



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:58 PM UTC

PDB ID : 7VG4 / pdb_00007vg4
Title : 10,5-methenyltetrahydrofolate cyclohydrolase from *Methylobacterium extorquens* AM1 strain
Authors : Kim, S.; Lee, S.; Kim, I.-K.; Seo, H.; Kim, K.-J.
Deposited on : 2021-09-14
Resolution : 2.77 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
Xtriage (Phenix) : 2.0
EDS : 3.0
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

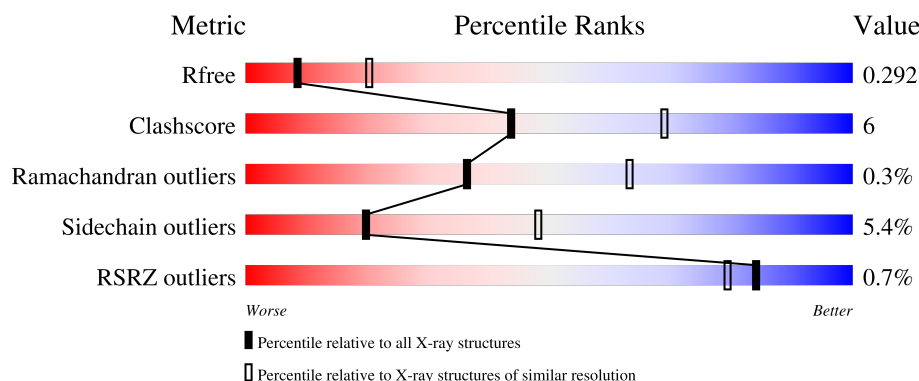
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.77 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	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

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	216	78%16%6%
1	B	216	83%12%5%
1	C	216	78%15%5%
1	D	216	75%19%6%
1	E	216	74%14%9%

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Mol	Chain	Length	Quality of chain
1	F	216	% 69% 17% 14%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8805 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 Methenyltetrahydrofolate cyclohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	204	Total	C	N	O	S	0	0	0
			1485	924	254	299	8			
1	B	206	Total	C	N	O	S	0	0	0
			1504	935	258	303	8			
1	C	205	Total	C	N	O	S	0	0	0
			1500	933	257	302	8			
1	D	204	Total	C	N	O	S	0	0	0
			1492	929	255	300	8			
1	E	196	Total	C	N	O	S	0	0	0
			1420	883	244	285	8			
1	F	186	Total	C	N	O	S	0	0	0
			1343	832	232	271	8			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	209	LEU	-	expression tag	UNP Q49135
A	210	GLU	-	expression tag	UNP Q49135
A	211	HIS	-	expression tag	UNP Q49135
A	212	HIS	-	expression tag	UNP Q49135
A	213	HIS	-	expression tag	UNP Q49135
A	214	HIS	-	expression tag	UNP Q49135
A	215	HIS	-	expression tag	UNP Q49135
A	216	HIS	-	expression tag	UNP Q49135
B	209	LEU	-	expression tag	UNP Q49135
B	210	GLU	-	expression tag	UNP Q49135
B	211	HIS	-	expression tag	UNP Q49135
B	212	HIS	-	expression tag	UNP Q49135
B	213	HIS	-	expression tag	UNP Q49135
B	214	HIS	-	expression tag	UNP Q49135
B	215	HIS	-	expression tag	UNP Q49135
B	216	HIS	-	expression tag	UNP Q49135
C	209	LEU	-	expression tag	UNP Q49135

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Chain	Residue	Modelled	Actual	Comment	Reference
C	210	GLU	-	expression tag	UNP Q49135
C	211	HIS	-	expression tag	UNP Q49135
C	212	HIS	-	expression tag	UNP Q49135
C	213	HIS	-	expression tag	UNP Q49135
C	214	HIS	-	expression tag	UNP Q49135
C	215	HIS	-	expression tag	UNP Q49135
C	216	HIS	-	expression tag	UNP Q49135
D	209	LEU	-	expression tag	UNP Q49135
D	210	GLU	-	expression tag	UNP Q49135
D	211	HIS	-	expression tag	UNP Q49135
D	212	HIS	-	expression tag	UNP Q49135
D	213	HIS	-	expression tag	UNP Q49135
D	214	HIS	-	expression tag	UNP Q49135
D	215	HIS	-	expression tag	UNP Q49135
D	216	HIS	-	expression tag	UNP Q49135
E	209	LEU	-	expression tag	UNP Q49135
E	210	GLU	-	expression tag	UNP Q49135
E	211	HIS	-	expression tag	UNP Q49135
E	212	HIS	-	expression tag	UNP Q49135
E	213	HIS	-	expression tag	UNP Q49135
E	214	HIS	-	expression tag	UNP Q49135
E	215	HIS	-	expression tag	UNP Q49135
E	216	HIS	-	expression tag	UNP Q49135
F	209	LEU	-	expression tag	UNP Q49135
F	210	GLU	-	expression tag	UNP Q49135
F	211	HIS	-	expression tag	UNP Q49135
F	212	HIS	-	expression tag	UNP Q49135
F	213	HIS	-	expression tag	UNP Q49135
F	214	HIS	-	expression tag	UNP Q49135
F	215	HIS	-	expression tag	UNP Q49135
F	216	HIS	-	expression tag	UNP Q49135

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	18	Total O 18 18	0	0
2	B	13	Total O 13 13	0	0
2	C	15	Total O 15 15	0	0
2	D	7	Total O 7 7	0	0

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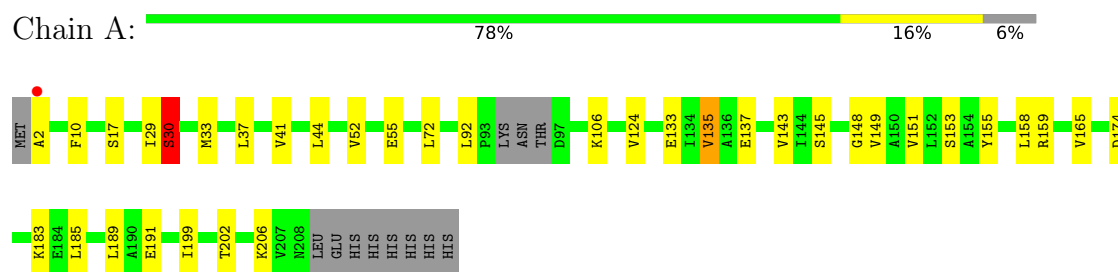
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	E	7	Total	O	0	0
			7	7		
2	F	1	Total	O	0	0
			1	1		

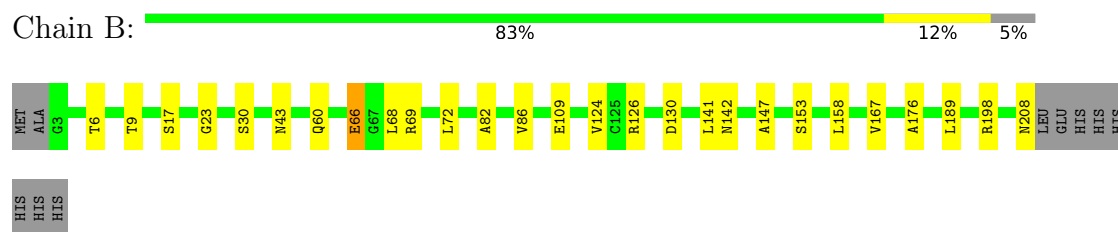
3 Residue-property plots

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.

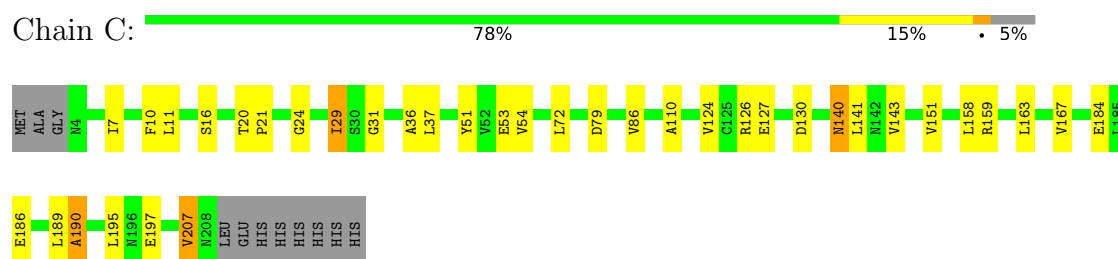
• Molecule 1: Methenyltetrahydrofolate cyclohydrolase



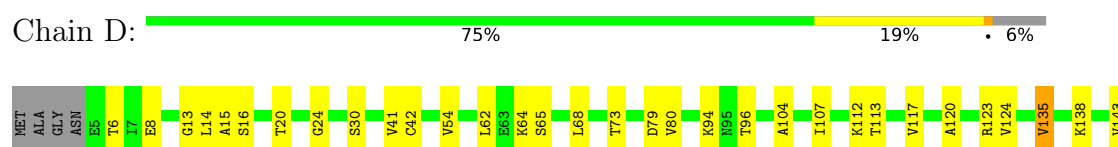
• Molecule 1: Methenyltetrahydrofolate cyclohydrolase



• Molecule 1: Methenyltetrahydrofolate cyclohydrolase

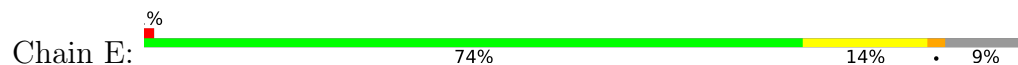


• Molecule 1: Methenyltetrahydrofolate cyclohydrolase

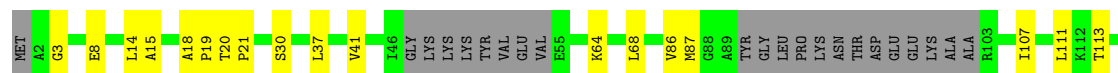




- Molecule 1: Methenyltetrahydrofolate cyclohydrolase



- Molecule 1: Methenyltetrahydrofolate cyclohydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	53.74Å 67.73Å 88.59Å 80.14° 79.31° 79.42°	Depositor
Resolution (Å)	30.59 – 2.77 30.59 – 2.77	Depositor EDS
% Data completeness (in resolution range)	94.5 (30.59-2.77) 94.5 (30.59-2.77)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.196 , 0.286 0.206 , 0.292	Depositor DCC
R_{free} test set	1405 reflections (4.64%)	wwPDB-VP
Wilson B-factor (Å ²)	47.7	Xtriage
Anisotropy	0.184	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.1	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	8805	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.75% 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](#)

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	1.05	0/1495	1.64	4/2017 (0.2%)
1	B	1.09	0/1515	1.70	4/2045 (0.2%)
1	C	1.07	0/1511	1.68	0/2040
1	D	1.09	0/1503	1.75	0/2029
1	E	1.08	0/1427	1.73	4/1922 (0.2%)
1	F	1.08	0/1349	1.74	6/1820 (0.3%)
All	All	1.08	0/8800	1.71	18/11873 (0.2%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	153	SER	CA-C-N	5.84	128.38	120.38
1	F	153	SER	C-N-CA	5.84	128.38	120.38
1	F	192	ALA	CA-C-N	5.68	126.25	119.94
1	F	192	ALA	C-N-CA	5.68	126.25	119.94
1	F	87	MET	CA-C-N	5.63	127.22	120.13
1	F	87	MET	C-N-CA	5.63	127.22	120.13
1	E	143	VAL	N-CA-C	-5.29	108.35	113.53
1	E	71	THR	CA-CB-OG1	-5.21	101.78	109.60
1	E	186	GLU	CA-C-N	5.21	125.72	119.94
1	E	186	GLU	C-N-CA	5.21	125.72	119.94
1	A	30	SER	CA-C-N	5.10	125.60	119.94
1	A	30	SER	C-N-CA	5.10	125.60	119.94
1	B	176	ALA	CA-C-N	5.09	127.36	120.38
1	B	176	ALA	C-N-CA	5.09	127.36	120.38
1	A	191	GLU	CA-C-N	5.05	127.04	120.28
1	A	191	GLU	C-N-CA	5.05	127.04	120.28
1	B	147	ALA	CA-C-N	5.02	125.42	119.94
1	B	147	ALA	C-N-CA	5.02	125.42	119.94

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	1485	0	1511	24	0
1	B	1504	0	1533	14	0
1	C	1500	0	1530	20	0
1	D	1492	0	1524	26	0
1	E	1420	0	1451	19	0
1	F	1343	0	1366	18	0
2	A	18	0	0	0	0
2	B	13	0	0	0	0
2	C	15	0	0	0	0
2	D	7	0	0	0	0
2	E	7	0	0	0	0
2	F	1	0	0	0	0
All	All	8805	0	8915	107	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 6.

All (107) 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:41:VAL:HG11	1:D:135:VAL:HG13	1.54	0.90
1:F:64:LYS:O	1:F:68:LEU:HG	1.77	0.85
1:A:41:VAL:HG11	1:A:135:VAL:HG13	1.73	0.70
1:D:155:TYR:CZ	1:D:159:ARG:HD2	2.28	0.68
1:E:6:THR:HG22	1:E:9:THR:H	1.62	0.64
1:C:197:GLU:OE1	1:D:159:ARG:NH2	2.31	0.64
1:E:6:THR:HG23	1:F:3:GLY:O	1.99	0.62
1:A:155:TYR:CZ	1:A:159:ARG:HD2	2.36	0.61
1:B:68:LEU:HD22	1:B:124:VAL:HG13	1.83	0.59
1:A:72:LEU:HG	1:A:124:VAL:HG11	1.83	0.59
1:A:72:LEU:HD11	1:A:124:VAL:HG12	1.85	0.59
1:F:133:GLU:O	1:F:137:GLU:HG2	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:163:LEU:O	1:C:167:VAL:HG23	2.07	0.55
1:D:20:THR:HG21	1:D:80:VAL:HG22	1.89	0.54
1:E:75:MET:HA	1:E:78:ASP:HB2	1.90	0.53
1:E:6:THR:CG2	1:F:3:GLY:O	2.57	0.53
1:C:51:TYR:CZ	1:C:140:ASN:HA	2.45	0.52
1:D:68:LEU:HD22	1:D:124:VAL:HG13	1.90	0.52
1:E:202:THR:HG22	1:E:206:LYS:HD2	1.91	0.52
1:C:86:VAL:HG22	1:C:110:ALA:HB1	1.92	0.51
1:F:203:VAL:O	1:F:207:VAL:HG23	2.10	0.51
1:A:44:LEU:HD21	1:B:17:SER:O	2.11	0.50
1:E:66:GLU:HA	1:E:66:GLU:OE2	2.11	0.50
1:A:148:GLY:O	1:A:151:VAL:HG12	2.11	0.50
1:C:207:VAL:HG13	1:D:167:VAL:HG12	1.94	0.49
1:D:113:THR:O	1:D:117:VAL:HG23	2.12	0.49
1:B:6:THR:HG22	1:B:9:THR:H	1.77	0.49
1:D:24:GLY:HA3	1:D:79:ASP:OD1	2.13	0.49
1:D:41:VAL:CG1	1:D:135:VAL:HG13	2.34	0.49
1:A:33:MET:O	1:A:37:LEU:HG	2.13	0.49
1:E:29:ILE:O	1:E:30:SER:C	2.55	0.49
1:D:13:GLY:O	1:D:16:SER:OG	2.27	0.48
1:F:14:LEU:HD12	1:F:21:PRO:O	2.13	0.48
1:E:72:LEU:HG	1:E:124:VAL:HG11	1.96	0.48
1:A:30:SER:O	1:A:153:SER:HB3	2.14	0.48
1:B:72:LEU:HG	1:B:124:VAL:HG11	1.94	0.48
1:B:66:GLU:CD	1:B:69:ARG:HH11	2.20	0.48
1:E:148:GLY:O	1:E:151:VAL:HG12	2.14	0.48
1:D:104:ALA:O	1:D:107:ILE:HB	2.14	0.48
1:B:6:THR:HB	1:B:9:THR:OG1	2.13	0.47
1:F:37:LEU:O	1:F:41:VAL:HG23	2.14	0.47
1:D:42:CYS:HB2	1:D:62:LEU:HD13	1.97	0.47
1:A:41:VAL:CG1	1:A:135:VAL:HG13	2.42	0.47
1:C:72:LEU:HG	1:C:124:VAL:HG11	1.97	0.47
1:E:3:GLY:O	1:F:8:GLU:HB2	2.15	0.47
1:B:126:ARG:NH2	1:B:130:ASP:OD1	2.44	0.47
1:F:197:GLU:O	1:F:201:GLU:HG2	2.15	0.47
1:F:18:ALA:HB1	1:F:19:PRO:HD2	1.96	0.46
1:A:133:GLU:OE1	1:A:137:GLU:OE1	2.33	0.46
1:A:155:TYR:CE1	1:A:159:ARG:HD2	2.50	0.46
1:E:141:LEU:C	1:E:143:VAL:H	2.23	0.46
1:D:120:ALA:HA	1:D:123:ARG:NH1	2.31	0.46
1:C:189:LEU:O	1:C:190:ALA:C	2.59	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:GLU:HA	1:B:66:GLU:OE1	2.15	0.45
1:A:202:THR:O	1:A:206:LYS:HG3	2.17	0.45
1:B:82:ALA:O	1:B:86:VAL:HG23	2.17	0.45
1:C:126:ARG:NH2	1:C:130:ASP:OD1	2.50	0.45
1:F:20:THR:HB	1:F:21:PRO:HA	1.98	0.45
1:D:54:VAL:CG1	1:D:138:LYS:HB3	2.47	0.45
1:D:165:VAL:HG11	1:D:185:LEU:HD22	1.97	0.45
1:C:24:GLY:HA3	1:C:79:ASP:OD1	2.17	0.44
1:B:30:SER:O	1:B:153:SER:HB3	2.17	0.44
1:E:43:ASN:ND2	1:F:15:ALA:O	2.50	0.44
1:A:37:LEU:HD13	1:A:149:VAL:HB	1.99	0.44
1:E:54:VAL:HG21	1:E:138:LYS:HD3	2.00	0.44
1:A:158:LEU:CD2	1:A:189:LEU:HD23	2.48	0.44
1:C:31:GLY:C	1:C:72:LEU:HD13	2.43	0.44
1:E:135:VAL:O	1:E:139:GLY:HA3	2.18	0.44
1:E:189:LEU:O	1:E:192:ALA:N	2.51	0.43
1:B:158:LEU:HD21	1:B:189:LEU:HD23	2.00	0.43
1:A:44:LEU:HD13	1:A:143:VAL:HG11	2.01	0.43
1:D:20:THR:CG2	1:D:80:VAL:HG22	2.49	0.43
1:A:149:VAL:HG21	1:B:23:GLY:HA3	2.01	0.43
1:C:20:THR:HA	1:C:21:PRO:C	2.44	0.42
1:A:2:ALA:N	1:E:81:GLU:OE2	2.53	0.42
1:E:126:ARG:NH2	1:E:130:ASP:OD1	2.36	0.42
1:F:158:LEU:C	1:F:158:LEU:HD23	2.44	0.42
1:F:111:LEU:HD23	1:F:111:LEU:HA	1.94	0.42
1:D:64:LYS:O	1:D:68:LEU:HG	2.19	0.42
1:A:10:PHE:CE2	1:A:29:ILE:CD1	3.02	0.42
1:A:10:PHE:CE2	1:A:29:ILE:HD12	2.54	0.42
1:A:17:SER:HB3	1:B:43:ASN:HB3	2.02	0.42
1:C:10:PHE:CZ	1:C:29:ILE:HD12	2.55	0.42
1:C:7:ILE:O	1:C:11:LEU:HG	2.20	0.42
1:C:124:VAL:O	1:C:127:GLU:HB2	2.20	0.42
1:F:41:VAL:HG11	1:F:135:VAL:HB	2.01	0.42
1:E:19:PRO:HB2	1:E:83:PHE:CZ	2.55	0.42
1:C:159:ARG:HH22	1:D:197:GLU:CD	2.28	0.41
1:D:54:VAL:HG11	1:D:138:LYS:HB3	2.02	0.41
1:A:174:ASP:C	1:A:174:ASP:OD1	2.61	0.41
1:F:37:LEU:CD1	1:F:149:VAL:HG12	2.50	0.41
1:A:151:VAL:HG21	1:A:199:ILE:HB	2.03	0.41
1:D:6:THR:HG22	1:D:8:GLU:N	2.35	0.41
1:E:195:LEU:O	1:E:199:ILE:HG13	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:112:LYS:N	1:D:172:LEU:HD21	2.36	0.41
1:A:165:VAL:HG11	1:A:185:LEU:HD22	2.03	0.41
1:D:175:ARG:NE	1:D:175:ARG:HA	2.36	0.41
1:D:135:VAL:HG12	1:D:147:ALA:HB2	2.02	0.41
1:F:173:ASP:O	1:F:175:ARG:HG2	2.20	0.41
1:C:141:LEU:HD12	1:C:141:LEU:HA	1.94	0.40
1:F:167:VAL:HG12	1:F:168:ASN:OD1	2.21	0.40
1:B:208:ASN:OD1	1:C:184:GLU:CD	2.64	0.40
1:C:36:ALA:O	1:D:15:ALA:HA	2.22	0.40
1:C:37:LEU:HD21	1:D:14:LEU:HD11	2.03	0.40
1:C:140:ASN:O	1:C:143:VAL:HG12	2.21	0.40
1:D:24:GLY:HA3	1:D:79:ASP:CG	2.45	0.40
1:A:92:LEU:HD11	1:A:106:LYS:HD2	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	200/216 (93%)	191 (96%)	9 (4%)	0	100	100
1	B	204/216 (94%)	191 (94%)	13 (6%)	0	100	100
1	C	203/216 (94%)	195 (96%)	7 (3%)	1 (0%)	24	52
1	D	202/216 (94%)	193 (96%)	9 (4%)	0	100	100
1	E	190/216 (88%)	176 (93%)	14 (7%)	0	100	100
1	F	180/216 (83%)	163 (91%)	15 (8%)	2 (1%)	11	32
All	All	1179/1296 (91%)	1109 (94%)	67 (6%)	3 (0%)	36	63

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	138	LYS
1	C	190	ALA
1	F	207	VAL

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	150/162 (93%)	144 (96%)	6 (4%)	28	60
1	B	153/162 (94%)	146 (95%)	7 (5%)	24	55
1	C	153/162 (94%)	143 (94%)	10 (6%)	15	40
1	D	152/162 (94%)	142 (93%)	10 (7%)	15	40
1	E	143/162 (88%)	135 (94%)	8 (6%)	19	46
1	F	136/162 (84%)	129 (95%)	7 (5%)	21	50
All	All	887/972 (91%)	839 (95%)	48 (5%)	20	48

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	SER
1	A	52	VAL
1	A	55	GLU
1	A	135	VAL
1	A	145	SER
1	A	183	LYS
1	B	60	GLN
1	B	66	GLU
1	B	109	GLU
1	B	141	LEU
1	B	142	ASN
1	B	167	VAL
1	B	198	ARG
1	C	16	SER
1	C	29	ILE
1	C	53	GLU

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Mol	Chain	Res	Type
1	C	54	VAL
1	C	140	ASN
1	C	151	VAL
1	C	158	LEU
1	C	186	GLU
1	C	195	LEU
1	C	207	VAL
1	D	30	SER
1	D	65	SER
1	D	73	THR
1	D	94	LYS
1	D	96	THR
1	D	135	VAL
1	D	143	VAL
1	D	153	SER
1	D	167	VAL
1	D	207	VAL
1	E	6	THR
1	E	9	THR
1	E	49	LYS
1	E	75	MET
1	E	117	VAL
1	E	137	GLU
1	E	141	LEU
1	E	145	SER
1	F	30	SER
1	F	86	VAL
1	F	107	ILE
1	F	113	THR
1	F	141	LEU
1	F	145	SER
1	F	184	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	168	ASN
1	C	168	ASN
1	D	168	ASN
1	D	208	ASN
1	E	4	ASN
1	F	43	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	204/216 (94%)	-0.22	1 (0%) 87 83	24, 45, 68, 103	0
1	B	206/216 (95%)	-0.18	0 100 100	31, 49, 73, 106	0
1	C	205/216 (94%)	-0.16	0 100 100	30, 48, 72, 106	0
1	D	204/216 (94%)	-0.11	1 (0%) 87 83	33, 48, 77, 89	0
1	E	196/216 (90%)	0.26	3 (1%) 72 65	39, 66, 95, 107	0
1	F	186/216 (86%)	0.39	3 (1%) 70 63	42, 70, 98, 108	0
All	All	1201/1296 (92%)	-0.01	8 (0%) 84 79	24, 52, 90, 108	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	169	ALA	2.5
1	A	2	ALA	2.2
1	E	207	VAL	2.2
1	F	134	ILE	2.2
1	E	203	VAL	2.1
1	D	208	ASN	2.1
1	E	137	GLU	2.1
1	F	138	LYS	2.1

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.