Full wwPDB X-ray Structure Validation Report ⓘ

Sep 2, 2023 – 11:52 AM EDT

PDB ID	:	3R8Y
Title	:	Structure of the Bacillus anthracis tetrahydropicolinate succinyltransferase
Authors	:	Anderson, S.M.; Wawrzak, Z.; Onopriyenko, O.; Peterson, S.N.; Anderson, W.F.; Savchenko, A.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on	:	2011-03-24
Resolution	:	1.70 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:

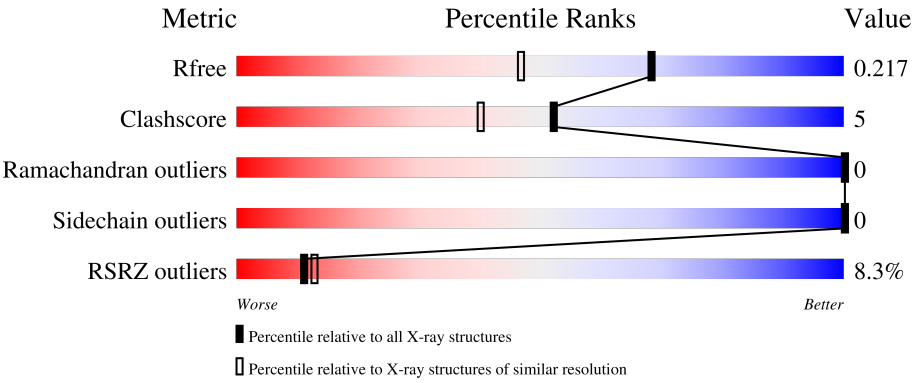
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	130704	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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 <=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	240	6% 79% 5% 15%
1	B	240	11% 81% 8% 11%
1	C	240	4% 73% 11% 16%
1	D	240	5% 78% 7% 15%
1	E	240	12% 81% 8% 12%

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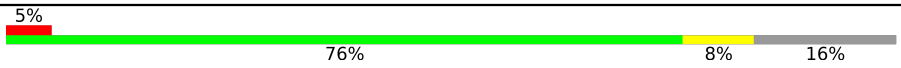
Mol	Chain	Length	Quality of chain
1	F	240	

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 10874 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 2,3,4,5-tetrahydropyridine-2,6-dicarboxylate N-acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	203	Total	C	N	O	S	0	1	0
			1501	955	253	287	6			
1	B	213	Total	C	N	O	S	0	0	0
			1574	1001	265	302	6			
1	C	201	Total	C	N	O	S	0	2	0
			1490	951	251	282	6			
1	D	203	Total	C	N	O	S	0	1	0
			1501	955	253	287	6			
1	E	212	Total	C	N	O	S	0	1	0
			1570	998	264	302	6			
1	F	202	Total	C	N	O	S	0	1	0
			1493	951	251	285	6			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Ca	0	0
			2	2		
2	B	1	Total	Ca	0	0
			1	1		
2	C	1	Total	Ca	0	0
			1	1		
2	D	2	Total	Ca	0	0
			2	2		
2	F	3	Total	Ca	0	0
			3	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	263	Total	O	0	2
			265	265		

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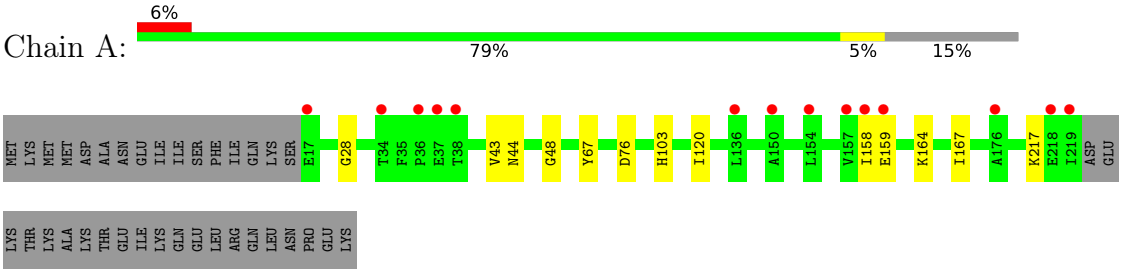
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	306	Total O 309 309	0	3
3	C	294	Total O 295 295	0	1
3	D	287	Total O 288 288	0	1
3	E	278	Total O 279 279	0	1
3	F	300	Total O 300 300	0	0

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.

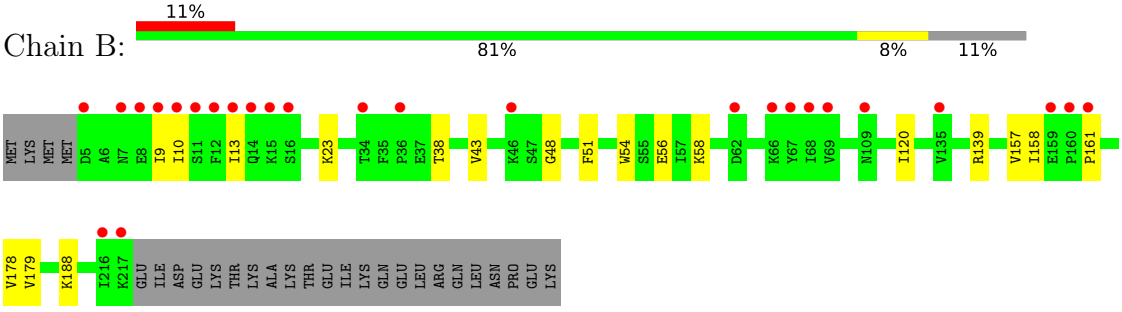
• Molecule 1: 2,3,4,5-tetrahydropyridine-2,6-dicarboxylate N-acetyltransferase

Chain A:



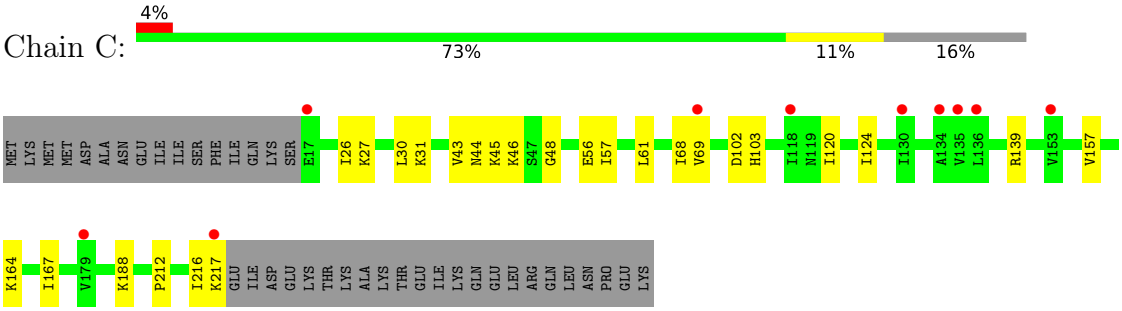
• Molecule 1: 2,3,4,5-tetrahydropyridine-2,6-dicarboxylate N-acetyltransferase

Chain B:



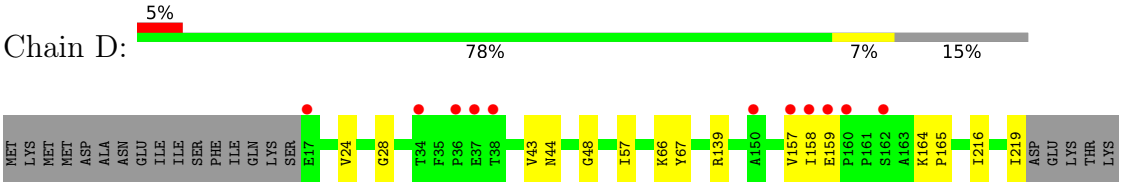
• Molecule 1: 2,3,4,5-tetrahydropyridine-2,6-dicarboxylate N-acetyltransferase

Chain C:



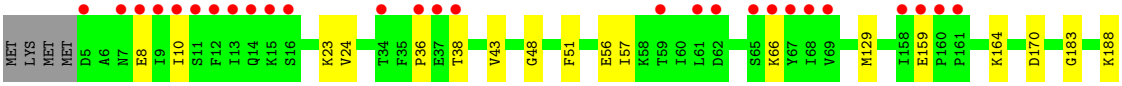
• Molecule 1: 2,3,4,5-tetrahydropyridine-2,6-dicarboxylate N-acetyltransferase

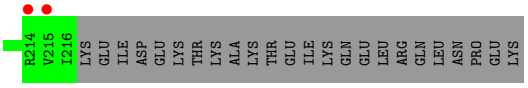
Chain D:



• Molecule 1: 2,3,4,5-tetrahydropyridine-2,6-dicarboxylate N-acetyltransferase

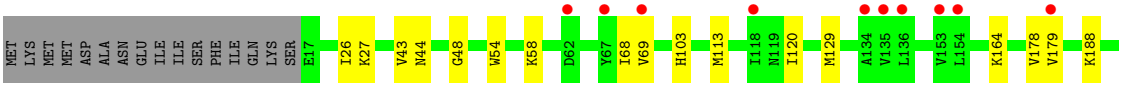
Chain E:

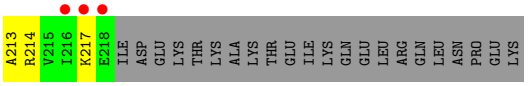




• Molecule 1: 2,3,4,5-tetrahydropyridine-2,6-dicarboxylate N-acetyltransferase

Chain F:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	52.42Å 52.22Å 134.55Å 97.41° 93.19° 115.71°	Depositor
Resolution (Å)	29.41 – 1.70 41.42 – 1.67	Depositor EDS
% Data completeness (in resolution range)	99.3 (29.41-1.70) 97.1 (41.42-1.67)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.04 (at 1.67Å)	Xtriage
Refinement program	PHENIX 1.7_650	Depositor
R, R_{free}	0.186 , 0.221 0.180 , 0.217	Depositor DCC
R_{free} test set	7127 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	18.0	Xtriage
Anisotropy	0.260	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 62.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.073 for k,h,-h-k-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	10874	wwPDB-VP
Average B, all atoms (Å^2)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.12% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: CA

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	0.57	0/1524	0.59	0/2072
1	B	0.44	0/1595	0.57	0/2167
1	C	0.47	0/1516	0.60	0/2061
1	D	0.48	0/1524	0.59	0/2072
1	E	0.36	0/1594	0.54	0/2167
1	F	0.37	0/1516	0.55	0/2061
All	All	0.45	0/9269	0.57	0/12600

There are no bond length outliers.

There are no bond angle outliers.

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	1501	0	1564	12	0
1	B	1574	0	1635	15	0
1	C	1490	0	1565	26	0
1	D	1501	0	1564	20	0
1	E	1570	0	1628	18	0
1	F	1493	0	1558	18	0
2	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	2	0	0	0	0
2	F	3	0	0	0	0
3	A	265	0	0	1	0
3	B	309	0	0	6	0
3	C	295	0	0	6	0
3	D	288	0	0	3	0
3	E	279	0	0	6	0
3	F	300	0	0	3	0
All	All	10874	0	9514	102	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 5.

All (102) 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:C:43:VAL:HG12	1:C:48:GLY:HA3	1.54	0.90
1:A:43:VAL:HG12	1:A:48:GLY:HA3	1.60	0.83
1:D:157:VAL:HG11	1:D:165:PRO:HA	1.66	0.78
1:D:157:VAL:O	1:D:157:VAL:HG13	1.86	0.76
1:E:164:LYS:HE3	1:E:183:GLY:O	1.89	0.72
1:B:43:VAL:HG12	1:B:48:GLY:HA3	1.71	0.71
1:C:26[A]:ILE:HD11	1:C:68:ILE:HD13	1.73	0.68
1:A:44:ASN:HB3	1:C:44:ASN:HD21	1.59	0.68
1:A:158:ILE:HG22	1:A:159:GLU:HG2	1.75	0.68
1:B:10:ILE:CD1	1:B:161:PRO:HD3	2.25	0.67
1:D:157:VAL:HG11	1:D:165:PRO:CA	2.25	0.65
1:B:188:LYS:HE2	3:B:713:HOH:O	1.97	0.65
1:D:44:ASN:HD22	1:F:44:ASN:ND2	1.96	0.64
1:C:103:HIS:HB2	1:C:120:ILE:HG13	1.81	0.62
1:E:43:VAL:HG12	1:E:48:GLY:HA3	1.80	0.62
1:E:66:LYS:HG2	3:E:1690:HOH:O	1.99	0.62
1:C:43:VAL:CG1	1:C:48:GLY:HA3	2.27	0.61
1:D:139:ARG:NH1	1:D:157:VAL:HG13	2.14	0.61
1:E:8:GLU:HG2	3:E:1492:HOH:O	2.00	0.61
1:D:158:ILE:HG22	1:D:159:GLU:HG3	1.82	0.61
1:C:27:LYS:HG3	1:C:69:VAL:CG2	2.31	0.61
1:F:213:ALA:O	1:F:214:ARG:HD3	2.00	0.60
1:D:157:VAL:CG1	1:D:165:PRO:HA	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:VAL:HG23	1:B:157:VAL:O	2.02	0.59
1:C:216:ILE:HD13	3:C:1719:HOH:O	2.02	0.59
1:D:44:ASN:ND2	1:F:44:ASN:ND2	2.52	0.58
1:E:10:ILE:CD1	1:E:159:GLU:HG2	2.34	0.57
1:F:27:LYS:HG3	1:F:69:VAL:CG2	2.34	0.57
1:F:103:HIS:HB2	1:F:120:ILE:HG13	1.89	0.55
1:B:9:ILE:O	1:B:13:ILE:HG13	2.06	0.55
1:B:23:LYS:HE2	1:B:51:PHE:CE1	2.42	0.54
1:F:43:VAL:HG12	1:F:48:GLY:HA3	1.88	0.54
1:D:44:ASN:HD22	1:F:44:ASN:HD21	1.54	0.54
1:E:164:LYS:NZ	3:E:955:HOH:O	2.28	0.54
1:F:113:MET:HG3	1:F:129:MET:HE2	1.89	0.54
1:F:188:LYS:HE2	3:F:974:HOH:O	2.08	0.53
1:C:164[B]:LYS:HE3	3:C:431:HOH:O	2.07	0.53
1:C:31:LYS:HE2	1:C:45:LYS:HG3	1.90	0.53
1:C:45:LYS:HG3	3:C:1193:HOH:O	2.07	0.53
1:A:164:LYS:HE2	1:A:167:ILE:HD12	1.90	0.53
1:A:28:GLY:HA3	1:A:67:TYR:O	2.09	0.52
1:B:158:ILE:HB	3:B:1732:HOH:O	2.09	0.52
1:D:157:VAL:O	1:D:157:VAL:CG1	2.58	0.51
1:E:66:LYS:CG	3:E:1690:HOH:O	2.57	0.51
1:A:158:ILE:C	1:A:159:GLU:HG2	2.32	0.50
1:C:188:LYS:HE2	3:C:597:HOH:O	2.11	0.50
1:B:9:ILE:HD12	1:B:120:ILE:HG12	1.93	0.50
1:C:164[B]:LYS:HE2	1:C:167:ILE:HD12	1.92	0.50
1:F:26[A]:ILE:HD11	1:F:68:ILE:HG12	1.94	0.50
1:E:23:LYS:HG2	1:E:51:PHE:CE1	2.46	0.50
1:E:188:LYS:HE2	3:E:956:HOH:O	2.11	0.50
1:C:27:LYS:HG3	1:C:69:VAL:HG22	1.94	0.50
1:F:27:LYS:HG3	1:F:69:VAL:HG23	1.94	0.49
1:E:164:LYS:CE	1:E:183:GLY:O	2.61	0.48
1:B:54:TRP:CZ2	1:B:58:LYS:HD3	2.49	0.48
1:E:10:ILE:HD11	1:E:159:GLU:HG2	1.95	0.48
1:F:178:VAL:CG1	1:F:179:VAL:N	2.77	0.48
1:B:10:ILE:HD11	1:B:161:PRO:HD3	1.96	0.47
1:E:36:PRO:HB2	1:E:38:THR:HG22	1.95	0.47
1:F:27:LYS:O	1:F:69:VAL:HG22	2.13	0.47
1:C:212:PRO:HD3	1:D:216:ILE:HD12	1.97	0.47
1:A:44:ASN:HB3	1:C:44:ASN:ND2	2.26	0.47
1:B:139:ARG:HD2	3:B:978:HOH:O	2.15	0.47
1:D:43:VAL:HG12	1:D:48:GLY:HA3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:216:ILE:HG22	1:C:217:LYS:HG3	1.98	0.46
1:C:57:ILE:O	1:C:61:LEU:HG	2.16	0.46
1:A:103:HIS:HB2	1:A:120:ILE:HG13	1.97	0.45
1:C:124:ILE:HD12	1:C:124:ILE:N	2.32	0.45
1:F:129:MET:HE1	3:F:279:HOH:O	2.17	0.45
1:B:38:THR:OG1	1:B:56:GLU:HG2	2.17	0.45
1:E:38:THR:CG2	1:E:56:GLU:HG2	2.47	0.44
1:C:102:ASP:O	1:C:103:HIS:HB2	2.18	0.44
1:C:139:ARG:HD3	1:C:157:VAL:O	2.18	0.44
1:E:24:VAL:HG21	1:E:57:ILE:HG21	2.00	0.44
1:F:164:LYS:HD3	3:F:1553:HOH:O	2.18	0.44
1:A:217:LYS:HD3	1:D:219:ILE:CD1	2.49	0.43
1:C:27:LYS:NZ	3:C:564:HOH:O	2.50	0.43
1:A:158:ILE:O	1:A:159:GLU:HG2	2.18	0.43
1:D:157:VAL:CG1	1:D:165:PRO:HB3	2.49	0.42
1:E:10:ILE:HD13	1:E:159:GLU:HG2	1.99	0.42
1:C:26[A]:ILE:HD11	1:C:68:ILE:CD1	2.44	0.42
1:D:66:LYS:HG2	3:D:1655:HOH:O	2.19	0.42
1:D:28:GLY:HA3	1:D:67:TYR:O	2.18	0.42
1:D:24:VAL:HG21	1:D:57:ILE:HG21	2.01	0.42
1:D:219:ILE:HG22	3:D:1093:HOH:O	2.18	0.42
1:A:44:ASN:HB2	3:A:1584:HOH:O	2.19	0.42
1:A:76:ASP:HB2	3:B:575:HOH:O	2.20	0.42
1:B:158:ILE:CD1	3:B:1518[A]:HOH:O	2.68	0.41
1:E:170:ASP:HB2	1:E:188:LYS:HE3	2.03	0.41
1:C:43:VAL:HG12	1:C:48:GLY:CA	2.39	0.41
1:B:178:VAL:CG1	1:B:179:VAL:N	2.83	0.41
1:C:56:GLU:HG3	3:C:1057:HOH:O	2.20	0.41
1:D:164:LYS:HE3	3:D:820:HOH:O	2.20	0.41
1:C:26[B]:ILE:HD12	1:C:30:LEU:HD22	2.01	0.41
1:B:158:ILE:HD12	3:B:904:HOH:O	2.21	0.41
1:F:54:TRP:CZ2	1:F:58:LYS:HD3	2.55	0.41
1:D:139:ARG:HH11	1:D:157:VAL:HG13	1.82	0.41
1:C:46:LYS:HE2	1:C:46:LYS:HB3	1.94	0.41
1:E:23:LYS:HG2	1:E:51:PHE:CD1	2.56	0.40
1:E:129:MET:HE1	3:E:1307:HOH:O	2.22	0.40
1:F:217:LYS:NZ	1:F:217:LYS:HB3	2.35	0.40
1:F:113:MET:HG3	1:F:129:MET:CE	2.51	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	202/240 (84%)	200 (99%)	2 (1%)	0	100	100
1	B	211/240 (88%)	209 (99%)	2 (1%)	0	100	100
1	C	201/240 (84%)	200 (100%)	1 (0%)	0	100	100
1	D	202/240 (84%)	201 (100%)	1 (0%)	0	100	100
1	E	211/240 (88%)	210 (100%)	1 (0%)	0	100	100
1	F	201/240 (84%)	199 (99%)	2 (1%)	0	100	100
All	All	1228/1440 (85%)	1219 (99%)	9 (1%)	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	163/197 (83%)	163 (100%)	0	100	100
1	B	171/197 (87%)	171 (100%)	0	100	100
1	C	162/197 (82%)	162 (100%)	0	100	100
1	D	163/197 (83%)	163 (100%)	0	100	100
1	E	171/197 (87%)	171 (100%)	0	100	100
1	F	162/197 (82%)	162 (100%)	0	100	100
All	All	992/1182 (84%)	992 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	103	HIS
1	D	44	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry

Of 9 ligands modelled in this entry, 9 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

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^{th} 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	203/240 (84%)	0.45	14 (6%) 16 19	9, 19, 47, 58	0
1	B	213/240 (88%)	0.55	26 (12%) 4 4	9, 20, 52, 80	0
1	C	201/240 (83%)	0.29	10 (4%) 28 32	9, 17, 43, 52	0
1	D	203/240 (84%)	0.40	11 (5%) 25 28	9, 19, 48, 62	0
1	E	212/240 (88%)	0.60	29 (13%) 3 3	10, 21, 54, 96	0
1	F	202/240 (84%)	0.30	13 (6%) 19 21	9, 18, 42, 70	0
All	All	1234/1440 (85%)	0.43	103 (8%) 11 13	9, 20, 50, 96	0

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	12	PHE	6.1
1	B	13	ILE	6.0
1	B	9	ILE	5.7
1	E	12	PHE	5.5
1	E	13	ILE	5.4
1	B	10	ILE	5.0
1	E	9	ILE	5.0
1	E	10	ILE	4.7
1	B	161	PRO	4.6
1	B	15	LYS	4.5
1	B	11	SER	4.4
1	E	16	SER	4.4
1	E	161	PRO	4.4
1	B	5	ASP	4.4
1	E	14	GLN	4.3
1	A	158	ILE	4.2
1	F	218	GLU	4.2
1	E	11	SER	4.2
1	E	160	PRO	4.1

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Mol	Chain	Res	Type	RSRZ
1	E	38	THR	4.1
1	F	217	LYS	4.1
1	B	14	GLN	3.9
1	E	7	ASN	3.9
1	E	67	TYR	3.9
1	B	69	VAL	3.8
1	B	34	THR	3.8
1	B	7	ASN	3.6
1	A	219	ILE	3.6
1	D	37	GLU	3.6
1	B	16	SER	3.5
1	B	160	PRO	3.5
1	D	159	GLU	3.5
1	E	5	ASP	3.4
1	B	68	ILE	3.3
1	E	36	PRO	3.3
1	E	8	GLU	3.3
1	B	67	TYR	3.3
1	E	66	LYS	3.3
1	E	62	ASP	3.3
1	D	158	ILE	3.3
1	D	36	PRO	3.2
1	D	150	ALA	3.2
1	E	159	GLU	3.2
1	F	135	VAL	3.1
1	F	179	VAL	3.1
1	F	153	VAL	3.1
1	A	34	THR	3.0
1	E	158	ILE	3.0
1	A	218	GLU	3.0
1	D	38	THR	3.0
1	D	17	GLU	2.8
1	E	214	ARG	2.8
1	A	37	GLU	2.8
1	C	153	VAL	2.8
1	D	157	VAL	2.8
1	E	59	THR	2.8
1	C	179	VAL	2.7
1	E	215	VAL	2.7
1	A	159	GLU	2.7
1	E	15	LYS	2.7
1	B	36	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	217	LYS	2.7
1	A	38	THR	2.7
1	E	68	ILE	2.6
1	E	65	SER	2.6
1	F	69	VAL	2.6
1	C	134	ALA	2.6
1	F	62	ASP	2.6
1	E	37	GLU	2.5
1	E	34	THR	2.5
1	F	118	ILE	2.5
1	D	34	THR	2.5
1	B	62	ASP	2.5
1	B	8	GLU	2.5
1	F	67	TYR	2.5
1	D	162	SER	2.4
1	B	135	VAL	2.3
1	A	36	PRO	2.3
1	F	136	LEU	2.3
1	F	216	ILE	2.3
1	A	157	VAL	2.3
1	F	134	ALA	2.3
1	A	17	GLU	2.3
1	C	69	VAL	2.3
1	A	154	LEU	2.3
1	C	136	LEU	2.3
1	E	61	LEU	2.2
1	B	66	LYS	2.2
1	B	217	LYS	2.2
1	B	46	LYS	2.1
1	C	118	ILE	2.1
1	C	130	ILE	2.1
1	C	135	VAL	2.1
1	A	136	LEU	2.1
1	B	159	GLU	2.1
1	F	154	LEU	2.1
1	B	216	ILE	2.1
1	E	69	VAL	2.1
1	C	17	GLU	2.1
1	A	176	ALA	2.0
1	D	160	PRO	2.0
1	A	150	ALA	2.0
1	B	109	ASN	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 monosaccharides 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, 95^{th} 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(Å^2)	Q<0.9
2	CA	B	241	1/1	0.87	0.07	68,68,68,68	0
2	CA	F	243	1/1	0.88	0.10	59,59,59,59	0
2	CA	F	242	1/1	0.91	0.07	54,54,54,54	0
2	CA	A	242	1/1	0.98	0.07	23,23,23,23	0
2	CA	D	241	1/1	0.98	0.07	23,23,23,23	0
2	CA	D	242	1/1	0.99	0.05	23,23,23,23	0
2	CA	F	241	1/1	1.00	0.06	19,19,19,19	0
2	CA	A	241	1/1	1.00	0.06	21,21,21,21	0
2	CA	C	241	1/1	1.00	0.06	19,19,19,19	0

6.5 Other polymers (i)

There are no such residues in this entry.