Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 08:08 PM GMT

PDB ID	:	1IV3
Title	:	Structure of 2C-Methyl-D-erythritol-2,4-cyclodiphosphate Synthase (bound form MG atoms)
Authors	:	Kishida, H.; Wada, T.; Unzai, S.; Kuzuyama, T.; Terada, T.; Sirouzu, M.; Yokoyama, S.; Tame, J.R.H.; Park, S.-Y.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on	:	2002-03-11
Resolution	:	1.52 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 (i) symbol.

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

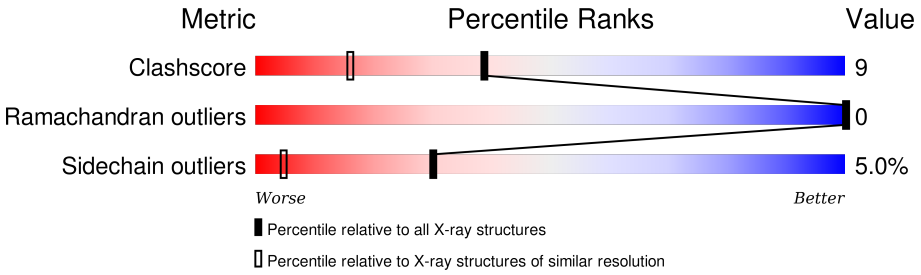
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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 1.52 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	102246	2887 (1.54-1.50)
Ramachandran outliers	100387	2818 (1.54-1.50)
Sidechain outliers	100360	2816 (1.54-1.50)

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 <=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	152	80% 18%
1	B	152	80% 16%
1	C	152	74% 20%
1	D	152	76% 20%
1	E	152	74% 18% 5%
1	F	152	80% 16%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7475 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	150	Total	C	N	O	S	0	0	0
			1152	726	213	209	4			
1	B	150	Total	C	N	O	S	0	0	0
			1152	726	213	209	4			
1	C	150	Total	C	N	O	S	0	0	0
			1152	726	213	209	4			
1	D	150	Total	C	N	O	S	0	0	0
			1152	726	213	209	4			
1	E	150	Total	C	N	O	S	0	0	0
			1152	726	213	209	4			
1	F	150	Total	C	N	O	S	0	0	0
			1152	726	213	209	4			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	41	MET	LEU	ENGINEERED	UNP Q8RQP5
A	81	MET	LEU	ENGINEERED	UNP Q8RQP5
A	120	MET	LEU	ENGINEERED	UNP Q8RQP5
B	241	MET	LEU	ENGINEERED	UNP Q8RQP5
B	281	MET	LEU	ENGINEERED	UNP Q8RQP5
B	320	MET	LEU	ENGINEERED	UNP Q8RQP5
C	441	MET	LEU	ENGINEERED	UNP Q8RQP5
C	481	MET	LEU	ENGINEERED	UNP Q8RQP5
C	520	MET	LEU	ENGINEERED	UNP Q8RQP5
D	1041	MET	LEU	ENGINEERED	UNP Q8RQP5
D	1081	MET	LEU	ENGINEERED	UNP Q8RQP5
D	1120	MET	LEU	ENGINEERED	UNP Q8RQP5
E	1241	MET	LEU	ENGINEERED	UNP Q8RQP5
E	1281	MET	LEU	ENGINEERED	UNP Q8RQP5
E	1320	MET	LEU	ENGINEERED	UNP Q8RQP5
F	1441	MET	LEU	ENGINEERED	UNP Q8RQP5
F	1481	MET	LEU	ENGINEERED	UNP Q8RQP5

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Chain	Residue	Modelled	Actual	Comment	Reference
F	1520	MET	LEU	ENGINEERED	UNP Q8RQP5

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	112	Total O 112 112	0	0
3	B	94	Total O 94 94	0	0
3	C	90	Total O 90 90	0	0
3	D	87	Total O 87 87	0	0
3	E	90	Total O 90 90	0	0
3	F	84	Total O 84 84	0	0

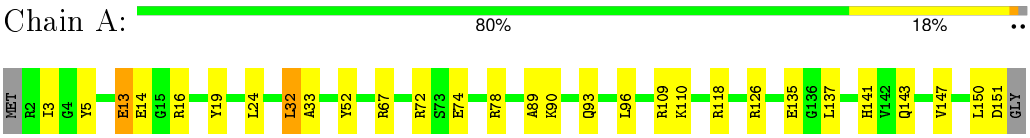
3 Residue-property plots (i)

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.

Note EDS was not executed.

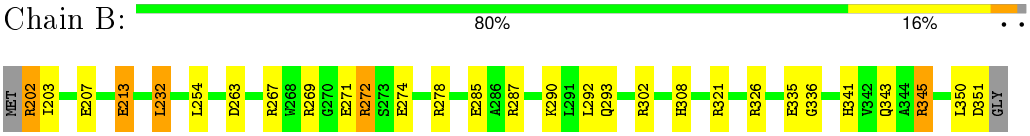
- Molecule 1: 2-C-methyl-D-erythritol 2,4-cyclodiphosphate synthase

Chain A:



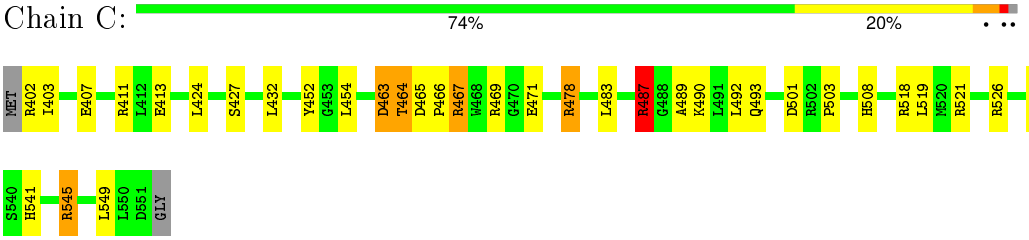
- Molecule 1: 2-C-methyl-D-erythritol 2,4-cyclodiphosphate synthase

Chain B:



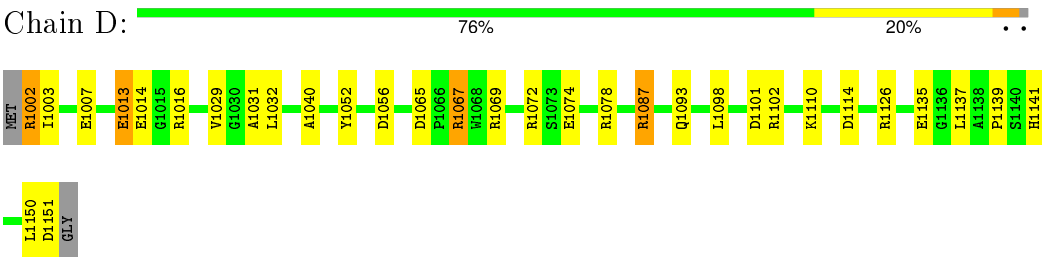
- Molecule 1: 2-C-methyl-D-erythritol 2,4-cyclodiphosphate synthase

Chain C:



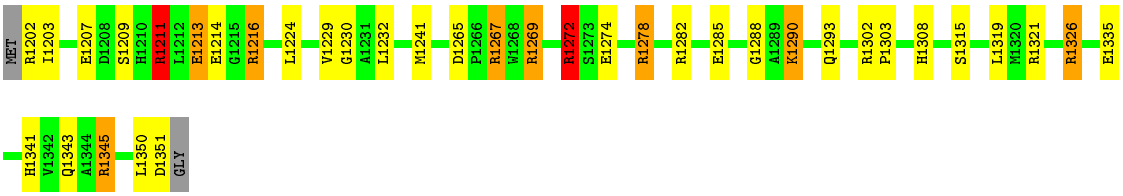
- Molecule 1: 2-C-methyl-D-erythritol 2,4-cyclodiphosphate synthase

Chain D:



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

Chain E:



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

Chain F:

4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	106.18Å 106.18Å 148.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.52	Depositor
% Data completeness (in resolution range)	93.5 (20.00-1.52)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.202 , 0.257	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7475	wwPDB-VP
Average B, all atoms (Å^2)	20.0	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: MG

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	0.39	0/1173	1.15	7/1587 (0.4%)
1	B	0.39	0/1173	1.12	6/1587 (0.4%)
1	C	0.41	0/1173	1.25	8/1587 (0.5%)
1	D	0.39	0/1173	1.17	10/1587 (0.6%)
1	E	0.40	0/1173	1.24	13/1587 (0.8%)
1	F	0.38	0/1173	1.14	4/1587 (0.3%)
All	All	0.39	0/7038	1.18	48/9522 (0.5%)

There are no bond length outliers.

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	1072	ARG	NE-CZ-NH2	-13.71	113.45	120.30
1	C	487	ARG	CD-NE-CZ	13.30	142.23	123.60
1	E	1272	ARG	NE-CZ-NH2	9.34	124.97	120.30
1	C	518	ARG	NE-CZ-NH2	-9.04	115.78	120.30
1	C	487	ARG	NE-CZ-NH1	8.99	124.80	120.30
1	F	1402	ARG	NE-CZ-NH1	8.65	124.62	120.30
1	B	326	ARG	NE-CZ-NH2	-8.60	116.00	120.30
1	A	126	ARG	NE-CZ-NH2	-8.43	116.08	120.30
1	C	545	ARG	NE-CZ-NH1	8.30	124.45	120.30
1	D	1126	ARG	NE-CZ-NH2	-7.86	116.37	120.30
1	B	326	ARG	NE-CZ-NH1	7.79	124.20	120.30
1	A	126	ARG	NE-CZ-NH1	7.72	124.16	120.30
1	A	67	ARG	NE-CZ-NH1	-7.70	116.45	120.30
1	C	545	ARG	NE-CZ-NH2	-7.66	116.47	120.30
1	E	1326	ARG	NE-CZ-NH2	-7.59	116.50	120.30
1	A	19	TYR	CB-CG-CD1	-7.50	116.50	121.00
1	C	518	ARG	NE-CZ-NH1	7.11	123.85	120.30
1	F	1526	ARG	NE-CZ-NH2	-6.88	116.86	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	478	ARG	NE-CZ-NH1	6.83	123.72	120.30
1	A	72	ARG	NE-CZ-NH2	-6.61	116.99	120.30
1	A	109	ARG	NE-CZ-NH1	-6.58	117.01	120.30
1	E	1345	ARG	NE-CZ-NH2	-6.53	117.03	120.30
1	E	1211	ARG	NE-CZ-NH1	6.49	123.55	120.30
1	C	521	ARG	NE-CZ-NH1	6.49	123.54	120.30
1	E	1345	ARG	NE-CZ-NH1	6.42	123.51	120.30
1	F	1518	ARG	NE-CZ-NH2	-6.41	117.09	120.30
1	D	1072	ARG	NH1-CZ-NH2	6.40	126.44	119.40
1	E	1351	ASP	CB-CG-OD1	-6.28	112.65	118.30
1	F	1551	ASP	CB-CG-OD2	-6.28	112.65	118.30
1	E	1211	ARG	CD-NE-CZ	6.26	132.37	123.60
1	D	1002	ARG	NE-CZ-NH1	6.11	123.35	120.30
1	E	1278	ARG	NE-CZ-NH1	5.98	123.29	120.30
1	B	202	ARG	NE-CZ-NH1	5.96	123.28	120.30
1	D	1151	ASP	CB-CG-OD1	-5.93	112.96	118.30
1	D	1002	ARG	NE-CZ-NH2	-5.85	117.38	120.30
1	E	1350	LEU	C-N-CA	5.84	136.29	121.70
1	A	118	ARG	NE-CZ-NH2	-5.67	117.47	120.30
1	E	1326	ARG	NE-CZ-NH1	5.64	123.12	120.30
1	D	1087	ARG	NE-CZ-NH2	-5.63	117.49	120.30
1	B	302	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	E	1278	ARG	CD-NE-CZ	5.38	131.13	123.60
1	E	1269	ARG	NE-CZ-NH1	-5.28	117.66	120.30
1	B	345	ARG	NE-CZ-NH2	-5.25	117.68	120.30
1	D	1150	LEU	O-C-N	5.18	130.99	122.70
1	D	1072	ARG	CD-NE-CZ	-5.08	116.48	123.60
1	D	1087	ARG	NE-CZ-NH1	5.08	122.84	120.30
1	B	272	ARG	NE-CZ-NH2	-5.08	117.76	120.30
1	E	1282	ARG	NE-CZ-NH1	-5.01	117.80	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1152	0	1182	24	0
1	B	1152	0	1182	18	0
1	C	1152	0	1182	28	0
1	D	1152	0	1182	30	0
1	E	1152	0	1182	31	0
1	F	1152	0	1180	21	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	112	0	0	1	0
3	B	94	0	0	1	0
3	C	90	0	0	1	0
3	D	87	0	0	1	0
3	E	90	0	0	4	0
3	F	84	0	0	1	0
All	All	7475	0	7090	126	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 9.

All (126) 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:E:1216:ARG:NH2	1:E:1232:LEU:HD12	1.63	1.13
1:E:1216:ARG:NH2	1:E:1232:LEU:CD1	2.28	0.97
1:E:1216:ARG:HH21	1:E:1232:LEU:HD12	1.28	0.90
1:C:483:LEU:HD22	1:C:487:ARG:HH22	1.39	0.86
1:D:1065:ASP:OD1	1:D:1067:ARG:HG3	1.84	0.78
1:F:1416:ARG:HB3	1:F:1416:ARG:HH11	1.50	0.77
1:E:1216:ARG:HH22	1:E:1232:LEU:CD1	2.02	0.72
1:C:463:ASP:O	1:C:469:ARG:HD3	1.90	0.71
1:E:1241:MET:HE1	1:E:1274:GLU:HA	1.73	0.69
1:A:13:GLU:OE1	1:A:32:LEU:HD13	1.92	0.69
1:C:454:LEU:HD21	1:C:487:ARG:HE	1.59	0.68
1:B:207:GLU:HG2	1:B:345:ARG:HG2	1.76	0.68
1:D:1101:ASP:OD2	1:D:1139:PRO:HA	1.96	0.66
1:E:1213:GLU:OE2	1:E:1232:LEU:HD13	1.96	0.64
1:C:501:ASP:OD2	1:C:539:PRO:HA	1.99	0.63
1:E:1207:GLU:HG2	1:E:1345:ARG:HG2	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:483:LEU:HB3	1:C:487:ARG:NH2	2.17	0.60
1:E:1343:GLN:NE2	1:E:1345:ARG:HH22	1.99	0.60
1:F:1474:GLU:O	1:F:1478:ARG:HG2	2.02	0.59
1:A:3:ILE:H	1:C:493:GLN:NE2	2.01	0.59
1:B:213:GLU:OE1	1:B:232:LEU:HD22	2.04	0.57
1:B:278:ARG:HG2	1:B:278:ARG:HH11	1.70	0.56
1:F:1501:ASP:OD2	1:F:1539:PRO:HA	2.06	0.56
1:E:1315:SER:O	1:E:1319:LEU:HD23	2.06	0.56
1:C:508:HIS:HD2	3:C:1573:HOH:O	1.90	0.55
1:D:1143:GLN:NE2	1:D:1145:ARG:HH22	2.05	0.54
1:C:465:ASP:OD1	1:C:467:ARG:HG3	2.07	0.54
1:E:1341:HIS:HD2	3:E:230:HOH:O	1.90	0.54
1:D:1003:ILE:HD13	1:F:1547:VAL:HG11	1.88	0.54
1:A:141:HIS:HE1	1:C:535:GLU:O	1.91	0.54
1:C:413:GLU:CD	1:C:432:LEU:HD13	2.28	0.53
1:A:150:LEU:O	1:A:151:ASP:HB2	2.09	0.53
1:B:350:LEU:O	1:B:351:ASP:HB2	2.08	0.53
1:A:135:GLU:O	1:B:341:HIS:HE1	1.92	0.53
1:A:93:GLN:HE22	1:B:202:ARG:HA	1.74	0.53
1:C:465:ASP:OD1	1:C:466:PRO:HD2	2.10	0.52
1:B:335:GLU:O	1:C:541:HIS:HE1	1.92	0.52
1:D:1135:GLU:HG3	1:E:1209:SER:O	2.10	0.52
1:E:1265:ASP:OD1	1:E:1267:ARG:HB2	2.08	0.52
1:B:336:GLY:HA3	1:C:411:ARG:HH21	1.74	0.52
1:D:1141:HIS:HE1	1:F:1535:GLU:O	1.92	0.51
1:C:492:LEU:HD23	1:C:549:LEU:HD23	1.93	0.51
1:A:110:LYS:HD2	1:D:1069:ARG:HD2	1.93	0.51
1:A:52:TYR:OH	1:A:89:ALA:HB2	2.11	0.51
1:E:1335:GLU:HA	1:E:1335:GLU:OE1	2.12	0.51
1:D:1135:GLU:O	1:E:1341:HIS:HE1	1.93	0.50
1:A:110:LYS:HD2	1:D:1069:ARG:CD	2.41	0.50
1:F:1412:LEU:O	1:F:1413:GLU:OE1	2.29	0.50
1:F:1416:ARG:NH1	1:F:1416:ARG:HB3	2.22	0.50
1:A:93:GLN:NE2	1:B:203:ILE:H	2.10	0.49
1:C:427:SER:OG	1:C:503:PRO:HD3	2.12	0.49
1:D:1078:ARG:HH11	1:D:1078:ARG:HG2	1.77	0.49
1:E:1216:ARG:NH2	1:E:1232:LEU:HD13	2.20	0.48
1:A:16:ARG:NH2	1:A:32:LEU:HD12	2.28	0.48
1:A:5:TYR:HD1	1:A:147:VAL:HG22	1.78	0.48
1:C:487:ARG:HG3	1:C:487:ARG:HH11	1.77	0.48
1:D:1016:ARG:NH2	1:D:1032:LEU:HD13	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:293:GLN:NE2	1:C:403:ILE:H	2.11	0.48
1:F:1407:GLU:HG2	1:F:1545:ARG:HG2	1.96	0.47
1:B:213:GLU:OE1	1:B:232:LEU:HD13	2.13	0.47
1:F:1416:ARG:NH1	1:F:1431:ALA:O	2.48	0.47
1:B:254:LEU:HD21	1:B:287:ARG:HE	1.79	0.47
1:F:1411:ARG:HG2	1:F:1413:GLU:OE1	2.15	0.47
1:E:1211:ARG:HH11	1:E:1211:ARG:HG3	1.79	0.47
1:E:1278:ARG:HH21	1:E:1319:LEU:HA	1.79	0.46
1:B:267:ARG:HG2	1:D:1087:ARG:NH2	2.30	0.46
1:A:16:ARG:HE	1:A:32:LEU:HB2	1.80	0.46
1:A:110:LYS:HG3	1:D:1069:ARG:CZ	2.45	0.46
1:A:52:TYR:CZ	1:A:89:ALA:HB2	2.50	0.46
1:C:492:LEU:HD21	1:C:526:ARG:NH1	2.31	0.46
1:A:16:ARG:NE	1:A:32:LEU:HB2	2.31	0.46
1:D:1014:GLU:OE2	1:D:1102:ARG:NH2	2.49	0.46
1:E:1272:ARG:NH1	3:E:438:HOH:O	2.49	0.46
1:D:1093:GLN:NE2	1:E:1203:ILE:H	2.14	0.46
1:A:32:LEU:O	1:A:33:ALA:HB2	2.16	0.46
1:E:1335:GLU:O	1:F:1541:HIS:HE1	1.98	0.46
1:D:1110:LYS:NZ	1:D:1114:ASP:OD2	2.49	0.46
1:D:1007:GLU:OE2	1:F:1545:ARG:HD2	2.17	0.45
1:E:1214:GLU:HA	1:E:1229:VAL:HG12	1.98	0.45
1:E:1211:ARG:HG3	1:E:1211:ARG:NH1	2.31	0.45
1:C:464:THR:O	1:C:469:ARG:NH1	2.50	0.45
1:E:1285:GLU:OE2	1:E:1321:ARG:NH2	2.50	0.45
1:B:269:ARG:O	1:B:269:ARG:HG2	2.17	0.45
1:E:1288:GLY:O	1:E:1290:LYS:HE3	2.17	0.45
1:A:110:LYS:HG3	1:D:1069:ARG:NH1	2.32	0.44
1:C:535:GLU:OE2	1:C:535:GLU:HA	2.17	0.44
1:B:308:HIS:HE1	3:B:1650:HOH:O	1.99	0.44
1:C:467:ARG:O	1:C:471:GLU:HG3	2.17	0.44
1:A:74:GLU:O	1:A:78:ARG:HG3	2.18	0.44
1:C:532:LYS:HB3	1:C:532:LYS:HE3	1.69	0.44
1:D:1074:GLU:OE2	1:D:1078:ARG:NH2	2.50	0.43
1:C:402:ARG:HH11	1:C:402:ARG:HD3	1.68	0.43
1:F:1551:ASP:OD2	1:F:1551:ASP:N	2.50	0.43
1:F:1431:ALA:HB3	1:F:1437:GLY:HA3	2.01	0.43
1:D:1014:GLU:OE2	1:D:1029:VAL:HG12	2.17	0.43
1:D:1074:GLU:OE2	1:D:1078:ARG:NE	2.50	0.43
1:A:143:GLN:HG2	3:A:1669:HOH:O	2.18	0.43
1:A:3:ILE:H	1:C:493:GLN:HE22	1.65	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274:GLU:OE2	1:B:278:ARG:NH2	2.51	0.43
1:F:1550:LEU:O	1:F:1551:ASP:HB2	2.19	0.43
1:D:1031:ALA:HB3	3:D:539:HOH:O	2.19	0.43
1:B:343:GLN:NE2	1:B:345:ARG:HH22	2.17	0.42
1:A:52:TYR:CE1	1:A:89:ALA:HB2	2.54	0.42
1:D:1013:GLU:OE1	1:D:1032:LEU:HB2	2.19	0.42
1:A:3:ILE:HD13	1:C:403:ILE:HD11	2.02	0.42
1:B:285:GLU:OE2	1:B:321:ARG:NH2	2.52	0.42
1:C:407:GLU:HG2	1:C:545:ARG:HG2	2.01	0.42
1:C:492:LEU:HD21	1:C:526:ARG:HH12	1.83	0.42
1:E:1216:ARG:O	1:E:1230:GLY:HA3	2.20	0.42
1:D:1003:ILE:H	1:F:1493:GLN:NE2	2.18	0.42
1:D:1040:ALA:HB2	1:D:1098:LEU:HD21	2.02	0.41
1:D:1002:ARG:HD2	1:D:1052:TYR:CZ	2.55	0.41
1:E:1302:ARG:HB2	1:E:1303:PRO:HA	2.01	0.41
1:D:1137:LEU:HD21	3:E:541:HOH:O	2.20	0.41
1:F:1410:HIS:HB3	1:F:1431:ALA:HB1	2.03	0.41
1:F:1541:HIS:HD2	3:F:334:HOH:O	2.03	0.41
1:D:1093:GLN:HE22	1:E:1202:ARG:HA	1.86	0.41
1:E:1269:ARG:HD3	1:E:1269:ARG:HH11	1.68	0.41
1:D:1056:ASP:HB3	1:F:1530:THR:HG22	2.02	0.41
1:A:96:LEU:HD23	1:A:96:LEU:N	2.36	0.41
1:F:1408:ASP:OD2	1:F:1410:HIS:CD2	2.74	0.40
1:E:1345:ARG:HD2	1:F:1407:GLU:OE2	2.21	0.40
1:D:1016:ARG:HH21	1:D:1032:LEU:HD13	1.86	0.40
1:E:1293:GLN:NE2	1:E:1326:ARG:HH11	2.19	0.40
1:C:452:TYR:CE1	1:C:489:ALA:HB2	2.56	0.40
1:E:1308:HIS:HE1	3:E:307:HOH:O	2.04	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	148/152 (97%)	145 (98%)	3 (2%)	0	100	100
1	B	148/152 (97%)	147 (99%)	1 (1%)	0	100	100
1	C	148/152 (97%)	143 (97%)	5 (3%)	0	100	100
1	D	148/152 (97%)	145 (98%)	3 (2%)	0	100	100
1	E	148/152 (97%)	146 (99%)	2 (1%)	0	100	100
1	F	148/152 (97%)	146 (99%)	2 (1%)	0	100	100
All	All	888/912 (97%)	872 (98%)	16 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	121/122 (99%)	115 (95%)	6 (5%)	30	5
1	B	121/122 (99%)	114 (94%)	7 (6%)	25	3
1	C	121/122 (99%)	113 (93%)	8 (7%)	21	2
1	D	121/122 (99%)	119 (98%)	2 (2%)	68	36
1	E	121/122 (99%)	114 (94%)	7 (6%)	25	3
1	F	121/122 (99%)	115 (95%)	6 (5%)	30	5
All	All	726/732 (99%)	690 (95%)	36 (5%)	30	5

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	GLU
1	A	14	GLU
1	A	24	LEU
1	A	32	LEU
1	A	90	LYS
1	A	137	LEU
1	B	213	GLU

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Mol	Chain	Res	Type
1	B	232	LEU
1	B	263	ASP
1	B	271	GLU
1	B	272	ARG
1	B	290	LYS
1	B	292	LEU
1	C	424	LEU
1	C	463	ASP
1	C	464	THR
1	C	467	ARG
1	C	478	ARG
1	C	487	ARG
1	C	490	LYS
1	C	519	LEU
1	D	1013	GLU
1	D	1067	ARG
1	E	1211	ARG
1	E	1213	GLU
1	E	1216	ARG
1	E	1224	LEU
1	E	1267	ARG
1	E	1272	ARG
1	E	1290	LYS
1	F	1413	GLU
1	F	1416	ARG
1	F	1432	LEU
1	F	1434	HIS
1	F	1463	ASP
1	F	1490	LYS

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	93	GLN
1	A	141	HIS
1	B	293	GLN
1	B	341	HIS
1	B	343	GLN
1	C	493	GLN
1	C	508	HIS
1	C	541	HIS
1	C	543	GLN

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Mol	Chain	Res	Type
1	D	1093	GLN
1	D	1141	HIS
1	D	1143	GLN
1	E	1293	GLN
1	E	1308	HIS
1	E	1341	HIS
1	E	1343	GLN
1	F	1493	GLN
1	F	1541	HIS
1	F	1543	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 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.