1:A:912:ASN:N	2.42	0.52
1:A:683:ARG:HD3	1:A:732:LEU:HD23	1.92	0.52
1:A:859:ASN:HD22	1:A:861:SER:N	2.08	0.52
1:A:1007:VAL:HG11	1:A:1012:GLU:OE2	2.10	0.51
1:A:672:THR:CG2	1:A:724:VAL:HG22	2.31	0.51
1:A:843:ALA:HB1	1:A:887:TYR:HB3	1.93	0.51
1:A:675:ILE:HD11	1:A:728:ALA:HB2	1.93	0.51
1:A:692:VAL:HG21	1:A:732:LEU:HD11	1.93	0.51
1:A:1144:GLU:HG3	1:A:1203:TYR:CE2	2.46	0.51
1:A:887:TYR:O	1:A:891:ARG:HB2	2.10	0.51
1:A:767:GLY:O	1:A:771:GLN:HG3	2.10	0.51
1:A:1112:TRP:CZ2	1:A:1183:SER:CB	2.93	0.51
1:A:717:GLN:O	1:A:720:LYS:HB3	2.10	0.51
1:A:748:LYS:HZ2	1:A:796:VAL:HG22	1.73	0.51
3:A:1248:B12:C53	3:A:1248:B12:C14	2.81	0.51
1:A:713:MET:O	1:A:714:PHE:CD1	2.64	0.51
1:A:832:ILE:HD13	1:A:846:ILE:CD1	2.41	0.51
1:A:1152:TRP:CE3	1:A:1165:LEU:HD11	2.45	0.51
1:A:893:GLN:NE2	1:A:893:GLN:O	2.44	0.51
3:A:1248:B12:C26	3:A:1248:B12:O28	2.47	0.50
1:A:696:PRO:HG2	1:A:697:LEU:H	1.76	0.50
1:A:1014:TYR:CD2	1:A:1019:ARG:HG2	2.46	0.50
1:A:1067:LEU:C	1:A:1069:GLU:N	2.65	0.50
1:A:1077:GLU:O	1:A:1078:ALA:O	2.29	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:943:ASN:N	1:A:943:ASN:HD22	2.08	0.50
1:A:768:VAL:HG22	1:A:1082:ASP:OD2	2.10	0.50
1:A:832:ILE:CD1	1:A:842:THR:HG23	2.42	0.50
1:A:1004:ALA:O	1:A:1005:ASN:ND2	2.45	0.50
1:A:826:PHE:CD1	1:A:828:ILE:CD1	2.94	0.50
1:A:790:LEU:CD2	1:A:801:ILE:CD1	2.89	0.50
1:A:757:ASP:OD1	1:A:806:LEU:N	2.44	0.50
1:A:1061:PHE:O	1:A:1178:SER:HA	2.12	0.50
1:A:754:VAL:CG1	1:A:755:LYS:N	2.74	0.50
1:A:1148:LYS:HE2	1:A:1169:PHE:O	2.11	0.50
1:A:1192:VAL:O	1:A:1193:ALA:HB3	2.10	0.50
1:A:859:ASN:HA	3:A:1248:B12:H2R	1.93	0.50
1:A:956:LEU:HG	1:A:969:VAL:CG2	2.41	0.50
1:A:807:ILE:HD11	1:A:1170:ALA:HB2	1.94	0.50
1:A:1202:ASP:OD1	1:A:1205:ARG:NH2	2.45	0.50
1:A:964:LEU:HD22	1:A:974:GLN:CG	2.39	0.50
1:A:762:GLY:H	3:A:1248:B12:C3P	2.24	0.49
1:A:707:LEU:HD13	1:A:713:MET:SD	2.52	0.49
1:A:983:MET:HE1	1:A:1087:MET:HB3	1.94	0.49
1:A:745:THR:HB	1:A:774:ASN:O	2.12	0.49
1:A:1177:VAL:HA	3:A:1248:B12:H401	1.75	0.49
1:A:857:VAL:HG11	1:A:863:THR:HA	1.95	0.49
1:A:1046:PHE:CE2	1:A:1187:SER:HA	2.47	0.49
1:A:654:TRP:H	1:A:654:TRP:HD1	1.61	0.49
1:A:1134:ARG:HG3	1:A:1134:ARG:O	2.13	0.49
1:A:993:LEU:CD1	1:A:1066:GLY:HA3	2.42	0.49
1:A:1211:VAL:O	1:A:1214:VAL:HG22	2.13	0.49
1:A:855:VAL:HG21	1:A:870:LEU:HD21	1.95	0.49
1:A:828:ILE:N	1:A:828:ILE:HD13	2.27	0.48
1:A:863:THR:O	1:A:867:VAL:HG23	2.14	0.48
1:A:1030:ARG:HG3	1:A:1042:CYS:HB2	1.94	0.48
1:A:701:MET:O	1:A:701:MET:HE3	2.14	0.48
1:A:1145:HIS:HD2	1:A:1148:LYS:HZ1	1.62	0.48
1:A:1186:ASP:O	1:A:1188:LYS:HG3	2.14	0.48
1:A:1211:VAL:HG13	1:A:1212:THR:N	2.28	0.48
1:A:768:VAL:CA	1:A:771:GLN:HE21	2.21	0.48
1:A:880:VAL:HG23	1:A:881:ALA:N	2.29	0.48
1:A:857:VAL:HB	3:A:1248:B12:HM61	1.95	0.48
1:A:799:ASP:O	1:A:829:PRO:HD2	2.13	0.48
1:A:1137:PRO:CB	1:A:1148:LYS:HG3	2.40	0.48
1:A:1009:ASP:HB3	1:A:1042:CYS:SG	2.53	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:983:MET:CE	1:A:1087:MET:HB3	2.43	0.48
1:A:950:PHE:O	1:A:953:THR:HB	2.14	0.48
1:A:934:VAL:HG12	1:A:935:GLU:H	1.78	0.48
1:A:971:VAL:HA	1:A:974:GLN:NE2	2.29	0.48
1:A:949:PRO:HA	1:A:952:MET:HG3	1.95	0.47
1:A:1048:ALA:CB	1:A:1056:ASP:HB2	2.44	0.47
1:A:696:PRO:O	1:A:699:ASP:HB2	2.15	0.47
1:A:898:LYS:C	1:A:900:ARG:N	2.67	0.47
1:A:714:PHE:O	1:A:718:VAL:HG23	2.15	0.47
1:A:910:ARG:HG3	1:A:1030:ARG:HE	1.78	0.47
1:A:978:LYS:HD3	1:A:982:ASP:OD1	2.15	0.47
1:A:946:ASP:HB2	1:A:1126:ILE:HD11	1.96	0.47
1:A:1016:ASP:OD1	1:A:1018:THR:HG23	2.14	0.47
1:A:870:LEU:HD13	1:A:879:PHE:CD1	2.50	0.47
1:A:1165:LEU:CD2	1:A:1170:ALA:O	2.53	0.47
1:A:654:TRP:N	1:A:654:TRP:CD1	2.83	0.47
1:A:752:ALA:O	1:A:804:SER:N	2.47	0.47
1:A:799:ASP:HB3	1:A:871:LEU:HD21	1.95	0.47
1:A:994:ASN:HA	1:A:995:PRO:HD3	1.74	0.47
1:A:662:ARG:HB3	1:A:678:ASP:OD2	2.14	0.47
1:A:1108:ARG:HG3	1:A:1114:TYR:CZ	2.49	0.46
1:A:1194:GLN:HB3	1:A:1223:GLY:CA	2.45	0.46
1:A:680:GLU:O	1:A:683:ARG:N	2.36	0.46
1:A:1108:ARG:HH22	1:A:1125:LEU:HD21	1.81	0.46
1:A:707:LEU:CD1	1:A:713:MET:SD	3.03	0.46
1:A:751:ILE:HD12	1:A:763:LYS:HG3	1.97	0.46
1:A:1211:VAL:O	1:A:1214:VAL:CG2	2.64	0.46
1:A:832:ILE:HG13	1:A:832:ILE:O	2.15	0.46
1:A:888:GLU:O	1:A:891:ARG:HB3	2.14	0.46
1:A:672:THR:O	1:A:673:GLU:HG3	2.16	0.46
1:A:749:MET:HE1	1:A:831:LEU:HD22	1.98	0.46
1:A:951:PHE:CZ	1:A:977:PHE:HB2	2.51	0.46
1:A:1091:LEU:O	1:A:1092:ALA:C	2.53	0.46
1:A:1108:ARG:HH22	1:A:1125:LEU:CD2	2.27	0.46
1:A:808:THR:HG22	1:A:809:PRO:N	2.31	0.46
1:A:768:VAL:HG21	1:A:1082:ASP:OD2	2.15	0.46
1:A:1184:HIS:CG	1:A:1185:PRO:HD2	2.51	0.46
1:A:838:SER:O	1:A:839:LYS:C	2.54	0.46
1:A:1095:LEU:HD23	1:A:1095:LEU:HA	1.73	0.45
1:A:692:VAL:O	1:A:696:PRO:HG2	2.16	0.45
1:A:857:VAL:HG22	1:A:866:VAL:HG21	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1205:ARG:O	1:A:1206:ARG:C	2.55	0.45
1:A:763:LYS:HZ3	1:A:779:ASP:CG	2.20	0.45
1:A:828:ILE:HB	1:A:829:PRO:HD2	1.99	0.45
1:A:762:GLY:H	3:A:1248:B12:H3P2	1.81	0.45
1:A:948:THR:O	1:A:949:PRO:C	2.53	0.45
1:A:1157:VAL:HG12	1:A:1163:MET:HB2	1.98	0.45
1:A:964:LEU:HD22	1:A:974:GLN:HA	1.97	0.45
1:A:1106:ARG:NE	1:A:1111:TYR:OH	2.48	0.45
1:A:943:ASN:N	1:A:943:ASN:ND2	2.64	0.45
1:A:964:LEU:CD2	1:A:974:GLN:HA	2.46	0.45
1:A:1031:GLN:OE1	1:A:1041:TYR:HD2	2.00	0.45
1:A:994:ASN:O	1:A:1066:GLY:HA2	2.16	0.45
1:A:1167:GLU:H	1:A:1167:GLU:CD	2.20	0.45
1:A:790:LEU:O	1:A:793:ALA:HB3	2.17	0.45
1:A:1106:ARG:NH2	1:A:1111:TYR:OH	2.48	0.45
1:A:1213:GLU:O	1:A:1217:TRP:HD1	1.99	0.45
1:A:1064:THR:HA	1:A:1175:ALA:O	2.16	0.44
1:A:720:LYS:O	1:A:723:ARG:HB3	2.17	0.44
1:A:879:PHE:O	1:A:882:ARG:N	2.50	0.44
1:A:1042:CYS:O	1:A:1044:ALA:N	2.50	0.44
1:A:1110:VAL:HB	1:A:1111:TYR:H	1.60	0.44
1:A:1218:LEU:C	1:A:1220:PRO:HD2	2.37	0.44
1:A:667:LEU:HD13	1:A:721:SER:HB3	2.00	0.44
1:A:732:LEU:HB3	1:A:736:ILE:HG13	1.98	0.44
1:A:823:ARG:HG2	1:A:823:ARG:HH11	1.82	0.44
1:A:1043:LEU:HA	1:A:1043:LEU:HD23	1.75	0.44
1:A:1040:ASN:O	1:A:1041:TYR:O	2.36	0.44
1:A:697:LEU:O	1:A:700:GLY:N	2.51	0.44
1:A:703:VAL:O	1:A:704:VAL:C	2.54	0.44
1:A:749:MET:O	1:A:778:VAL:HG12	2.18	0.44
1:A:832:ILE:HD13	1:A:846:ILE:HD12	1.99	0.44
1:A:854:THR:O	1:A:887:TYR:HE1	2.00	0.44
1:A:765:ILE:HG13	1:A:1086:ILE:HD13	1.99	0.44
1:A:821:MET:HB3	1:A:850:TYR:CE1	2.52	0.44
1:A:980:ALA:O	1:A:983:MET:N	2.51	0.44
1:A:1061:PHE:CD1	1:A:1061:PHE:N	2.86	0.43
1:A:805:GLY:O	1:A:834:GLY:HA3	2.17	0.43
1:A:839:LYS:N	1:A:856:TYR:CE2	2.85	0.43
1:A:950:PHE:HB2	1:A:1094:ARG:HD2	1.99	0.43
1:A:716:PRO:CG	1:A:976:LEU:HD12	2.48	0.43
1:A:908:ALA:O	1:A:911:ASP:HB3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1139:TYR:CE1	3:A:1248:B12:H352	2.54	0.43
1:A:723:ARG:HG3	1:A:727:GLN:HE21	1.84	0.43
1:A:1005:ASN:OD1	1:A:1019:ARG:NH1	2.51	0.43
1:A:1206:ARG:HH11	1:A:1206:ARG:HG3	1.83	0.43
1:A:806:LEU:HD12	1:A:835:ALA:H	1.84	0.43
1:A:1136:ALA:HB1	1:A:1178:SER:O	2.19	0.43
1:A:1137:PRO:HB2	1:A:1148:LYS:CG	2.46	0.43
1:A:1155:LEU:O	1:A:1156:GLU:C	2.56	0.43
1:A:689:PRO:HB2	1:A:736:ILE:HA	2.00	0.43
1:A:763:LYS:NZ	1:A:779:ASP:OD1	2.50	0.43
1:A:1064:THR:C	1:A:1066:GLY:H	2.22	0.43
1:A:1126:ILE:O	1:A:1126:ILE:HG22	2.18	0.43
1:A:860:ALA:N	3:A:1248:B12:O8R	2.46	0.43
1:A:784:VAL:HA	1:A:785:PRO:HD3	1.80	0.43
1:A:1200:VAL:HG12	1:A:1214:VAL:HG21	1.99	0.43
1:A:1085:LYS:O	1:A:1088:VAL:HG12	2.18	0.43
1:A:1106:ARG:NE	1:A:1111:TYR:CZ	2.84	0.43
1:A:978:LYS:HE2	1:A:982:ASP:OD1	2.18	0.43
1:A:674:PHE:O	1:A:678:ASP:HB2	2.18	0.43
1:A:897:LYS:O	1:A:898:LYS:O	2.37	0.43
1:A:1112:TRP:CH2	1:A:1183:SER:HB3	2.54	0.42
1:A:1183:SER:O	1:A:1184:HIS:C	2.57	0.42
1:A:1202:ASP:O	1:A:1206:ARG:CG	2.64	0.42
1:A:850:TYR:CE2	1:A:852:GLY:HA3	2.53	0.42
1:A:707:LEU:HB2	1:A:713:MET:SD	2.59	0.42
1:A:761:ILE:O	1:A:765:ILE:HD12	2.19	0.42
1:A:1145:HIS:CD2	1:A:1148:LYS:HZ1	2.38	0.42
1:A:1188:LYS:HD2	1:A:1190:TYR:CZ	2.55	0.42
1:A:782:VAL:O	1:A:783:MET:HB2	2.19	0.42
1:A:826:PHE:HD1	1:A:828:ILE:CD1	2.31	0.42
1:A:983:MET:CE	1:A:1084:ASN:HD22	2.32	0.42
1:A:941:LEU:O	1:A:942:ARG:C	2.55	0.42
1:A:1010:ASP:OD1	1:A:1027:HIS:HA	2.20	0.42
1:A:715:LEU:HB3	1:A:716:PRO:HD3	2.02	0.42
1:A:689:PRO:HG2	1:A:735:PHE:O	2.19	0.42
1:A:963:ILE:HG23	1:A:964:LEU:N	2.34	0.42
1:A:1128:GLU:HA	1:A:1130:TYR:CZ	2.54	0.42
1:A:1064:THR:HG22	1:A:1163:MET:SD	2.60	0.42
1:A:1135:PRO:O	1:A:1180:TRP:N	2.40	0.42
1:A:1031:GLN:OE1	1:A:1041:TYR:CD2	2.73	0.42
1:A:1210:SER:O	1:A:1211:VAL:C	2.58	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:895:GLY:O	1:A:897:LYS:NZ	2.46	0.41
1:A:1194:GLN:HB3	1:A:1223:GLY:HA3	2.02	0.41
1:A:948:THR:HB	1:A:949:PRO:CD	2.41	0.41
1:A:1018:THR:O	1:A:1020:THR:HG23	2.20	0.41
1:A:659:VAL:HG22	1:A:682:ALA:HB2	2.01	0.41
1:A:697:LEU:HD21	1:A:728:ALA:HB3	2.02	0.41
1:A:733:GLU:N	1:A:734:PRO:HD2	2.35	0.41
1:A:929:LEU:HD23	1:A:929:LEU:HA	1.85	0.41
1:A:929:LEU:CD2	1:A:1003:PRO:HD3	2.50	0.41
1:A:1018:THR:O	1:A:1019:ARG:C	2.58	0.41
1:A:761:ILE:HG22	1:A:762:GLY:N	2.35	0.41
1:A:1047:VAL:HA	1:A:1056:ASP:OD2	2.20	0.41
1:A:1122:ASN:O	1:A:1126:ILE:HG12	2.20	0.41
1:A:733:GLU:HB2	1:A:734:PRO:CD	2.50	0.41
1:A:805:GLY:HA3	1:A:814:MET:CE	2.50	0.41
1:A:917:ASP:OD2	1:A:917:ASP:C	2.58	0.41
1:A:944:TYR:HA	1:A:1122:ASN:OD1	2.21	0.41
1:A:1045:ASP:C	1:A:1047:VAL:H	2.23	0.41
1:A:1201:GLU:O	1:A:1204:ALA:CB	2.66	0.41
1:A:897:LYS:HA	1:A:900:ARG:NH2	2.28	0.41
1:A:912:ASN:CG	1:A:912:ASN:O	2.58	0.41
1:A:942:ARG:O	1:A:945:ILE:HG13	2.21	0.41
1:A:1147:GLU:C	1:A:1149:ALA:N	2.74	0.41
1:A:749:MET:HB3	1:A:777:ILE:CD1	2.51	0.41
1:A:794:LYS:HE2	1:A:826:PHE:CZ	2.55	0.41
1:A:806:LEU:HD12	1:A:806:LEU:HA	1.87	0.41
1:A:1157:VAL:O	1:A:1158:GLU:C	2.59	0.41
1:A:1148:LYS:CE	1:A:1169:PHE:O	2.69	0.41
1:A:1150:THR:O	1:A:1153:GLU:HB3	2.21	0.41
1:A:749:MET:HE3	1:A:800:LEU:CD1	2.51	0.41
1:A:749:MET:HB3	1:A:777:ILE:HD13	2.03	0.41
1:A:1078:ALA:C	1:A:1080:HIS:N	2.65	0.40
1:A:880:VAL:CG2	1:A:881:ALA:N	2.84	0.40
1:A:960:TYR:CG	1:A:961:PRO:HA	2.57	0.40
1:A:906:LEU:O	1:A:910:ARG:HB2	2.21	0.40
1:A:1002:PHE:HB2	1:A:1013:ILE:HG23	2.03	0.40
1:A:847:GLU:CD	1:A:887:TYR:HD1	2.24	0.40
1:A:755:LYS:HB3	1:A:813:GLU:CD	2.42	0.40
1:A:811:LEU:HD12	1:A:841:HIS:CE1	2.57	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	575/577 (100%)	445 (77%)	111 (19%)	19 (3%)	4	39

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	754	VAL
1	A	1041	TYR
1	A	1043	LEU
1	A	1078	ALA
1	A	1079	GLN
1	A	848	GLN
1	A	1017	GLU
1	A	1068	GLU
1	A	652	ALA
1	A	899	PRO
1	A	1110	VAL
1	A	727	GLN
1	A	873	ASP
1	A	898	LYS
1	A	1046	PHE
1	A	1111	TYR
1	A	828	ILE
1	A	1049	PRO
1	A	939	GLU

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	481/481 (100%)	448 (93%)	33 (7%)	18	55

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	660	ASN
1	A	678	ASP
1	A	715	LEU
1	A	740	LYS
1	A	741	GLU
1	A	751	ILE
1	A	754	VAL
1	A	800	LEU
1	A	807	ILE
1	A	808	THR
1	A	828	ILE
1	A	836	THR
1	A	842	THR
1	A	858	GLN
1	A	859	ASN
1	A	870	LEU
1	A	873	ASP
1	A	893	GLN
1	A	900	ARG
1	A	904	VAL
1	A	939	GLU
1	A	962	ARG
1	A	978	LYS
1	A	979	ASP
1	A	1052	SER
1	A	1064	THR
1	A	1082	ASP
1	A	1099	PHE
1	A	1144	GLU
1	A	1152	TRP
1	A	1185	PRO
1	A	1197	ARG
1	A	1221	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	660	ASN
1	A	702	ASN
1	A	727	GLN
1	A	771	GLN
1	A	849	ASN
1	A	858	GLN
1	A	859	ASN
1	A	893	GLN
1	A	943	ASN
1	A	974	GLN
1	A	981	ASN
1	A	1024	ASN
1	A	1027	HIS
1	A	1084	ASN
1	A	1117	ASN
1	A	1129	ASN
1	A	1131	GLN
1	A	1145	HIS
1	A	1184	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](#)

2 ligands 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	$\# Z > 2$
3	B12	A	1248	-	73,101,101	4.74	41 (56%)	111,166,166	3.50	51 (45%)
2	SO4	A	249	-	4,4,4	1.29	0	6,6,6	0.59	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	B12	A	1248	-	1/1/36/38	0/51/223/223	0/3/11/11
2	SO4	A	249	-	-	0/0/0/0	0/0/0/0

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1248	B12	C56-C57	-13.79	1.24	1.51
3	A	1248	B12	C60-C61	-7.36	1.31	1.51
3	A	1248	B12	O3-C2P	-6.46	1.27	1.45
3	A	1248	B12	C5M-C5B	-6.41	1.38	1.51
3	A	1248	B12	C61-N62	-4.52	1.17	1.32
3	A	1248	B12	C32-N33	-3.02	1.22	1.32
3	A	1248	B12	C43-N45	-2.96	1.23	1.32
3	A	1248	B12	C41-C42	-2.75	1.43	1.52
3	A	1248	B12	C50-N52	-2.60	1.24	1.32
3	A	1248	B12	C1-C19	-2.58	1.49	1.55
3	A	1248	B12	P-O5	-2.51	1.42	1.55
3	A	1248	B12	C2-C3	-2.43	1.54	1.58
3	A	1248	B12	O8R-C5R	-2.37	1.32	1.42
3	A	1248	B12	C42-C43	-2.23	1.42	1.51
3	A	1248	B12	C4B-C5B	2.24	1.43	1.37
3	A	1248	B12	C1P-C2P	2.31	1.57	1.51
3	A	1248	B12	C1-C2	2.43	1.64	1.58
3	A	1248	B12	C37-C7	2.81	1.65	1.56
3	A	1248	B12	C31-C32	3.15	1.63	1.51
3	A	1248	B12	C30-C3	3.67	1.59	1.54
3	A	1248	B12	C26-C2	3.98	1.65	1.55
3	A	1248	B12	O6R-C1R	4.32	1.47	1.41
3	A	1248	B12	O6R-C4R	4.51	1.55	1.45
3	A	1248	B12	C6M-C6B	4.56	1.60	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1248	B12	C4B-C9B	4.60	1.48	1.41
3	A	1248	B12	C5R-C4R	5.06	1.69	1.51
3	A	1248	B12	C49-C50	5.19	1.72	1.51
3	A	1248	B12	C8B-C9B	5.36	1.51	1.40
3	A	1248	B12	C25-C2	5.36	1.65	1.54
3	A	1248	B12	C55-C17	5.51	1.68	1.54
3	A	1248	B12	C37-C38	5.91	1.71	1.51
3	A	1248	B12	C7B-C6B	6.58	1.55	1.37
3	A	1248	B12	C6B-C5B	6.67	1.57	1.41
3	A	1248	B12	C7B-C8B	7.04	1.55	1.40
3	A	1248	B12	C48-C13	7.09	1.65	1.54
3	A	1248	B12	C20-C1	7.75	1.69	1.53
3	A	1248	B12	C18-C19	8.27	1.72	1.53
3	A	1248	B12	C26-C27	8.43	1.80	1.51
3	A	1248	B12	C55-C56	9.54	1.75	1.53
3	A	1248	B12	C60-C18	11.65	1.80	1.54
3	A	1248	B12	C53-C15	16.13	1.88	1.51

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1248	B12	C4R-O6R-C1R	-6.78	102.55	109.77
3	A	1248	B12	C60-C61-N62	-6.49	100.17	116.21
3	A	1248	B12	C20-C1-C19	-6.45	103.09	109.34
3	A	1248	B12	O28-C27-C26	-5.23	105.26	121.98
3	A	1248	B12	C60-C18-C17	-4.62	107.29	115.73
3	A	1248	B12	O3-C2P-C1P	-4.44	97.91	106.91
3	A	1248	B12	C5M-C5B-C4B	-4.44	109.21	120.34
3	A	1248	B12	C16-C15-C14	-4.43	116.62	124.00
3	A	1248	B12	C54-C17-C55	-4.17	102.40	109.23
3	A	1248	B12	C35-C5-C4	-4.05	112.57	117.85
3	A	1248	B12	O51-C50-N52	-3.76	112.11	122.47
3	A	1248	B12	C5-C6-N22	-3.49	118.37	124.92
3	A	1248	B12	C56-C57-N59	-3.36	110.69	116.49
3	A	1248	B12	C55-C56-C57	-3.26	104.60	111.01
3	A	1248	B12	C47-C12-C46	-3.05	102.91	109.66
3	A	1248	B12	C48-C49-C50	-3.03	102.67	112.65
3	A	1248	B12	O34-C32-N33	-2.54	115.46	122.47
3	A	1248	B12	C26-C2-C3	-2.53	102.72	107.51
3	A	1248	B12	C53-C15-C14	-2.52	114.56	117.85
3	A	1248	B12	O44-C43-N45	-2.36	115.95	122.47
3	A	1248	B12	C9B-C8B-N1B	-2.24	101.20	108.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1248	B12	C2P-C1P-N59	-2.09	109.96	112.96
3	A	1248	B12	C9-C10-C11	-2.02	126.01	131.90
3	A	1248	B12	O28-C27-N29	2.27	128.73	122.47
3	A	1248	B12	O58-C57-N59	2.66	128.04	122.97
3	A	1248	B12	C26-C27-N29	2.87	125.92	116.55
3	A	1248	B12	C1-C2-C3	2.97	105.41	101.60
3	A	1248	B12	C2-C26-C27	2.97	123.65	115.29
3	A	1248	B12	C31-C32-N33	2.98	126.14	116.55
3	A	1248	B12	O5-P-O3	3.02	118.61	106.49
3	A	1248	B12	C49-C50-N52	3.13	126.62	116.55
3	A	1248	B12	C19-C1-N21	3.17	105.33	102.16
3	A	1248	B12	C30-C3-C2	3.23	125.23	119.07
3	A	1248	B12	C42-C43-N45	3.37	127.38	116.55
3	A	1248	B12	C41-C8-C7	3.63	123.98	114.16
3	A	1248	B12	C8-C7-C6	3.75	107.37	101.04
3	A	1248	B12	C5M-C5B-C6B	4.48	130.12	120.72
3	A	1248	B12	C54-C17-C18	4.67	120.20	112.94
3	A	1248	B12	O63-C61-N62	4.91	135.99	122.47
3	A	1248	B12	C60-C18-C19	5.09	128.17	114.65
3	A	1248	B12	C7-C37-C38	5.25	130.20	114.25
3	A	1248	B12	O3-C2P-C3P	5.90	132.88	108.73
3	A	1248	B12	C49-C48-C13	6.35	133.29	114.76
3	A	1248	B12	C2-C1-C19	6.37	128.84	118.60
3	A	1248	B12	C56-C55-C17	6.81	129.08	115.56
3	A	1248	B12	C1-C19-C18	7.09	133.72	121.90
3	A	1248	B12	C53-C15-C16	7.77	127.99	117.85
3	A	1248	B12	C3P-C2P-C1P	8.17	127.75	111.40
3	A	1248	B12	C17-C18-C19	9.13	116.61	102.38
3	A	1248	B12	O6R-C4R-C5R	9.78	130.09	109.16
3	A	1248	B12	C18-C60-C61	12.55	146.59	113.98

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	1248	B12	C2P

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1248	B12	20	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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.