



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 09:41 PM UTC

PDB ID : 1R31 / pdb_00001r31
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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

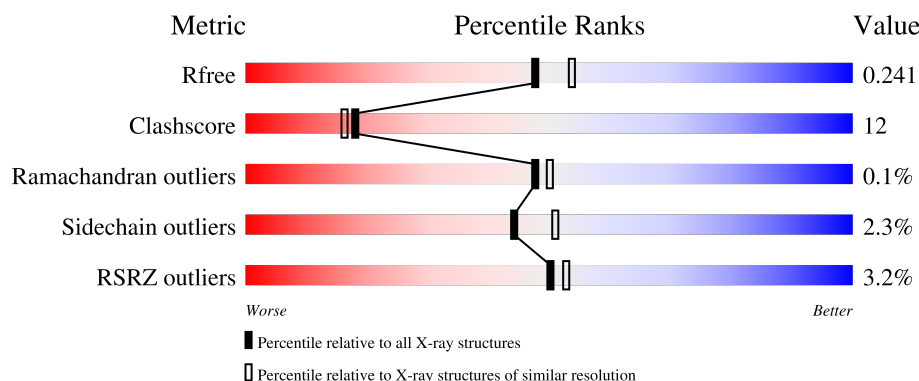
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

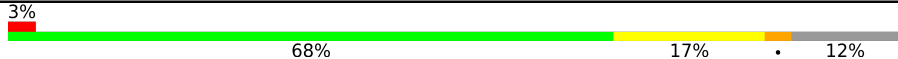
The reported resolution of this entry is 2.10 Å.

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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	
1	B	428	

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 (CCD ID: SO4) (formula: O₄S).



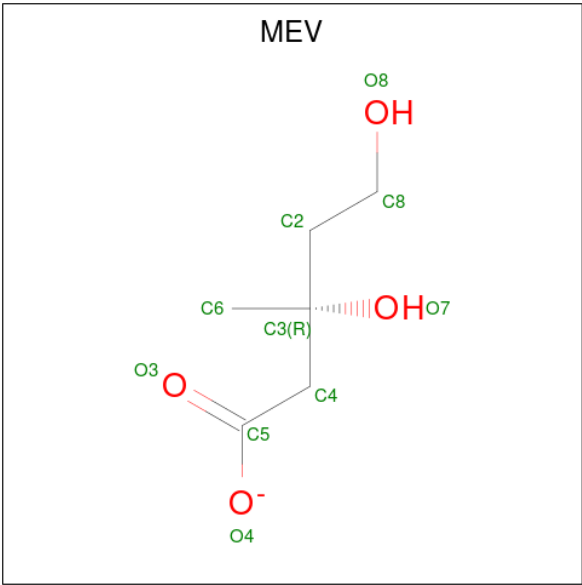
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is COENZYME A (CCD ID: COA) (formula: C₂₁H₃₆N₇O₁₆P₃S).



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

- Molecule 4 is (R)-MEVALONATE (CCD ID: MEV) (formula: C₆H₁₁O₄).



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

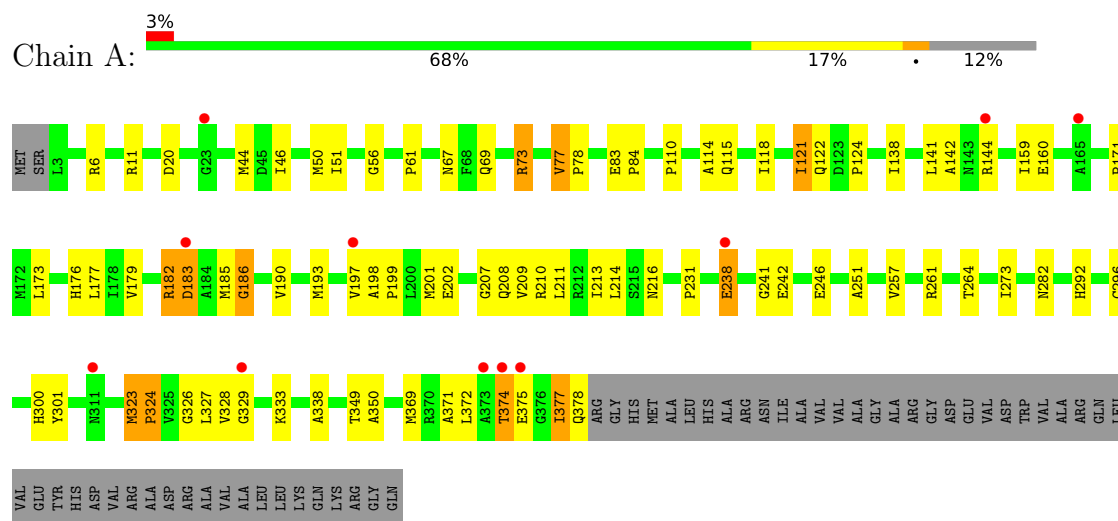
- Molecule 5 is water.

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

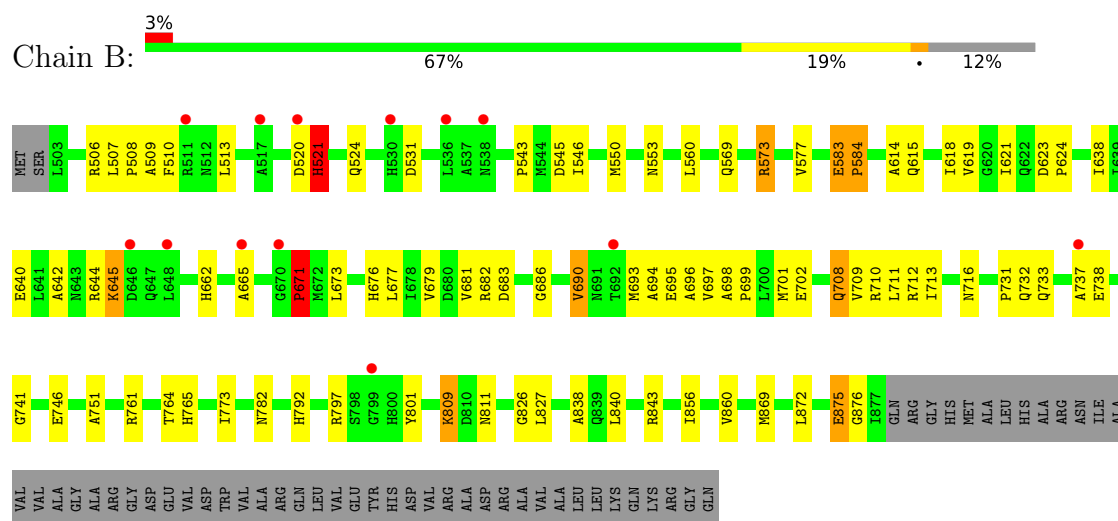
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

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



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4 Data and refinement statistics

Property	Value	Source
Space group	I 41 3 2	Depositor
Cell constants a, b, c, α , β , γ	227.37Å 227.37Å 227.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.10 30.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	93.7 (30.00-2.10) 91.7 (30.00-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	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%)	wwPDB-VP
Wilson B-factor (Å ²)	31.4	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.45 , 59.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5970	wwPDB-VP
Average B, all atoms (Å ²)	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.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: SO4, COA, MEV

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	0.75	4/2834 (0.1%)	1.04	17/3855 (0.4%)
1	B	0.75	7/2825 (0.2%)	1.06	22/3843 (0.6%)
All	All	0.75	11/5659 (0.2%)	1.05	39/7698 (0.5%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	584	PRO	N-CD	-19.32	1.20	1.47
1	A	324	PRO	N-CD	-18.31	1.22	1.47
1	A	84	PRO	N-CD	-17.89	1.22	1.47
1	B	521	HIS	C-N	12.65	1.49	1.34
1	B	875	GLU	C-N	-9.33	1.19	1.33
1	B	809	LYS	C-N	6.55	1.42	1.33
1	A	323	MET	N-CA	6.10	1.52	1.46
1	B	583	GLU	C-N	5.67	1.40	1.33
1	B	694	ALA	CA-C	-5.63	1.45	1.52
1	B	584	PRO	CA-C	-5.31	1.47	1.52
1	A	377	ILE	C-N	-5.22	1.26	1.33

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	671	PRO	CA-N-CD	-12.48	94.53	112.00
1	A	122	GLN	OE1-CD-NE2	-9.89	112.71	122.60
1	B	708	GLN	OE1-CD-NE2	-9.73	112.87	122.60
1	B	569	GLN	OE1-CD-NE2	-9.48	113.12	122.60
1	A	69	GLN	OE1-CD-NE2	-9.34	113.26	122.60
1	B	732	GLN	OE1-CD-NE2	-9.34	113.26	122.60
1	B	733	GLN	OE1-CD-NE2	-9.19	113.41	122.60
1	B	619	VAL	CB-CA-C	-8.79	96.88	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	521	HIS	O-C-N	8.20	130.86	122.08
1	A	374	THR	CA-C-N	-8.07	108.29	122.26
1	A	374	THR	C-N-CA	-8.07	108.29	122.26
1	B	619	VAL	N-CA-C	7.80	125.57	109.34
1	A	257	VAL	CB-CA-C	-7.72	104.27	111.06
1	A	122	GLN	CG-CD-NE2	7.57	127.75	116.40
1	B	708	GLN	CG-CD-NE2	7.05	126.98	116.40
1	A	69	GLN	CG-CD-NE2	6.81	126.61	116.40
1	A	182	ARG	CB-CA-C	-6.63	107.06	116.34
1	B	679	VAL	N-CA-C	6.54	117.26	108.11
1	A	273	ILE	N-CA-C	6.52	117.15	110.82
1	A	83	GLU	N-CA-C	6.46	117.79	109.72
1	A	179	VAL	N-CA-C	6.40	117.07	108.11
1	B	733	GLN	CG-CD-NE2	6.22	125.74	116.40
1	B	569	GLN	CG-CD-NE2	6.21	125.72	116.40
1	B	577	VAL	N-CA-C	6.12	117.38	107.71
1	A	122	GLN	CA-C-N	-6.10	115.66	123.16
1	A	122	GLN	C-N-CA	-6.10	115.66	123.16
1	B	732	GLN	CG-CD-NE2	6.05	125.48	116.40
1	A	186	GLY	N-CA-C	5.96	121.57	114.48
1	B	773	ILE	N-CA-C	5.93	116.57	110.82
1	A	77	VAL	N-CA-C	5.91	117.05	107.71
1	B	827	LEU	N-CA-C	-5.71	106.30	113.72
1	A	327	LEU	N-CA-C	-5.61	106.43	113.72
1	B	584	PRO	N-CD-CG	5.58	111.57	103.20
1	B	686	GLY	N-CA-C	5.54	121.08	114.48
1	B	838	ALA	N-CA-C	-5.41	105.46	111.36
1	B	531	ASP	CB-CA-C	-5.32	102.29	110.81
1	A	338	ALA	N-CA-C	-5.20	105.69	111.36
1	B	619	VAL	CA-C-N	-5.17	111.27	121.41
1	B	619	VAL	C-N-CA	-5.17	111.27	121.41

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	72	0
1	B	2782	0	2825	74	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	5	0
5	B	131	0	0	5	0
All	All	5970	0	5710	138	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 12.

All (138) 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	1.04
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:A:61:PRO:HG2	1:B:550:MET:HE2	1.64	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:44:MET:HE1	1:A:56:GLY:HA2	1.67	0.76
1:A:238:GLU:CD	1:A:375:GLU:OE2	2.30	0.75
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: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:HH11	1:A:73:ARG:HB3	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:B:614:ALA:CB	1:B:690:VAL:HG13	2.20	0.69
1:B:506:ARG:O	1:B:508:PRO:HD3	1.92	0.68
1:B:695:GLU:HG2	1:B:712:ARG:HD2	1.75	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:238:GLU:CD	1:A:238:GLU:H	2.07	0.62
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: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:HB3	1:B:573:ARG:NH1	2.18	0.57
1:A:61:PRO:CG	1:B:550:MET:HE2	2.35	0.56
1:A:238:GLU:OE1	1:A:375:GLU:OE2	2.24	0.56
1:B:708:GLN:CD	1:B:710:ARG:HH12	2.14	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:521:HIS:O	1:B:524:GLN:HB3	2.08	0.54
1:B:583:GLU:HG3	1:B:584:PRO:HD2	1.90	0.54
1:B:573:ARG:HH11	1:B:573:ARG:CB	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:238:GLU:CD	1:A:238:GLU:N	2.68	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:A:296:CYS:O	1:A:296:CYS:SG	2.68	0.51
3:A:1002:COA:H2A	1:B:553:ASN:OD1	2.10	0.51
1:B:840:LEU:HA	1:B:843:ARG:NH1	2.25	0.51
1:A:333:LYS:NZ	5:A:1129:HOH:O	2.43	0.51
1:B:681:VAL:O	1:B:682:ARG:HB2	2.10	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:618:ILE:CG2	1:B:621:ILE:HD11	2.41	0.51
1:A:197:VAL:HG11	1:A:201:MET:HE3	1.94	0.50
1:B:615:GLN:HE21	1:B:676:HIS:HE1	1.59	0.50
1:A:186:GLY:O	1:A:190:VAL:HG22	2.12	0.50
1:A:231:PRO:O	1:A:241:GLY:HA3	2.12	0.50
1:B:509:ALA:O	1:B:513:LEU:HD13	2.11	0.49
1:A:44:MET:HE1	1:A:56:GLY:CA	2.42	0.49
1:B:642:ALA:O	1:B:693:MET:HE2	2.13	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:A:121:ILE:HG23	1:A:207:GLY:CA	2.42	0.48
1:B:624:PRO:HG3	1:B:671:PRO:HB3	1.94	0.48
1:B:665:ALA:O	1:B:671:PRO:CD	2.61	0.48
1:B:797:ARG:HH11	1:B:797:ARG:HG2	1.79	0.48
1:A:6:ARG:HG2	1:A:67:ASN:OD1	2.14	0.48
1:B:702:GLU:CG	1:B:709:VAL:HG23	2.42	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:B:875:GLU:O	1:B:876:GLY:C	2.51	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:A:142:ALA:O	1:A:193:MET:HE2	2.14	0.47
1:B:765:HIS:HD2	5:B:81:HOH:O	1.96	0.46
1:B:782:ASN:ND2	1:B:826:GLY:H	2.13	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:115:GLN:HE21	1:A:176:HIS:HE1	1.61	0.46
1:A:242:GLU:O	1:A:246:GLU:HG2	2.15	0.46
1:B:665:ALA:O	1:B:671:PRO:CG	2.64	0.46
1:A:110:PRO:HD2	1:A:182:ARG:HH21	1.80	0.46
1:A:208:GLN:CD	1:A:210:ARG:HH12	2.23	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:A:173:LEU:C	1:A:173:LEU:HD23	2.41	0.45
1:B:662:HIS:CE1	5:B:187:HOH:O	2.69	0.45
1:B:697:VAL:HG11	1:B:701:MET:HE3	1.98	0.45
1:B:507:LEU:O	1:B:510:PHE:HB2	2.16	0.45
1:B:621:ILE:HG22	1:B:623:ASP:H	1.82	0.45
1:A:159:ILE:HG12	1:A:177:LEU:HD23	1.98	0.45
1:B:560:LEU:HD23	1:B:560:LEU:HA	1.86	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:738:GLU:N	1:B:738:GLU:CD	2.75	0.45
1:A:115:GLN:HE21	1:A:176:HIS:CE1	2.34	0.45
1:A:77:VAL:HA	1:A:78:PRO:HD3	1.88	0.44
1:B:583:GLU:HA	1:B:584:PRO:HD3	1.64	0.44
1:B:645:LYS:HD2	1:B:696:ALA:HB2	2.00	0.44
1:A:185:MET:HE1	5:A:1017:HOH:O	2.18	0.43
1:A:213:ILE:HG13	1:A:214:LEU:N	2.33	0.43
1:B:682:ARG:HB3	1:B:683:ASP:H	1.49	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:673:LEU:C	1:B:673:LEU:HD23	2.43	0.43
1:A:251:ALA:HB3	1:A:369:MET:CE	2.49	0.43
1:A:328:VAL:HG22	1:A:329:GLY:N	2.34	0.43
1:A:121:ILE:HG23	1:A:207:GLY:HA2	2.01	0.42
5:A:1188:HOH:O	1:B:713:ILE:HD11	2.19	0.42
1:B:638:ILE:HG23	1:B:697:VAL:HG11	2.01	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:B:751:ALA:HB3	1:B:869:MET:HE2	2.03	0.41
1:A:138:ILE:HG23	1:A:197:VAL:HG11	2.03	0.41
1:A:251:ALA:HB3	1:A:369:MET:HE2	2.02	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:506:ARG:C	1:B:508:PRO:HD3	2.44	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%)	36	36
1	B	373/428 (87%)	360 (96%)	13 (4%)	0	100	100
All	All	747/856 (87%)	719 (96%)	27 (4%)	1 (0%)	48	50

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%)	53	62
1	B	287/327 (88%)	279 (97%)	8 (3%)	38	43
All	All	575/654 (88%)	562 (98%)	13 (2%)	44	51

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	ASP
1	A	73	ARG
1	A	121	ILE
1	A	238	GLU
1	A	261	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	690	VAL

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Mol	Chain	Res	Type
1	B	761	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	100	ASN
1	A	147	GLN
1	A	150	ASN
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	B	567	ASN
1	B	647	GLN
1	B	650	ASN
1	B	676	HIS
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 oligosaccharides 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$
4	MEV	B	1004	-	8,9,9	1.75	2 (25%)	7,12,12	1.61	1 (14%)
2	SO4	B	1001	-	4,4,4	0.35	0	6,6,6	0.08	0
4	MEV	A	1003	-	8,9,9	1.85	3 (37%)	7,12,12	1.58	1 (14%)
2	SO4	A	1000	-	4,4,4	0.37	0	6,6,6	0.14	0
3	COA	A	1002	-	47,50,50	1.46	5 (10%)	69,75,75	1.51	8 (11%)

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
4	MEV	B	1004	-	-	5/9/9/9	-
4	MEV	A	1003	-	-	3/9/9/9	-
3	COA	A	1002	-	-	7/48/64/64	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1002	COA	C2A-N3A	6.76	1.45	1.33
4	B	1004	MEV	C6-C3	3.22	1.56	1.52
4	A	1003	MEV	O8-C8	-2.92	1.27	1.42
4	A	1003	MEV	C6-C3	2.81	1.55	1.52
4	B	1004	MEV	O8-C8	-2.75	1.28	1.42
3	A	1002	COA	C4A-N3A	2.71	1.39	1.34
3	A	1002	COA	P3B-O3B	2.62	1.64	1.59
3	A	1002	COA	C5A-C6A	2.39	1.47	1.41
3	A	1002	COA	O4B-C4B	2.22	1.49	1.45
4	A	1003	MEV	O4-C5	-2.14	1.23	1.30

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1002	COA	O6A-CCP-CBP	6.17	120.47	110.55
3	A	1002	COA	C3B-C2B-C1B	5.26	111.46	99.89
4	B	1004	MEV	O8-C8-C2	3.62	121.50	111.33
4	A	1003	MEV	O8-C8-C2	3.58	121.38	111.33
3	A	1002	COA	O6A-P2A-O4A	-2.84	97.67	108.94
3	A	1002	COA	CDP-CBP-CAP	2.82	113.57	108.77
3	A	1002	COA	C3P-N4P-C5P	2.76	127.96	122.82
3	A	1002	COA	O3B-C3B-C4B	2.67	119.44	110.03
3	A	1002	COA	C1B-N9A-C8A	2.20	131.98	127.09
3	A	1002	COA	C7P-N8P-C9P	2.02	126.17	122.55

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1002	COA	O4B-C4B-C5B-O5B
3	A	1002	COA	CBP-CCP-O6A-P2A
3	A	1002	COA	C5P-C6P-C7P-N8P
4	A	1003	MEV	C8-C2-C3-C4
4	B	1004	MEV	C3-C2-C8-O8
4	B	1004	MEV	C8-C2-C3-C4
4	B	1004	MEV	C8-C2-C3-C6
3	A	1002	COA	C3B-C4B-C5B-O5B
3	A	1002	COA	P2A-O3A-P1A-O1A
4	A	1003	MEV	C8-C2-C3-C6
4	B	1004	MEV	C3-C4-C5-O4
4	B	1004	MEV	C3-C4-C5-O3
4	A	1003	MEV	C3-C2-C8-O8
3	A	1002	COA	C2B-C3B-O3B-P3B
3	A	1002	COA	P2A-O3A-P1A-O2A

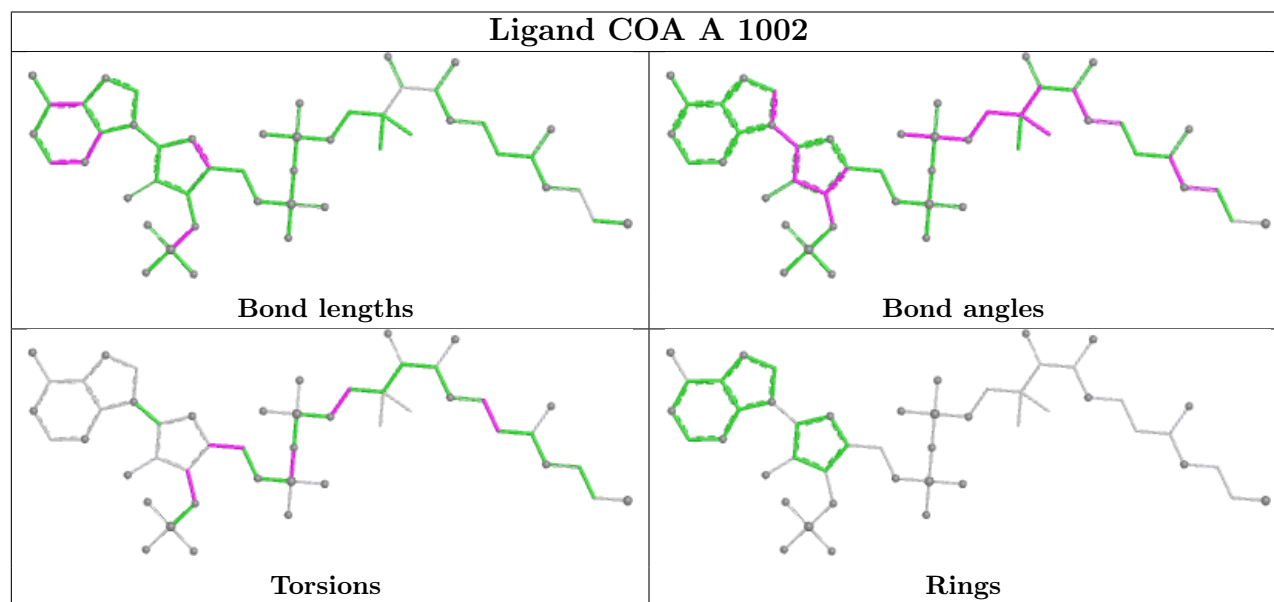
There are no ring outliers.

1 monomer is involved in 2 short contacts:

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

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is

within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	875:GLU	C	876:GLY	N	1.19

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	376/428 (87%)	-0.06	11 (2%) 53 56	21, 32, 58, 84	0
1	B	375/428 (87%)	0.03	13 (3%) 47 49	23, 33, 59, 84	0
All	All	751/856 (87%)	-0.02	24 (3%) 50 53	21, 32, 59, 84	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	375	GLU	5.1
1	A	238	GLU	4.1
1	A	329	GLY	3.3
1	B	648	LEU	3.2
1	A	311	ASN	3.1
1	A	144	ARG	3.0
1	B	670	GLY	3.0
1	B	517	ALA	2.8
1	A	183	ASP	2.7
1	B	799	GLY	2.7
1	A	165	ALA	2.7
1	B	530	HIS	2.6
1	B	665	ALA	2.5
1	B	737	ALA	2.5
1	A	23	GLY	2.5
1	B	538	ASN	2.3
1	A	374	THR	2.3
1	B	536	LEU	2.2
1	B	692	THR	2.2
1	A	197	VAL	2.2
1	B	646	ASP	2.1
1	A	373	ALA	2.1
1	B	520	ASP	2.0
1	B	511	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

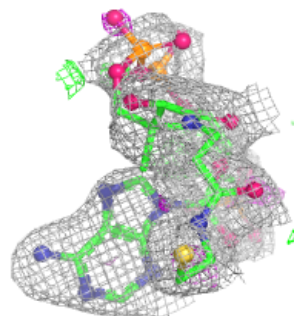
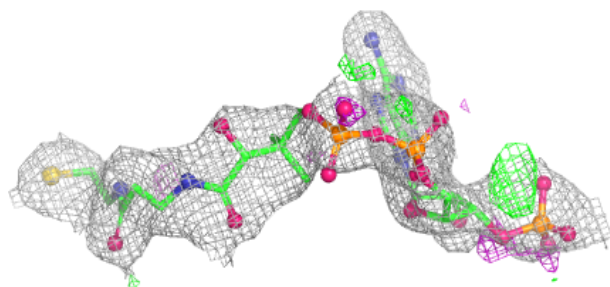
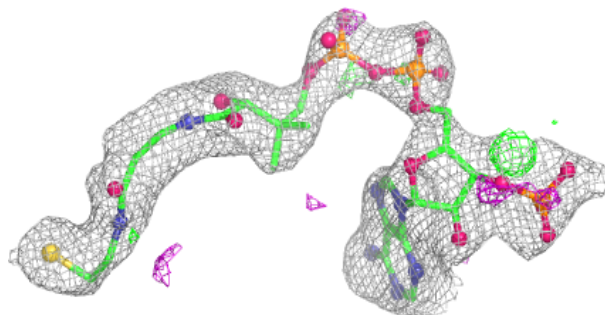
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MEV	B	1004	10/10	0.73	0.12	37,41,44,45	0
4	MEV	A	1003	10/10	0.82	0.09	28,32,40,42	0
2	SO4	A	1000	5/5	0.85	0.17	44,47,49,49	0
3	COA	A	1002	48/48	0.86	0.10	41,62,79,82	0
2	SO4	B	1001	5/5	0.96	0.08	44,45,46,48	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around COA A 1002:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.