



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:29 AM GMT

PDB ID : 2AMT  
Title : Structure of 2C-Methyl-D-Erythritol 2,4-Clycodiphosphate Synthase complexed with a CDP derived fluorescent inhibitor  
Authors : Ramsden, N.L.; Hunter, W.N.  
Deposited on : 2005-08-10  
Resolution : 2.30 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

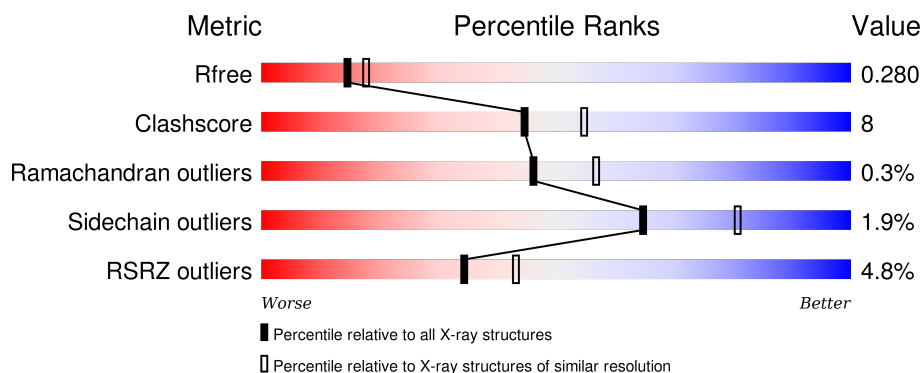
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	159	<div> <div>6%</div> <div>78%18%...</div> </div>
1	B	159	<div> <div>4%</div> <div>79%19%..</div> </div>
1	C	159	<div> <div>5%</div> <div>81%15%..</div> </div>
1	D	159	<div> <div>3%</div> <div>80%16%...</div> </div>
1	E	159	<div> <div>8%</div> <div>82%16%..</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	159	

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	1AA	E	4901	-	-	-	X
4	GPP	A	7903	-	-	-	X
4	GPP	D	8903	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7495 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.

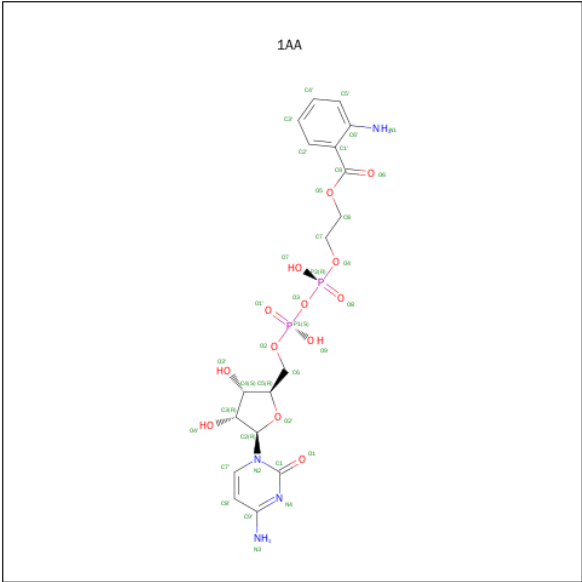
- Molecule 1 is a protein called 2-C-methyl-D-erythritol 2,4-cyclodiphosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	156	Total	C	N	O	S	0	0	0
			1166	741	206	213	6			
1	B	156	Total	C	N	O	S	0	0	0
			1166	741	206	213	6			
1	C	156	Total	C	N	O	S	0	0	0
			1166	741	206	213	6			
1	D	156	Total	C	N	O	S	0	0	0
			1166	741	206	213	6			
1	E	156	Total	C	N	O	S	0	0	0
			1166	741	206	213	6			
1	F	156	Total	C	N	O	S	0	0	0
			1166	741	206	213	6			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

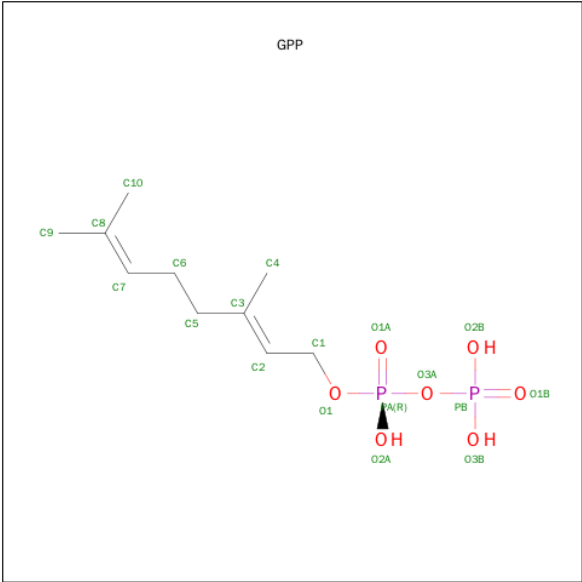
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total	Zn	0	0
			1	1		
2	E	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		
2	F	1	Total	Zn	0	0
			1	1		

- Molecule 3 is 5'-O-([((2-((2-AMINOPHENYL)CARBONYL)OXY)ETHYL)OXY]PHOSPHINATO)OXY)PHOSPHINATO]CYTIDINE (three-letter code: 1AA) (formula: C<sub>18</sub>H<sub>24</sub>N<sub>4</sub>O<sub>13</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	11	0
			37	18	4	13	2		
3	C	1	Total	C	N	O	P	0	0
			37	18	4	13	2		
3	C	1	Total	C	N	O	P	0	0
			37	18	4	13	2		
3	E	1	Total	C	N	O	P	0	0
			37	18	4	13	2		
3	E	1	Total	C	N	O	P	0	0
			37	18	4	13	2		
3	D	1	Total	C	N	O	P	0	0
			37	18	4	13	2		

- Molecule 4 is GERANYL DIPHOSPHATE (three-letter code: GPP) (formula: C<sub>10</sub>H<sub>20</sub>O<sub>7</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	P	0	0
			19	10	7	2		
4	D	1	Total	C	O	P	0	0
			19	10	7	2		

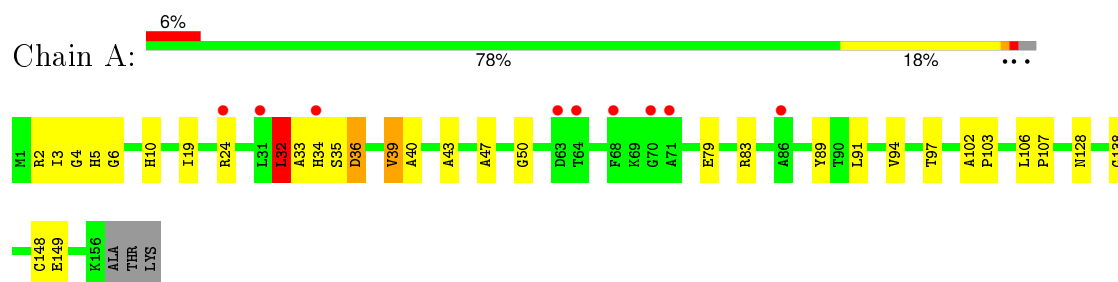
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	37	Total	O	0	0
			37	37		
5	B	37	Total	O	0	0
			37	37		
5	C	43	Total	O	0	0
			43	43		
5	D	44	Total	O	0	0
			44	44		
5	E	36	Total	O	0	0
			36	36		
5	F	36	Total	O	0	0
			36	36		

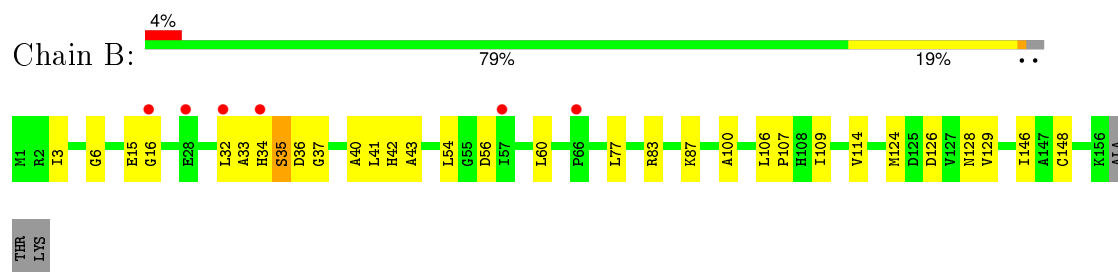
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). 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.

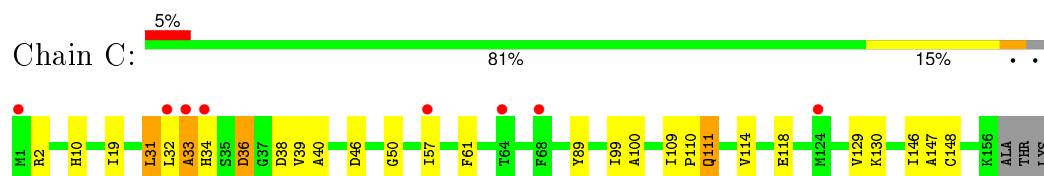
- Molecule 1: 2-C-methyl-D-erythritol 2,4-cyclodiphosphate synthase



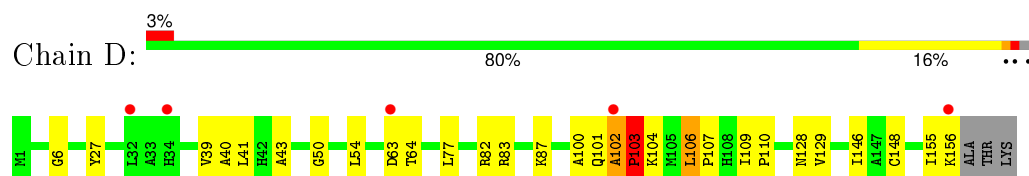
- Molecule 1: 2-C-methyl-D-erythritol 2,4-cyclodiphosphate synthase



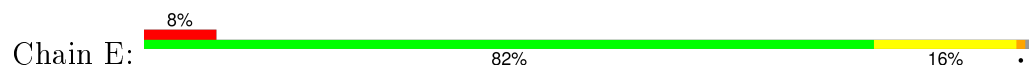
- Molecule 1: 2-C-methyl-D-erythritol 2,4-cyclodiphosphate synthase

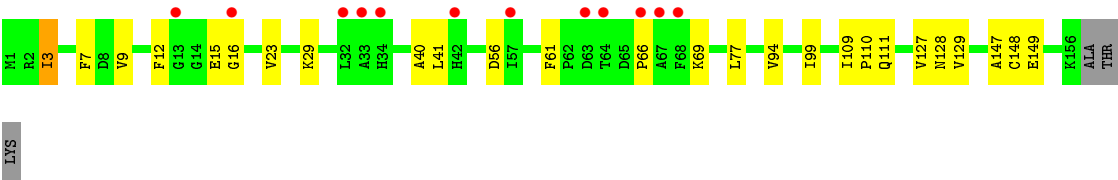


- Molecule 1: 2-C-methyl-D-erythritol 2,4-cyclodiphosphate synthase

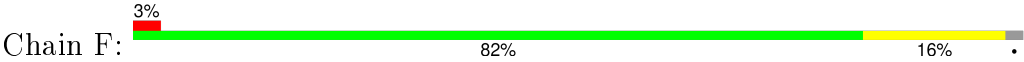


- Molecule 1: 2-C-methyl-D-erythritol 2,4-cyclodiphosphate synthase





● Molecule 1: 2-C-methyl-D-erythritol 2,4-cyclodiphosphate synthase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	54.03Å 115.27Å 87.61Å 90.00° 90.18° 90.00°	Depositor
Resolution (Å)	29.50 – 2.30 29.50 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.5 (29.50-2.30) 98.7 (29.50-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.69 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.243 , 0.280 0.242 , 0.280	Depositor DCC
$R_{free}$ test set	2397 reflections (5.32%)	DCC
Wilson B-factor (Å <sup>2</sup> )	25.6	Xtriage
Anisotropy	1.274	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 21.4	EDS
Estimated twinning fraction	0.138 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 47421 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	7495	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.45% 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: GPP, ZN, 1AA

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	0.31	0/1188	0.50	1/1606 (0.1%)
1	B	0.31	0/1188	0.52	1/1606 (0.1%)
1	C	0.30	0/1188	0.49	2/1606 (0.1%)
1	D	0.43	1/1188 (0.1%)	0.55	1/1606 (0.1%)
1	E	0.35	0/1188	0.46	0/1606
1	F	0.30	0/1188	0.46	0/1606
All	All	0.34	1/7128 (0.0%)	0.50	5/9636 (0.1%)

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
1	B	0	3
1	C	0	2
1	D	0	1
All	All	0	7

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	103	PRO	N-CD	8.94	1.60	1.47

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	36	ASP	CB-CG-OD2	5.73	123.45	118.30
1	A	36	ASP	CB-CG-OD2	5.39	123.15	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	103	PRO	CA-N-CD	-5.37	103.98	111.50
1	C	36	ASP	CB-CG-OD2	5.26	123.04	118.30
1	C	38	ASP	CB-CG-OD2	5.21	122.99	118.30

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	32	LEU	Peptide
1	B	32	LEU	Peptide
1	B	33	ALA	Peptide
1	B	35	SER	Peptide
1	C	31	LEU	Peptide
1	C	34	HIS	Peptide
1	D	101	GLN	Peptide

## 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	1166	0	1182	33	0
1	B	1166	0	1182	24	0
1	C	1166	0	1182	17	0
1	D	1166	0	1182	20	0
1	E	1166	0	1182	18	0
1	F	1166	0	1182	18	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	37	0	22	0	0
3	C	74	0	44	4	0
3	D	37	0	22	0	0
3	E	74	0	44	1	0
4	A	19	0	17	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	19	0	17	1	0
5	A	37	0	0	0	0
5	B	37	0	0	0	0
5	C	43	0	0	1	0
5	D	44	0	0	0	0
5	E	36	0	0	0	0
5	F	36	0	0	1	0
All	All	7495	0	7258	115	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 8.

All (115) 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:D:102:ALA:HB1	1:D:103:PRO:HD3	1.28	1.11
1:D:102:ALA:CB	1:D:103:PRO:CD	2.30	1.10
1:D:102:ALA:HB1	1:D:103:PRO:CD	1.87	1.05
1:D:102:ALA:HB3	1:D:103:PRO:HD2	1.51	0.93
1:E:3:ILE:HD12	1:F:3:ILE:HD11	1.57	0.87
1:B:34:HIS:O	1:B:34:HIS:CD2	2.30	0.84
1:A:102:ALA:HB1	1:A:103:PRO:HA	1.59	0.83
1:D:102:ALA:HB3	1:D:103:PRO:CD	2.09	0.80
1:E:3:ILE:HD12	1:F:3:ILE:CD1	2.12	0.79
1:C:19:ILE:HG21	1:C:36:ASP:HB2	1.63	0.79
1:B:35:SER:HB2	1:B:37:GLY:H	1.47	0.78
1:F:18:ILE:HD12	1:F:20:ILE:HD13	1.69	0.75
1:C:10:HIS:HE1	3:C:3901:1AA:H72	1.53	0.73
1:A:10:HIS:HE1	1:A:34:HIS:NE2	1.90	0.69
1:D:102:ALA:CB	1:D:103:PRO:HD2	2.09	0.66
1:C:19:ILE:CG2	1:C:36:ASP:HB2	2.27	0.63
1:F:12:PHE:HB3	1:F:29:LYS:HD2	1.81	0.62
1:A:40:ALA:HA	1:A:148:CYS:SG	2.43	0.59
1:A:106:LEU:N	1:A:107:PRO:CD	2.66	0.58
1:C:99:ILE:HB	1:C:147:ALA:HB3	1.85	0.58
1:D:100:ALA:HB2	1:D:146:ILE:HG23	1.85	0.57
1:A:10:HIS:CE1	1:A:34:HIS:NE2	2.70	0.57
1:B:60:LEU:HD11	1:B:83:ARG:HD2	1.87	0.56
1:A:102:ALA:CB	1:A:103:PRO:HA	2.29	0.55
1:B:34:HIS:O	1:B:34:HIS:HD2	1.88	0.55
1:D:27:TYR:CD1	1:D:103:PRO:HG3	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:ILE:HD12	1:B:3:ILE:HD13	1.88	0.54
1:C:40:ALA:HA	1:C:148:CYS:SG	2.48	0.54
1:F:40:ALA:HA	1:F:148:CYS:SG	2.48	0.54
1:A:3:ILE:HD12	1:B:3:ILE:CD1	2.38	0.53
1:B:35:SER:OG	3:C:2901:1AA:O5	2.26	0.53
1:E:40:ALA:HA	1:E:148:CYS:SG	2.49	0.52
1:D:40:ALA:HA	1:D:148:CYS:SG	2.50	0.52
1:B:54:LEU:HD21	1:B:87:LYS:HD2	1.91	0.52
1:B:34:HIS:O	1:B:34:HIS:CG	2.62	0.52
1:D:82:ARG:NH2	1:D:83:ARG:HG2	2.24	0.52
1:E:56:ASP:HB3	1:F:130:LYS:HB2	1.92	0.51
1:A:19:ILE:HG12	1:A:24:ARG:HG2	1.93	0.50
1:A:33:ALA:C	1:A:34:HIS:CD2	2.84	0.50
1:B:109:ILE:HG23	1:B:129:VAL:HB	1.94	0.49
1:B:56:ASP:HB3	1:C:130:LYS:HB2	1.94	0.49
1:C:32:LEU:O	1:C:33:ALA:O	2.31	0.49
1:A:6:GLY:HA3	1:A:43:ALA:O	2.13	0.49
1:D:102:ALA:O	1:D:103:PRO:C	2.49	0.49
1:A:33:ALA:HB1	1:A:34:HIS:HD2	1.79	0.48
1:A:34:HIS:CG	1:A:35:SER:H	2.29	0.48
3:E:5901:1AA:H2'	3:E:5901:1AA:H82	1.94	0.48
1:A:19:ILE:HD12	1:A:36:ASP:HB2	1.95	0.48
1:F:23:VAL:HG11	1:F:108:HIS:HB3	1.96	0.48
1:C:114:VAL:O	1:C:118:GLU:HG2	2.14	0.47
1:D:155:ILE:HG22	1:D:156:LYS:HG2	1.95	0.47
1:A:3:ILE:CD1	1:B:3:ILE:HD11	2.45	0.47
1:E:66:PRO:HA	1:E:69:LYS:HG3	1.95	0.47
1:B:41:LEU:HD22	1:B:77:LEU:HD22	1.97	0.47
1:A:5:HIS:NE2	1:A:149:GLU:OE1	2.46	0.46
1:B:106:LEU:HB3	1:B:107:PRO:HD3	1.98	0.46
1:A:50:GLY:HA2	1:B:128:ASN:ND2	2.30	0.46
1:D:6:GLY:HA3	1:D:43:ALA:O	2.15	0.46
1:A:138:GLY:HA3	4:A:7903:GPP:H12	1.97	0.46
1:D:128:ASN:ND2	1:F:50:GLY:HA2	2.31	0.46
1:D:109:ILE:HG23	1:D:129:VAL:HB	1.98	0.46
1:A:106:LEU:HB2	3:C:3901:1AA:O1	2.15	0.45
1:A:3:ILE:HG12	1:A:4:GLY:N	2.31	0.45
1:C:2:ARG:HD3	1:C:89:TYR:CE1	2.51	0.45
1:B:15:GLU:HB2	1:B:16:GLY:H	1.60	0.45
1:C:46:ASP:OD1	1:C:57:ILE:HG22	2.17	0.45
1:D:50:GLY:HA2	1:E:128:ASN:ND2	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:114:VAL:HA	1:B:124:MET:CE	2.46	0.45
1:F:13:GLY:HA3	1:F:32:LEU:HG	1.99	0.45
1:B:40:ALA:HA	1:B:148:CYS:SG	2.57	0.45
1:E:15:GLU:HB3	1:E:16:GLY:H	1.58	0.45
1:B:35:SER:OG	3:C:2901:1AA:C9	2.65	0.45
1:E:41:LEU:HD22	1:E:77:LEU:HD22	1.99	0.44
1:A:106:LEU:HB3	1:A:107:PRO:HD3	2.00	0.44
1:A:33:ALA:CB	1:A:34:HIS:HD2	2.31	0.44
1:F:2:ARG:HD3	1:F:89:TYR:CE1	2.53	0.43
1:F:35:SER:HB2	5:F:6917:HOH:O	2.18	0.43
1:A:2:ARG:HH22	1:B:126:ASP:HA	1.83	0.43
1:C:57:ILE:HG13	1:C:61:PHE:HD2	1.84	0.43
1:E:3:ILE:HD12	1:F:3:ILE:HD13	1.94	0.43
1:E:7:PHE:HD1	1:E:149:GLU:HG2	1.83	0.43
1:E:12:PHE:HB3	1:E:29:LYS:HD3	2.01	0.43
1:A:128:ASN:ND2	1:C:50:GLY:HA2	2.34	0.43
1:B:100:ALA:HB2	1:B:146:ILE:HG23	2.01	0.43
1:A:79:GLU:O	1:A:83:ARG:HG2	2.19	0.43
1:F:33:ALA:C	1:F:35:SER:H	2.22	0.42
1:A:4:GLY:HA3	1:A:47:ALA:O	2.18	0.42
1:A:32:LEU:HD12	1:A:32:LEU:HA	1.80	0.42
4:D:8903:GPP:H11	1:F:139:PHE:CE2	2.55	0.42
1:D:54:LEU:HD21	1:D:87:LYS:HD2	2.01	0.42
1:E:23:VAL:HG22	1:E:111:GLN:HG3	2.01	0.42
1:D:41:LEU:HD22	1:D:77:LEU:HD22	2.02	0.42
1:E:94:VAL:HG23	1:E:127:VAL:HA	2.02	0.42
1:E:99:ILE:HB	1:E:147:ALA:HB3	2.02	0.42
1:C:109:ILE:HG23	1:C:129:VAL:HB	2.02	0.42
1:E:109:ILE:N	1:E:110:PRO:HD2	2.35	0.41
1:F:6:GLY:HA3	1:F:43:ALA:O	2.19	0.41
1:F:97:THR:HB	1:F:149:GLU:HB2	2.02	0.41
1:A:10:HIS:CD2	1:A:39:VAL:HG13	2.54	0.41
1:D:106:LEU:HB3	1:D:107:PRO:HD3	2.01	0.41
1:B:35:SER:HB3	1:B:42:HIS:CE1	2.56	0.41
1:A:2:ARG:HD2	1:A:89:TYR:CE2	2.55	0.41
1:A:91:LEU:HD11	1:A:94:VAL:HG13	2.01	0.41
1:A:33:ALA:O	1:A:34:HIS:CD2	2.74	0.41
1:A:97:THR:HB	1:A:149:GLU:HB2	2.01	0.41
1:E:9:VAL:HB	1:F:134:THR:HA	2.02	0.41
1:E:109:ILE:HG23	1:E:129:VAL:HB	2.02	0.40
1:C:111:GLN:HE21	1:C:111:GLN:HA	1.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:ILE:CD1	1:B:3:ILE:CD1	2.99	0.40
1:B:6:GLY:HA3	1:B:43:ALA:O	2.21	0.40
1:E:3:ILE:CD1	1:F:3:ILE:HD11	2.40	0.40
1:D:109:ILE:N	1:D:110:PRO:CD	2.84	0.40
1:C:109:ILE:HB	1:C:110:PRO:HD3	2.04	0.40
1:C:31:LEU:HD12	5:C:3913:HOH:O	2.21	0.40
1:C:100:ALA:HB2	1:C:146:ILE:HG23	2.03	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	154/159 (97%)	148 (96%)	6 (4%)	0	100	100
1	B	154/159 (97%)	147 (96%)	7 (4%)	0	100	100
1	C	154/159 (97%)	147 (96%)	6 (4%)	1 (1%)	30	36
1	D	154/159 (97%)	148 (96%)	4 (3%)	2 (1%)	15	15
1	E	154/159 (97%)	148 (96%)	6 (4%)	0	100	100
1	F	154/159 (97%)	151 (98%)	3 (2%)	0	100	100
All	All	924/954 (97%)	889 (96%)	32 (4%)	3 (0%)	46	57

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	102	ALA
1	D	103	PRO
1	C	33	ALA

### 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	117/119 (98%)	115 (98%)	2 (2%)	68	83
1	B	117/119 (98%)	117 (100%)	0	100	100
1	C	117/119 (98%)	115 (98%)	2 (2%)	68	83
1	D	117/119 (98%)	112 (96%)	5 (4%)	35	47
1	E	117/119 (98%)	115 (98%)	2 (2%)	68	83
1	F	117/119 (98%)	115 (98%)	2 (2%)	68	83
All	All	702/714 (98%)	689 (98%)	13 (2%)	65	81

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

Mol	Chain	Res	Type
1	A	32	LEU
1	A	39	VAL
1	C	39	VAL
1	C	111	GLN
1	D	39	VAL
1	D	63	ASP
1	D	64	THR
1	D	104	LYS
1	D	106	LEU
1	E	3	ILE
1	E	61	PHE
1	F	39	VAL
1	F	64	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	B	34	HIS
1	C	111	GLN



### 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 ⓘ

Of 14 ligands modelled in this entry, 6 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$
3	1AA	A	1901	2	32,39,39	1.39	3 (9%)	44,57,57	1.22	4 (9%)
4	GPP	A	7903	-	16,18,18	0.63	0	21,25,25	1.32	2 (9%)
3	1AA	C	2901	2	32,39,39	1.41	3 (9%)	44,57,57	1.24	4 (9%)
3	1AA	C	3901	2	32,39,39	1.39	3 (9%)	44,57,57	1.33	4 (9%)
3	1AA	D	6901	2	32,39,39	1.38	3 (9%)	44,57,57	1.36	4 (9%)
4	GPP	D	8903	-	16,18,18	0.62	0	21,25,25	1.19	2 (9%)
3	1AA	E	4901	-	32,39,39	1.37	3 (9%)	44,57,57	1.30	4 (9%)
3	1AA	E	5901	2	32,39,39	1.41	3 (9%)	44,57,57	1.43	5 (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
3	1AA	A	1901	2	-	0/24/44/44	0/3/3/3
4	GPP	A	7903	-	-	0/19/19/19	0/0/0/0
3	1AA	C	2901	2	-	0/24/44/44	0/3/3/3
3	1AA	C	3901	2	-	0/24/44/44	0/3/3/3
3	1AA	D	6901	2	-	0/24/44/44	0/3/3/3
4	GPP	D	8903	-	-	0/19/19/19	0/0/0/0
3	1AA	E	4901	-	-	0/24/44/44	0/3/3/3
3	1AA	E	5901	2	-	0/24/44/44	0/3/3/3

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	3901	1AA	C9'-N4	2.67	1.40	1.35
3	D	6901	1AA	C9'-N4	2.68	1.40	1.35
3	A	1901	1AA	C9'-N4	2.70	1.40	1.35
3	E	4901	1AA	C9'-N4	2.71	1.40	1.35
3	E	5901	1AA	C9'-N4	2.71	1.40	1.35
3	C	2901	1AA	C9'-N4	2.73	1.40	1.35
3	E	4901	1AA	C7'-N2	4.07	1.41	1.35
3	D	6901	1AA	C7'-N2	4.13	1.41	1.35
3	D	6901	1AA	O2'-C2	4.13	1.46	1.41
3	C	2901	1AA	C7'-N2	4.15	1.41	1.35
3	C	3901	1AA	O2'-C2	4.17	1.46	1.41
3	E	4901	1AA	O2'-C2	4.17	1.46	1.41
3	A	1901	1AA	C7'-N2	4.21	1.41	1.35
3	C	3901	1AA	C7'-N2	4.23	1.41	1.35
3	A	1901	1AA	O2'-C2	4.24	1.46	1.41
3	E	5901	1AA	C7'-N2	4.31	1.41	1.35
3	E	5901	1AA	O2'-C2	4.33	1.46	1.41
3	C	2901	1AA	O2'-C2	4.39	1.46	1.41

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	6901	1AA	P2-O3-P1	-3.56	122.72	132.73
3	E	5901	1AA	P2-O3-P1	-3.53	122.82	132.73
3	E	4901	1AA	P2-O3-P1	-3.41	123.15	132.73
4	A	7903	GPP	PA-O3A-PB	-3.07	122.39	132.67
3	A	1901	1AA	P2-O3-P1	-2.89	124.62	132.73
4	D	8903	GPP	PA-O3A-PB	-2.75	123.45	132.67
3	C	3901	1AA	O5-C9-O6	-2.56	118.96	123.66
3	E	5901	1AA	O5-C9-O6	-2.51	119.05	123.66
3	C	3901	1AA	P2-O3-P1	-2.44	125.89	132.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	6901	1AA	O5-C9-O6	-2.36	119.33	123.66
3	C	2901	1AA	P2-O3-P1	-2.28	126.32	132.73
3	E	4901	1AA	C5'-C6'-C1'	2.06	120.15	118.06
3	C	2901	1AA	C5'-C6'-C1'	2.14	120.23	118.06
3	E	5901	1AA	O5-C8-C7	2.19	115.04	108.54
3	A	1901	1AA	C5'-C6'-C1'	2.21	120.30	118.06
4	D	8903	GPP	C4-C3-C5	2.25	118.85	115.41
4	A	7903	GPP	C4-C3-C5	2.90	119.84	115.41
3	E	4901	1AA	C1-N4-C9'	3.35	120.34	115.61
3	A	1901	1AA	C1-N4-C9'	3.37	120.36	115.61
3	C	3901	1AA	C1-N4-C9'	3.42	120.43	115.61
3	D	6901	1AA	C1-N4-C9'	3.51	120.56	115.61
3	C	2901	1AA	C1-N4-C9'	3.51	120.56	115.61
3	E	5901	1AA	C1-N4-C9'	3.55	120.62	115.61
3	A	1901	1AA	O5-C9-C1'	3.91	119.28	112.16
3	E	4901	1AA	O5-C9-C1'	3.96	119.36	112.16
3	C	2901	1AA	O5-C9-C1'	4.23	119.85	112.16
3	C	3901	1AA	O5-C9-C1'	4.57	120.48	112.16
3	D	6901	1AA	O5-C9-C1'	4.66	120.64	112.16
3	E	5901	1AA	O5-C9-C1'	5.05	121.34	112.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	7903	GPP	1	0
3	C	2901	1AA	2	0
3	C	3901	1AA	2	0
4	D	8903	GPP	1	0
3	E	5901	1AA	1	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, 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	156/159 (98%)	0.58	9 (5%) 26 35	27, 36, 49, 52	0
1	B	156/159 (98%)	0.36	6 (3%) 44 53	19, 29, 42, 45	0
1	C	156/159 (98%)	0.37	8 (5%) 32 41	19, 28, 44, 48	0
1	D	156/159 (98%)	0.27	5 (3%) 51 60	20, 30, 41, 44	1 (0%)
1	E	156/159 (98%)	0.41	12 (7%) 16 23	19, 28, 45, 49	1 (0%)
1	F	156/159 (98%)	0.49	5 (3%) 51 60	22, 33, 46, 50	0
All	All	936/954 (98%)	0.42	45 (4%) 34 43	19, 30, 45, 52	2 (0%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	64	THR	8.6
1	C	64	THR	5.9
1	B	32	LEU	5.8
1	E	32	LEU	5.7
1	D	34	HIS	5.2
1	E	34	HIS	5.0
1	A	70	GLY	4.8
1	A	34	HIS	4.6
1	D	63	ASP	4.1
1	A	63	ASP	4.0
1	E	64	THR	4.0
1	B	34	HIS	4.0
1	C	68	PHE	4.0
1	E	68	PHE	4.0
1	D	32	LEU	3.9
1	C	33	ALA	3.7
1	C	34	HIS	3.7
1	C	32	LEU	3.7
1	F	34	HIS	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	64	THR	3.5
1	E	67	ALA	3.0
1	A	86	ALA	3.0
1	F	68	PHE	2.8
1	F	83	ARG	2.7
1	E	33	ALA	2.7
1	B	16	GLY	2.6
1	E	66	PRO	2.6
1	C	57	ILE	2.6
1	E	42	HIS	2.6
1	A	24	ARG	2.6
1	A	71	ALA	2.5
1	B	57	ILE	2.5
1	D	156	LYS	2.4
1	E	13	GLY	2.4
1	E	63	ASP	2.4
1	C	124	MET	2.4
1	E	57	ILE	2.3
1	A	68	PHE	2.3
1	B	66	PRO	2.2
1	D	102	ALA	2.1
1	A	31	LEU	2.1
1	F	65	ASP	2.1
1	B	28	GLU	2.1
1	E	16	GLY	2.0
1	C	1	MET	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( $\text{\AA}^2$ )	Q<0.9
4	GPP	D	8903	19/19	0.76	0.47	17.85	6,9,12,12	19
4	GPP	A	7903	19/19	0.76	0.47	13.27	2,7,10,10	19
3	1AA	E	4901	37/37	0.66	0.30	2.08	49,57,67,67	0
3	1AA	D	6901	37/37	0.70	0.29	1.77	62,67,75,75	0
3	1AA	E	5901	37/37	0.82	0.27	1.25	42,51,61,61	0
3	1AA	C	2901	37/37	0.81	0.25	1.18	35,46,57,57	0
3	1AA	A	1901	37/37	0.70	0.24	0.70	60,66,72,72	11
3	1AA	C	3901	37/37	0.81	0.23	0.59	39,45,56,56	0
2	ZN	E	5900	1/1	0.99	0.07	-1.67	34,34,34,34	0
2	ZN	B	2900	1/1	0.99	0.07	-1.67	36,36,36,36	0
2	ZN	D	4900	1/1	0.99	0.09	-1.82	37,37,37,37	0
2	ZN	C	3900	1/1	0.99	0.09	-3.01	30,30,30,30	0
2	ZN	F	6900	1/1	0.99	0.05	-	46,46,46,46	0
2	ZN	A	1900	1/1	0.99	0.06	-	45,45,45,45	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.