



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:54 PM GMT

PDB ID : 1R31  
Title : HMG-CoA reductase from *Pseudomonas mevalonii* complexed with HMG-CoA  
Authors : Watson, J.M.; Steussy, C.N.; Burgner, J.W.; Lawrence, C.M.; Tabernero, L.;  
Rodwell, V.W.; Stauffacher, C.V.  
Deposited on : 2003-09-30  
Resolution : 2.10 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](mailto: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.

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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

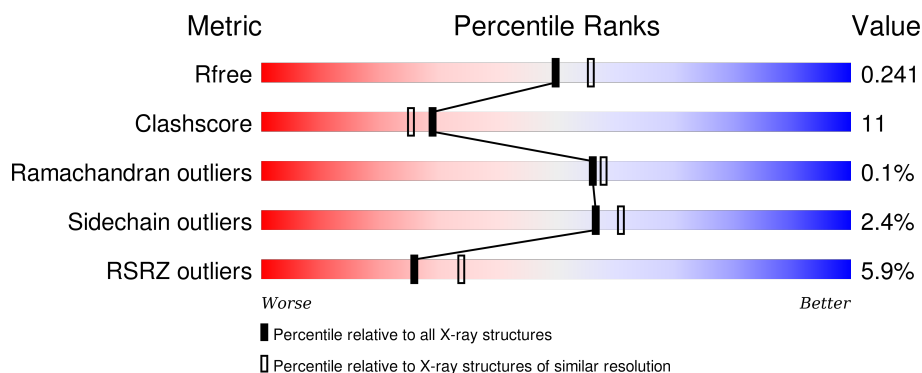
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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  $\geq 3$ , 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	428	<div> <div>4%</div> <div>70% 16% • 12%</div> </div>
1	B	428	<div> <div>7%</div> <div>69% 17% • 12%</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
4	MEV	B	1004	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5970 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 3-hydroxy-3-methylglutaryl-coenzyme A reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	376	Total	C	N	O	S	0	0	0
			2791	1748	504	525	14			
1	B	375	Total	C	N	O	S	0	0	0
			2782	1743	502	523	14			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



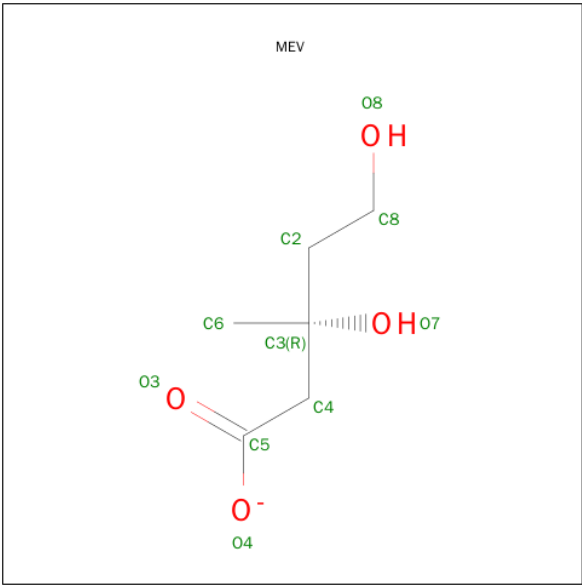
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is COENZYME A (three-letter code: COA) (formula: C<sub>21</sub>H<sub>36</sub>N<sub>7</sub>O<sub>16</sub>P<sub>3</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	A	1	48	21	7	16	3	0	0

- Molecule 4 is (R)-MEVALONATE (three-letter code: MEV) (formula: C<sub>6</sub>H<sub>11</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	10	6	4	0	0
4	B	1	10	6	4	0	0

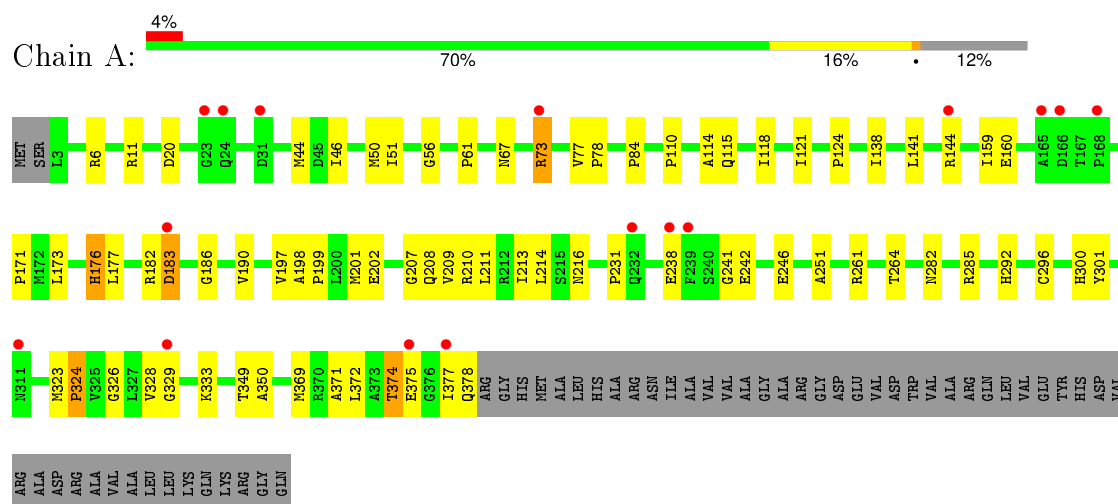
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	188	Total 188	O 188	0	0
5	B	131	Total 131	O 131	0	0

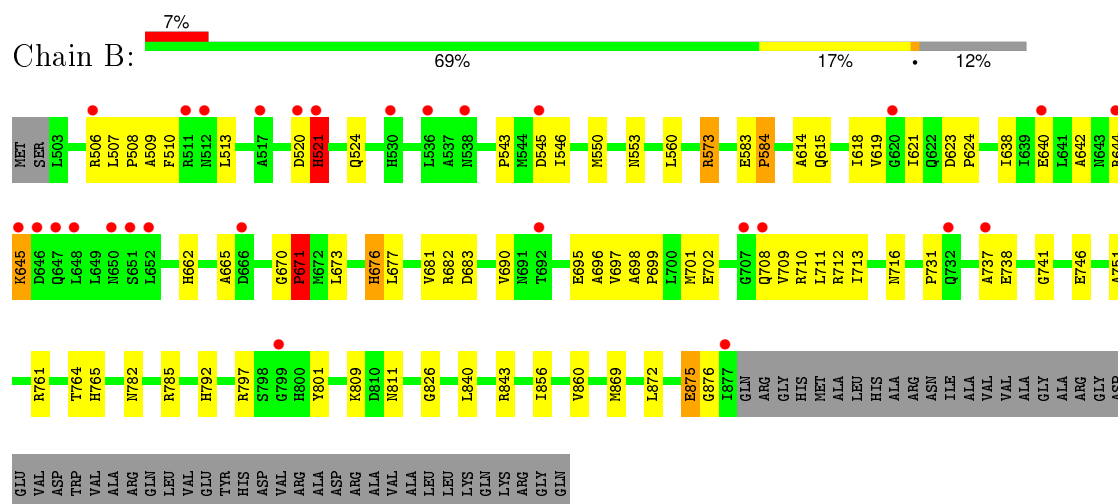
### 3 Residue-property plots

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: 3-hydroxy-3-methylglutaryl-coenzyme A reductase



- Molecule 1: 3-hydroxy-3-methylglutaryl-coenzyme A reductase



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 41 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	227.37Å 227.37Å 227.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.10 8.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	93.7 (30.00-2.10) 93.7 (8.00-2.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.51 (at 2.11Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.220 , 0.241 0.220 , 0.241	Depositor DCC
$R_{free}$ test set	4332 reflections (8.12%)	DCC
Wilson B-factor (Å <sup>2</sup> )	31.4	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.45 , 63.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 53329 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5970	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.84% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: COA, MEV, 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	0.64	2/2834 (0.1%)	0.66	1/3855 (0.0%)
1	B	0.58	3/2825 (0.1%)	0.73	7/3843 (0.2%)
All	All	0.61	5/5659 (0.1%)	0.70	8/7698 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	584	PRO	N-CD	-19.39	1.20	1.47
1	A	324	PRO	N-CD	-18.38	1.22	1.47
1	A	84	PRO	N-CD	-17.96	1.22	1.47
1	B	875	GLU	C-N	-7.31	1.19	1.33
1	B	521	HIS	C-N	6.89	1.49	1.34

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	671	PRO	CA-N-CD	-12.12	94.53	111.50
1	B	619	VAL	CB-CA-C	-7.64	96.88	111.40
1	B	670	GLY	C-N-CD	-5.76	107.93	120.60
1	B	584	PRO	N-CD-CG	5.58	111.57	103.20
1	B	619	VAL	N-CA-C	5.39	125.57	111.00
1	A	374	THR	C-N-CA	-5.36	108.29	121.70
1	B	619	VAL	C-N-CA	-5.25	111.27	122.30
1	B	521	HIS	O-C-N	5.10	130.86	122.70

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2791	0	2833	69	0
1	B	2782	0	2824	69	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	48	0	32	2	0
4	A	10	0	10	0	0
4	B	10	0	10	0	0
5	A	188	0	0	4	0
5	B	131	0	0	5	0
All	All	5970	0	5709	130	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 11.

All (130) 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:374:THR:HG23	1:A:375:GLU:N	1.61	1.07
1:A:374:THR:CG2	1:A:375:GLU:H	1.70	1.03
1:A:374:THR:HG23	1:A:375:GLU:H	0.86	1.01
1:B:614:ALA:HB2	1:B:690:VAL:HG13	1.46	0.93
1:B:618:ILE:HD12	1:B:673:LEU:HD22	1.65	0.79
1:B:645:LYS:HE3	1:B:645:LYS:HA	1.65	0.79
1:A:211:LEU:HD22	1:B:872:LEU:HB3	1.67	0.76
1:A:374:THR:CG2	1:A:375:GLU:N	2.34	0.75
1:A:264:THR:HA	1:B:716:ASN:HD22	1.52	0.74
1:A:61:PRO:HG2	1:B:550:MET:HE2	1.68	0.74
1:B:746:GLU:OE1	1:B:809:LYS:NZ	2.22	0.72
1:B:573:ARG:HH11	1:B:573:ARG:HB3	1.54	0.72
1:B:702:GLU:HG3	1:B:709:VAL:HG23	1.72	0.72
1:A:73:ARG:HB3	1:A:73:ARG:HH11	1.54	0.72
1:A:372:LEU:HB3	1:B:711:LEU:HD22	1.73	0.71
1:B:665:ALA:O	1:B:671:PRO:HD3	1.91	0.71
1:A:238:GLU:CD	1:A:375:GLU:OE2	2.30	0.70
1:B:614:ALA:CB	1:B:690:VAL:HG13	2.20	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:695:GLU:HG2	1:B:712:ARG:HD2	1.75	0.68
1:B:506:ARG:O	1:B:508:PRO:HD3	1.92	0.68
1:A:183:ASP:OD1	1:A:349:THR:HA	1.95	0.67
1:A:44:MET:CE	1:A:56:GLY:HA2	2.25	0.67
1:A:216:ASN:HD22	1:B:764:THR:HA	1.60	0.65
1:A:372:LEU:HD23	1:A:377:ILE:HB	1.79	0.65
1:A:124:PRO:HG2	1:A:171:PRO:HB2	1.78	0.65
1:A:118:ILE:HD12	1:A:173:LEU:HD22	1.77	0.64
1:A:183:ASP:OD1	1:A:350:ALA:N	2.30	0.64
1:A:292:HIS:HD2	1:A:301:TYR:OH	1.80	0.64
1:B:792:HIS:HD2	1:B:801:TYR:OH	1.81	0.63
1:B:731:PRO:O	1:B:741:GLY:HA3	2.00	0.62
1:B:665:ALA:O	1:B:671:PRO:HG3	1.99	0.61
1:A:202:GLU:HG3	1:A:209:VAL:HG23	1.81	0.60
1:B:583:GLU:CG	1:B:584:PRO:HD2	2.32	0.59
1:A:51:ILE:HG22	1:B:583:GLU:O	2.03	0.59
1:A:372:LEU:CD2	1:A:377:ILE:HB	2.32	0.59
1:A:73:ARG:HB3	1:A:73:ARG:NH1	2.18	0.58
1:A:282:ASN:HD21	1:A:326:GLY:H	1.52	0.57
1:A:238:GLU:H	1:A:238:GLU:CD	2.07	0.57
1:A:197:VAL:HG12	1:A:201:MET:HG3	1.86	0.57
1:B:640:GLU:O	1:B:644:ARG:HG2	2.04	0.57
1:B:573:ARG:NH1	1:B:573:ARG:HB3	2.18	0.57
1:A:44:MET:HE3	1:A:56:GLY:HA2	1.85	0.56
1:A:238:GLU:OE1	1:A:375:GLU:OE2	2.24	0.56
1:B:737:ALA:HB3	1:B:738:GLU:OE2	2.05	0.56
1:B:697:VAL:HG12	1:B:701:MET:HG3	1.86	0.56
1:A:377:ILE:HG23	1:A:378:GLN:HG2	1.87	0.55
1:B:543:PRO:HG2	1:B:546:ILE:CG1	2.37	0.54
1:B:583:GLU:HG3	1:B:584:PRO:HD2	1.90	0.54
1:B:521:HIS:O	1:B:524:GLN:HB3	2.08	0.54
1:B:573:ARG:CB	1:B:573:ARG:HH11	2.21	0.54
1:A:73:ARG:HH11	1:A:73:ARG:CB	2.20	0.53
1:A:377:ILE:HG22	5:B:283:HOH:O	2.09	0.53
1:B:782:ASN:HD21	1:B:826:GLY:H	1.56	0.53
1:A:198:ALA:HB3	1:A:199:PRO:HD3	1.91	0.53
1:A:44:MET:HE1	1:A:56:GLY:HA2	1.90	0.52
1:B:642:ALA:O	1:B:645:LYS:HB3	2.09	0.52
1:B:614:ALA:HB3	1:B:677:LEU:HB2	1.91	0.52
1:B:681:VAL:O	1:B:682:ARG:HB2	2.10	0.52
3:A:1002:COA:H2A	1:B:553:ASN:OD1	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:840:LEU:HA	1:B:843:ARG:NH1	2.25	0.51
1:A:296:CYS:SG	1:A:296:CYS:O	2.68	0.51
1:A:333:LYS:NZ	5:A:1129:HOH:O	2.43	0.51
1:B:708:GLN:CD	1:B:710:ARG:HH12	2.14	0.51
1:B:698:ALA:HB3	1:B:699:PRO:HD3	1.92	0.51
1:B:546:ILE:O	1:B:550:MET:HG3	2.11	0.51
1:B:618:ILE:CG2	1:B:621:ILE:HD11	2.41	0.51
1:A:61:PRO:CG	1:B:550:MET:HE2	2.41	0.50
1:B:615:GLN:HE21	1:B:676:HIS:HE1	1.59	0.50
1:A:231:PRO:O	1:A:241:GLY:HA3	2.12	0.50
1:A:186:GLY:O	1:A:190:VAL:HG22	2.12	0.50
1:B:509:ALA:O	1:B:513:LEU:HD13	2.11	0.49
1:A:251:ALA:HB3	1:A:369:MET:HE2	1.93	0.49
1:A:114:ALA:HB2	1:A:190:VAL:HB	1.94	0.48
1:A:46:ILE:O	1:A:50:MET:HG3	2.13	0.48
1:B:665:ALA:O	1:B:671:PRO:CD	2.61	0.48
1:B:624:PRO:HG3	1:B:671:PRO:HB3	1.94	0.48
1:B:797:ARG:HG2	1:B:797:ARG:HH11	1.79	0.48
1:A:121:ILE:HG23	1:A:207:GLY:CA	2.42	0.48
1:B:702:GLU:CG	1:B:709:VAL:HG23	2.42	0.48
1:A:6:ARG:HG2	1:A:67:ASN:OD1	2.14	0.48
1:A:371:ALA:HB2	3:A:1002:COA:H71	1.95	0.48
1:B:615:GLN:HE21	1:B:676:HIS:CE1	2.31	0.48
1:A:238:GLU:CD	1:A:238:GLU:N	2.68	0.48
1:A:114:ALA:HB3	1:A:177:LEU:HB2	1.96	0.48
1:B:645:LYS:HD2	1:B:696:ALA:CB	2.44	0.47
1:A:282:ASN:ND2	1:A:326:GLY:H	2.12	0.47
1:B:782:ASN:ND2	1:B:826:GLY:H	2.13	0.46
1:B:765:HIS:HD2	5:B:81:HOH:O	1.96	0.46
1:A:183:ASP:OD1	1:A:349:THR:CA	2.62	0.46
1:B:543:PRO:HG2	1:B:546:ILE:HG12	1.96	0.46
1:A:242:GLU:O	1:A:246:GLU:HG2	2.15	0.46
1:A:115:GLN:HE21	1:A:176:HIS:HE1	1.61	0.46
1:B:665:ALA:O	1:B:671:PRO:CG	2.64	0.46
1:B:875:GLU:O	1:B:876:GLY:C	2.51	0.46
1:A:110:PRO:HD2	1:A:182:ARG:HH21	1.80	0.46
1:B:645:LYS:HE3	1:B:645:LYS:CA	2.43	0.46
1:A:141:LEU:O	1:A:144:ARG:HG2	2.16	0.46
1:B:662:HIS:CE1	5:B:187:HOH:O	2.69	0.45
1:B:621:ILE:HG22	1:B:623:ASP:H	1.82	0.45
1:B:507:LEU:O	1:B:510:PHE:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:ILE:HG12	1:A:177:LEU:HD23	1.98	0.45
1:B:560:LEU:HA	1:B:560:LEU:HD23	1.86	0.45
1:A:115:GLN:HE21	1:A:176:HIS:CE1	2.34	0.45
1:B:583:GLU:HA	1:B:584:PRO:HD3	1.64	0.44
1:A:77:VAL:HA	1:A:78:PRO:HD3	1.88	0.44
1:B:645:LYS:HD2	1:B:696:ALA:HB2	2.00	0.44
1:B:682:ARG:HB3	1:B:683:ASP:H	1.49	0.43
1:A:213:ILE:HG13	1:A:214:LEU:N	2.33	0.43
1:A:371:ALA:O	1:A:374:THR:HG22	2.18	0.43
1:B:506:ARG:NH1	5:B:237:HOH:O	2.47	0.43
1:A:323:MET:N	1:A:324:PRO:HD3	2.33	0.43
1:B:751:ALA:HB3	1:B:869:MET:HE2	2.01	0.43
1:A:328:VAL:HG22	1:A:329:GLY:N	2.34	0.43
1:A:251:ALA:HB3	1:A:369:MET:CE	2.49	0.42
1:A:121:ILE:HG23	1:A:207:GLY:HA2	2.01	0.42
1:B:638:ILE:HG23	1:B:697:VAL:HG11	2.01	0.42
5:A:1188:HOH:O	1:B:713:ILE:HD11	2.19	0.42
1:A:124:PRO:CG	1:A:171:PRO:HB2	2.45	0.42
1:A:160:GLU:HB2	1:A:176:HIS:HB2	2.02	0.42
1:B:856:ILE:O	1:B:860:VAL:HG23	2.20	0.42
1:A:11:ARG:HG3	5:A:1055:HOH:O	2.19	0.42
1:A:264:THR:HG23	1:B:716:ASN:ND2	2.35	0.41
1:B:751:ALA:HB3	1:B:869:MET:CE	2.50	0.41
1:A:208:GLN:CD	1:A:210:ARG:HH12	2.23	0.41
1:A:138:ILE:HG23	1:A:197:VAL:HG11	2.03	0.41
1:A:173:LEU:C	1:A:173:LEU:HD23	2.41	0.41
1:A:300:HIS:HE1	5:A:1015:HOH:O	2.04	0.41
1:A:377:ILE:HD12	1:A:377:ILE:O	2.21	0.40
1:B:507:LEU:CB	1:B:510:PHE:HB2	2.51	0.40
1:B:811:ASN:ND2	5:B:358:HOH:O	2.54	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	374/428 (87%)	359 (96%)	14 (4%)	1 (0%)	46	45
1	B	373/428 (87%)	360 (96%)	13 (4%)	0	100	100
All	All	747/856 (87%)	719 (96%)	27 (4%)	1 (0%)	56	58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	183	ASP

### 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	288/327 (88%)	283 (98%)	5 (2%)	68	74
1	B	287/327 (88%)	278 (97%)	9 (3%)	47	50
All	All	575/654 (88%)	561 (98%)	14 (2%)	57	61

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	ASP
1	A	73	ARG
1	A	176	HIS
1	A	261	ARG
1	A	285	ARG
1	B	520	ASP
1	B	521	HIS
1	B	545	ASP
1	B	573	ARG
1	B	645	LYS
1	B	671	PRO
1	B	676	HIS

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Mol	Chain	Res	Type
1	B	761	ARG
1	B	785	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	147	GLN
1	A	176	HIS
1	A	216	ASN
1	A	265	HIS
1	A	282	ASN
1	A	292	HIS
1	A	300	HIS
1	A	311	ASN
1	A	364	GLN
1	A	378	GLN
1	B	647	GLN
1	B	676	HIS
1	B	688	ASN
1	B	716	ASN
1	B	765	HIS
1	B	782	ASN
1	B	792	HIS
1	B	800	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

5 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	$\# Z  > 2$
2	SO4	A	1000	-	4,4,4	0.26	0	6,6,6	0.13	0
3	COA	A	1002	-	40,50,50	1.53	3 (7%)	50,75,75	1.87	13 (26%)
4	MEV	A	1003	-	5,9,9	1.92	2 (40%)	2,12,12	3.17	1 (50%)
2	SO4	B	1001	-	4,4,4	0.19	0	6,6,6	0.07	0
4	MEV	B	1004	-	5,9,9	2.00	2 (40%)	2,12,12	3.22	1 (50%)

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	SO4	A	1000	-	-	0/0/0/0	0/0/0/0
3	COA	A	1002	-	-	0/44/64/64	0/3/3/3
4	MEV	A	1003	-	-	0/7/9/9	0/0/0/0
2	SO4	B	1001	-	-	0/0/0/0	0/0/0/0
4	MEV	B	1004	-	-	0/7/9/9	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1003	MEV	O8-C8	-2.74	1.27	1.42
4	B	1004	MEV	O8-C8	-2.58	1.28	1.42
3	A	1002	COA	O4B-C4B	2.13	1.49	1.45
3	A	1002	COA	C4A-N3A	2.62	1.39	1.35
4	A	1003	MEV	C6-C3	2.85	1.55	1.52
4	B	1004	MEV	C6-C3	3.27	1.56	1.52
3	A	1002	COA	C2A-N3A	7.78	1.45	1.32

All (15) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1002	COA	N3A-C2A-N1A	-3.14	126.49	128.89
3	A	1002	COA	C2B-C1B-N9A	-3.12	109.53	114.29
3	A	1002	COA	O6A-P2A-O4A	-3.08	97.67	109.62
3	A	1002	COA	OAP-CAP-C9P	-3.01	103.47	110.38
3	A	1002	COA	C1B-N9A-C4A	-2.65	122.94	126.94
3	A	1002	COA	C6P-C5P-N4P	-2.07	112.87	116.46
3	A	1002	COA	O3B-P3B-O7A	2.24	112.69	107.11
3	A	1002	COA	CDP-CBP-CAP	2.32	113.57	109.34
3	A	1002	COA	O3B-C3B-C4B	2.41	119.44	109.99
3	A	1002	COA	C4B-O4B-C1B	2.45	112.41	109.72
3	A	1002	COA	C3P-N4P-C5P	2.63	127.96	122.79
4	A	1003	MEV	O8-C8-C2	4.49	121.38	111.14
4	B	1004	MEV	O8-C8-C2	4.54	121.50	111.14
3	A	1002	COA	C3B-C2B-C1B	4.78	111.46	99.98
3	A	1002	COA	O6A-CCP-CBP	6.17	120.47	110.55

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1002	COA	2	0

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	376/428 (87%)	-0.10	16 (4%)	39 48	21, 32, 58, 84	0
1	B	375/428 (87%)	0.10	28 (7%)	17 23	23, 33, 59, 84	0
All	All	751/856 (87%)	-0.00	44 (5%)	26 34	21, 32, 59, 84	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	646	ASP	5.0
1	A	375	GLU	4.8
1	B	520	ASP	4.6
1	B	530	HIS	4.2
1	A	238	GLU	4.1
1	A	144	ARG	4.0
1	B	648	LEU	3.9
1	A	166	ASP	3.8
1	B	652	LEU	3.7
1	B	511	ARG	3.2
1	B	645	LYS	2.9
1	B	521	HIS	2.8
1	B	666	ASP	2.8
1	B	644	ARG	2.8
1	B	538	ASN	2.8
1	A	311	ASN	2.8
1	B	651	SER	2.8
1	B	647	GLN	2.7
1	B	545	ASP	2.7
1	A	329	GLY	2.7
1	B	737	ALA	2.6
1	A	24	GLN	2.6
1	B	732	GLN	2.6
1	A	73	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	692	THR	2.5
1	B	506	ARG	2.5
1	A	232	GLN	2.5
1	B	517	ALA	2.4
1	A	23	GLY	2.3
1	A	31	ASP	2.3
1	A	168	PRO	2.3
1	B	708	GLN	2.2
1	B	512	ASN	2.2
1	A	239	PHE	2.2
1	B	536	LEU	2.2
1	A	183	ASP	2.2
1	B	707	GLY	2.2
1	A	377	ILE	2.2
1	B	799	GLY	2.1
1	B	640	GLU	2.1
1	A	165	ALA	2.1
1	B	877	ILE	2.0
1	B	650	ASN	2.0
1	B	620	GLY	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
4	MEV	B	1004	10/10	0.84	0.16	3.29	37,41,44,45	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	SO4	A	1000	5/5	0.89	0.21	1.07	44,47,49,49	0
2	SO4	B	1001	5/5	0.95	0.18	0.81	44,45,46,48	0
4	MEV	A	1003	10/10	0.88	0.12	0.59	28,32,40,42	0
3	COA	A	1002	48/48	0.88	0.14	0.43	41,62,79,82	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.