Full wwPDB X-ray Structure Validation Report (i)

Jan 30, 2021 – 06:28 PM EST

PDB ID	:	3KKE
Title	:	Crystal structure of a LacI family transcriptional regulator from Mycobacterium smegmatis
Authors	:	Bonanno, J.B.; Rutter, M.; Bain, K.T.; Do, J.; Ozyurt, S.; Wasserman, S.; Sauder, J.M.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on	:	2009-11-05
Resolution	:	2.20 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

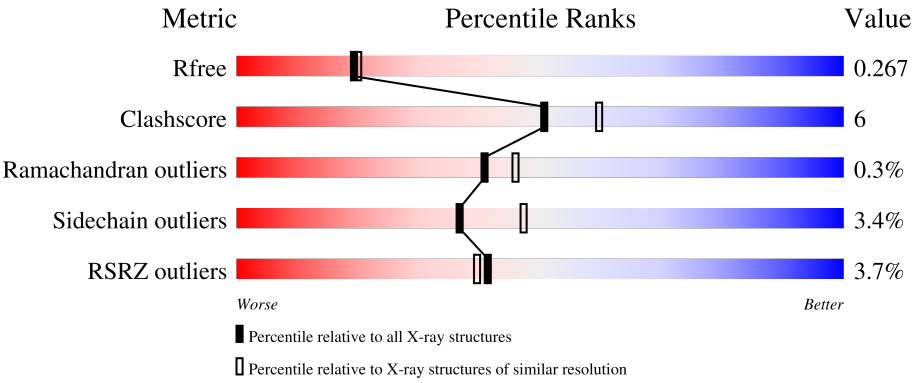
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.16
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.16

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 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	303	% 79% 12% • 8%
1	B	303	86% 10% • •
1	C	303	3% 84% 10% • 6%
1	D	303	10% 74% 16% • 9%

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
2	ACT	A	350	-	-	X	-
2	ACT	D	1	-	-	X	-

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 8251 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 LacI family Transcriptional regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	278	Total	C	N	O	S	0	0	0
			1992	1246	348	394	4			
1	B	291	Total	C	N	O	S	0	0	0
			2096	1307	372	413	4			
1	C	285	Total	C	N	O	S	0	0	0
			2046	1277	360	405	4			
1	D	277	Total	C	N	O	S	0	0	0
			1970	1238	345	383	4			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	MET	-	expression tag	UNP A0QPR6
A	48	SER	-	expression tag	UNP A0QPR6
A	49	LEU	-	expression tag	UNP A0QPR6
A	342	GLU	-	expression tag	UNP A0QPR6
A	343	GLY	-	expression tag	UNP A0QPR6
A	344	HIS	-	expression tag	UNP A0QPR6
A	345	HIS	-	expression tag	UNP A0QPR6
A	346	HIS	-	expression tag	UNP A0QPR6
A	347	HIS	-	expression tag	UNP A0QPR6
A	348	HIS	-	expression tag	UNP A0QPR6
A	349	HIS	-	expression tag	UNP A0QPR6
B	47	MET	-	expression tag	UNP A0QPR6
B	48	SER	-	expression tag	UNP A0QPR6
B	49	LEU	-	expression tag	UNP A0QPR6
B	342	GLU	-	expression tag	UNP A0QPR6
B	343	GLY	-	expression tag	UNP A0QPR6
B	344	HIS	-	expression tag	UNP A0QPR6
B	345	HIS	-	expression tag	UNP A0QPR6
B	346	HIS	-	expression tag	UNP A0QPR6
B	347	HIS	-	expression tag	UNP A0QPR6
B	348	HIS	-	expression tag	UNP A0QPR6

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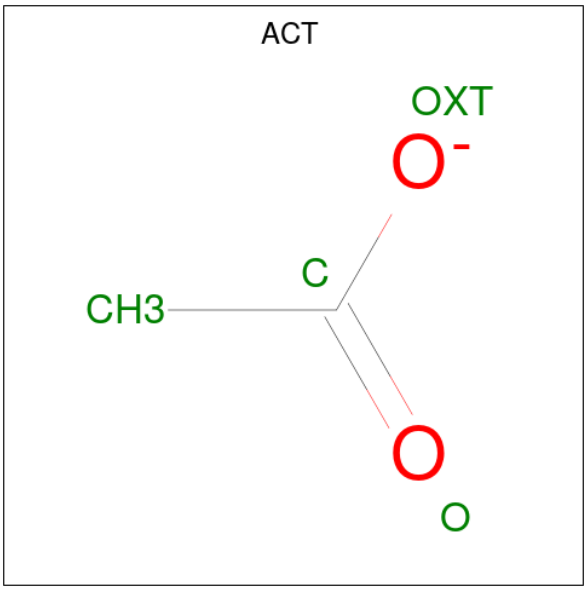
Chain	Residue	Modelled	Actual	Comment	Reference
B	349	HIS	-	expression tag	UNP A0QPR6
C	47	MET	-	expression tag	UNP A0QPR6
C	48	SER	-	expression tag	UNP A0QPR6
C	49	LEU	-	expression tag	UNP A0QPR6
C	342	GLU	-	expression tag	UNP A0QPR6
C	343	GLY	-	expression tag	UNP A0QPR6
C	344	HIS	-	expression tag	UNP A0QPR6
C	345	HIS	-	expression tag	UNP A0QPR6
C	346	HIS	-	expression tag	UNP A0QPR6
C	347	HIS	-	expression tag	UNP A0QPR6
C	348	HIS	-	expression tag	UNP A0QPR6
C	349	HIS	-	expression tag	UNP A0QPR6
D	47	MET	-	expression tag	UNP A0QPR6
D	48	SER	-	expression tag	UNP A0QPR6
D	49	LEU	-	expression tag	UNP A0QPR6
D	342	GLU	-	expression tag	UNP A0QPR6
D	343	GLY	-	expression tag	UNP A0QPR6
D	344	HIS	-	expression tag	UNP A0QPR6
D	345	HIS	-	expression tag	UNP A0QPR6
D	346	HIS	-	expression tag	UNP A0QPR6
D	347	HIS	-	expression tag	UNP A0QPR6
D	348	HIS	-	expression tag	UNP A0QPR6
D	349	HIS	-	expression tag	UNP A0QPR6

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0

- Molecule 3 is water.

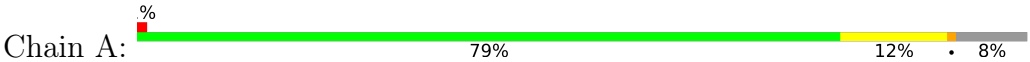
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	44	Total O 44 44	0	0
3	B	60	Total O 60 60	0	0
3	C	19	Total O 19 19	0	0
3	D	4	Total O 4 4	0	0

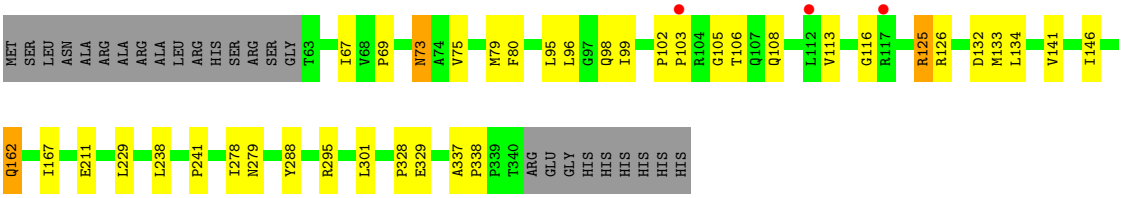
3 Residue-property plots (i)

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: LacI family Transcriptional regulator

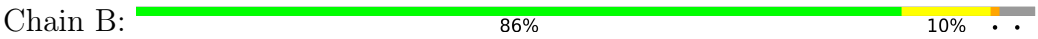
Chain A:

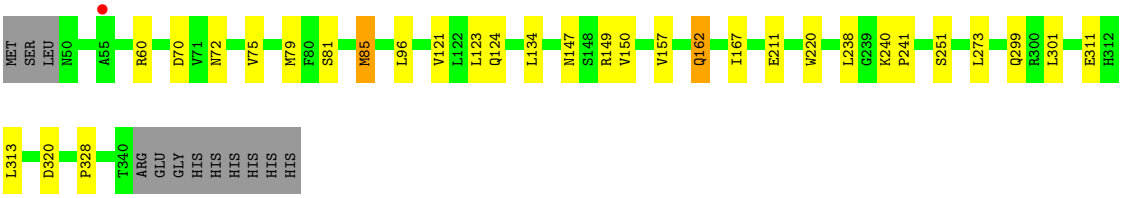




• Molecule 1: LacI family Transcriptional regulator

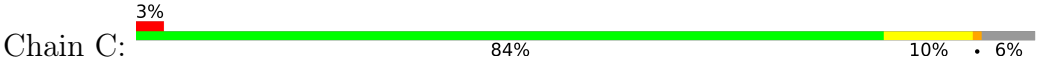
Chain B:

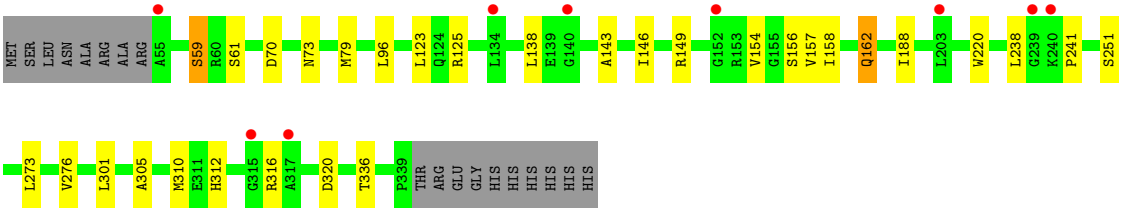




• Molecule 1: LacI family Transcriptional regulator

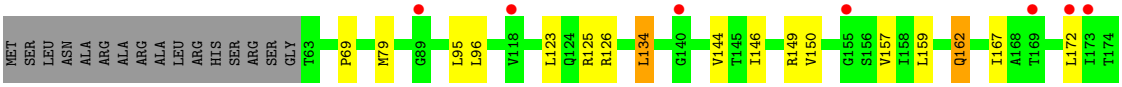
Chain C:

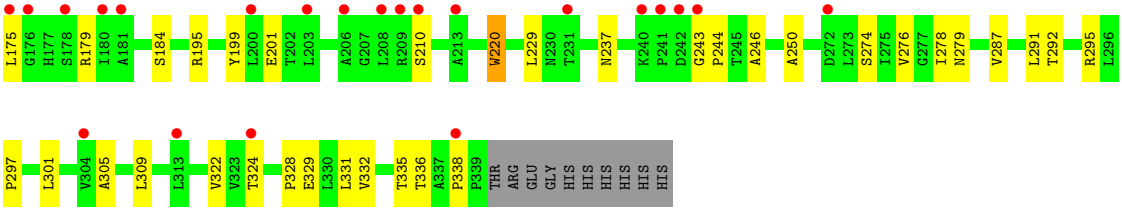




• Molecule 1: LacI family Transcriptional regulator

Chain D:




L175
G176
H177
S178
R179
I180
A181
S184
R195
Y199
L200
E201
T202
L203
A206
G207
L208
R209
S210
A213
W220
L229
N230
T231
N237
K240
P241
D242
G243
P244
T245
A246
A250
D272
L273
S274
I275
V276
G277
I278
N279
V287
L291
T292
R295
L296
P297
L301
V304
A305
L309
L313
V322
V323
T324
P328
E329
L330
L331
V332
T335
T336
A337
P338
P339
THR
ARG
GLU
GLY
HIS
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	64.60Å 65.47Å 116.03Å 90.00° 90.72° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20 33.37 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.2 (20.00-2.20) 99.2 (33.37-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	4.45 (at 2.20Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.194 , 0.247 0.214 , 0.267	Depositor DCC
R_{free} test set	2460 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	34.7	Xtriage
Anisotropy	0.464	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 44.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.011 for -k,-h,-l 0.009 for k,h,-l 0.032 for h,-k,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	8251	wwPDB-VP
Average B, all atoms (Å^2)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.65% 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: ACT

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	0.73	0/2022	0.78	0/2772
1	B	0.80	0/2127	0.84	5/2911 (0.2%)
1	C	0.64	0/2077	0.72	0/2844
1	D	0.54	0/2000	0.63	0/2743
All	All	0.69	0/8226	0.75	5/11270 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	149	ARG	NE-CZ-NH1	6.66	123.63	120.30
1	B	79	MET	CG-SD-CE	-5.99	90.62	100.20
1	B	149	ARG	NE-CZ-NH2	-5.90	117.35	120.30
1	B	320	ASP	CB-CG-OD1	5.78	123.50	118.30
1	B	85	MET	CA-CB-CG	5.05	121.88	113.30

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	1992	0	1968	25	0
1	B	2096	0	2089	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2046	0	2021	18	0
1	D	1970	0	1949	35	0
2	A	8	0	6	3	0
2	B	4	0	3	0	0
2	C	4	0	3	0	0
2	D	4	0	3	2	0
3	A	44	0	0	0	0
3	B	60	0	0	0	0
3	C	19	0	0	0	0
3	D	4	0	0	0	0
All	All	8251	0	8042	92	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 (92) 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:162:GLN:HE21	1:D:162:GLN:H	1.07	0.98
1:C:162:GLN:H	1:C:162:GLN:HE21	1.18	0.90
1:B:134:LEU:HD23	1:B:150:VAL:HG22	1.52	0.89
2:A:350:ACT:H1	1:D:287:VAL:HB	1.57	0.87
1:B:162:GLN:HE21	1:B:162:GLN:H	1.22	0.87
1:A:162:GLN:H	1:A:162:GLN:HE21	1.23	0.85
1:A:105:GLY:HA2	1:A:108:GLN:NE2	1.96	0.80
1:A:102:PRO:HG3	1:A:133:MET:HE1	1.65	0.78
1:D:79:MET:SD	1:D:146:ILE:HD12	2.28	0.73
1:D:95:LEU:HD23	1:D:95:LEU:C	2.07	0.73
1:D:162:GLN:NE2	1:D:162:GLN:H	1.87	0.72
1:D:295:ARG:NH1	1:D:331:LEU:HD11	2.10	0.67
1:D:69:PRO:HB3	1:D:126:ARG:HG3	1.77	0.66
1:C:79:MET:SD	1:C:146:ILE:HD12	2.37	0.65
1:B:96:LEU:HD22	1:C:96:LEU:HD22	1.80	0.64
1:D:123:LEU:HD22	1:D:134:LEU:HD21	1.81	0.63
1:B:238:LEU:C	1:B:238:LEU:HD13	2.20	0.62
1:D:179:ARG:CB	1:D:244:PRO:HA	2.30	0.62
1:B:134:LEU:CD2	1:B:150:VAL:HG22	2.27	0.61
1:D:172:LEU:HD13	1:D:246:ALA:HB1	1.82	0.61
1:A:162:GLN:H	1:A:162:GLN:NE2	1.97	0.60
1:A:105:GLY:HA2	1:A:108:GLN:HE21	1.64	0.59
1:B:70:ASP:OD2	1:B:72:ASN:HB2	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:162:GLN:H	1:B:162:GLN:NE2	1.97	0.59
2:A:350:ACT:H1	1:D:287:VAL:CB	2.33	0.58
1:A:73:ASN:OD1	1:A:75:VAL:HG22	2.03	0.57
1:D:134:LEU:HD13	1:D:150:VAL:HG22	1.87	0.56
1:D:220:TRP:CE2	2:D:1:ACT:H2	2.40	0.56
1:A:125:ARG:HB2	1:A:134:LEU:HD11	1.90	0.54
1:D:322:VAL:HG12	1:D:324:THR:HG23	1.89	0.54
1:A:125:ARG:HG3	1:A:125:ARG:O	2.08	0.53
1:A:98:GLN:C	1:A:99:ILE:HD12	2.29	0.53
1:A:337:ALA:HB1	1:A:338:PRO:HD2	1.91	0.52
1:A:96:LEU:HD22	1:D:96:LEU:HD22	1.91	0.52
1:D:162:GLN:HE21	1:D:162:GLN:N	1.91	0.52
1:C:146:ILE:HG22	1:C:305:ALA:CB	2.39	0.52
1:D:95:LEU:C	1:D:95:LEU:CD2	2.76	0.52
1:D:295:ARG:NH2	1:D:329:GLU:OE1	2.43	0.52
1:D:95:LEU:HD23	1:D:96:LEU:N	2.25	0.51
1:D:69:PRO:HB3	1:D:126:ARG:CG	2.39	0.51
1:C:123:LEU:HD22	1:C:138:LEU:HD21	1.93	0.50
1:B:167:ILE:HD12	1:B:328:PRO:HB2	1.93	0.50
1:A:105:GLY:HA2	1:A:108:GLN:HE22	1.77	0.49
1:C:138:LEU:HD22	1:C:143:ALA:CB	2.42	0.49
1:C:70:ASP:O	1:C:73:ASN:ND2	2.46	0.49
1:D:159:LEU:HD22	1:D:297:PRO:HD2	1.94	0.49
1:B:81:SER:OG	1:B:299:GLN:NE2	2.43	0.48
1:D:175:LEU:HD22	1:D:338:PRO:HD3	1.94	0.48
1:A:103:PRO:O	1:A:106:THR:HG22	2.14	0.48
1:B:157:VAL:HG12	1:B:301:LEU:HD12	1.95	0.48
1:D:279:ASN:ND2	2:D:1:ACT:OXT	2.46	0.48
1:A:69:PRO:HB3	1:A:126:ARG:HG2	1.95	0.48
1:B:238:LEU:HD21	1:B:273:LEU:CD1	2.44	0.47
1:D:278:ILE:O	1:D:279:ASN:HB2	2.14	0.47
1:B:123:LEU:C	1:B:123:LEU:HD23	2.35	0.47
1:D:167:ILE:HD12	1:D:328:PRO:HB2	1.95	0.47
1:B:124:GLN:NE2	1:B:147:ASN:HD22	2.12	0.47
1:C:220:TRP:CE3	1:C:251:SER:HB2	2.49	0.46
1:D:157:VAL:O	1:D:301:LEU:HD11	2.15	0.46
1:D:144:VAL:HG23	1:D:309:LEU:HD13	1.97	0.46
1:B:75:VAL:HG21	1:B:124:GLN:HE22	1.80	0.46
1:A:113:VAL:CG1	1:A:141:VAL:HG21	2.46	0.46
1:C:238:LEU:HD11	1:C:273:LEU:HD11	1.98	0.46
1:D:291:LEU:HD23	1:D:335:THR:HG21	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:162:GLN:H	1:C:162:GLN:NE2	2.00	0.45
1:C:154:VAL:HG23	1:C:320:ASP:OD2	2.17	0.45
1:B:60:ARG:NH2	1:B:313:LEU:O	2.49	0.44
1:A:105:GLY:CA	1:A:108:GLN:HE21	2.29	0.44
1:B:220:TRP:CE3	1:B:251:SER:HB2	2.52	0.44
1:C:310:MET:HE3	1:C:310:MET:HB2	1.92	0.44
1:C:149:ARG:HG2	1:C:156:SER:OG	2.18	0.43
1:A:79:MET:HE2	1:A:301:LEU:HD23	1.99	0.43
1:A:278:ILE:O	1:A:279:ASN:HB2	2.18	0.43
1:C:59:SER:OG	1:C:61:SER:HB2	2.18	0.43
1:C:276:VAL:HG22	1:C:336:THR:HG21	2.01	0.43
1:D:237:ASN:O	1:D:243:GLY:HA3	2.18	0.43
1:D:292:THR:HG23	1:D:332:VAL:HA	2.00	0.43
1:D:184:SER:HA	1:D:250:ALA:HB2	2.01	0.42
1:C:312:HIS:HA	1:C:316:ARG:O	2.19	0.42
1:A:167:ILE:CD1	1:A:328:PRO:HB2	2.50	0.42
1:B:240:LYS:HB3	1:B:241:PRO:CD	2.49	0.42
1:D:157:VAL:HG21	1:D:305:ALA:HA	2.01	0.42
1:A:67:ILE:HG23	1:A:99:ILE:HD13	2.02	0.42
1:A:105:GLY:CA	1:A:108:GLN:NE2	2.77	0.41
1:A:288:TYR:H	2:A:350:ACT:C	2.34	0.41
1:C:149:ARG:HG3	1:C:158:ILE:HG21	2.03	0.41
1:A:80:PHE:CZ	1:D:96:LEU:HD23	2.56	0.41
1:A:337:ALA:HB1	1:A:338:PRO:CD	2.50	0.41
1:A:295:ARG:NH2	1:A:329:GLU:OE2	2.54	0.41
1:D:195:ARG:HD3	1:D:278:ILE:HG21	2.02	0.41
1:D:276:VAL:HG22	1:D:336:THR:HG21	2.04	0.40
1:C:157:VAL:HG12	1:C:301:LEU:HD12	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	276/303 (91%)	266 (96%)	8 (3%)	2 (1%)	22 22
1	B	289/303 (95%)	283 (98%)	6 (2%)	0	100 100
1	C	283/303 (93%)	271 (96%)	11 (4%)	1 (0%)	34 37
1	D	275/303 (91%)	263 (96%)	12 (4%)	0	100 100
All	All	1123/1212 (93%)	1083 (96%)	37 (3%)	3 (0%)	41 46

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	116	GLY
1	A	241	PRO
1	C	241	PRO

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	205/237 (86%)	196 (96%)	9 (4%)	28 35
1	B	217/237 (92%)	212 (98%)	5 (2%)	50 63
1	C	211/237 (89%)	207 (98%)	4 (2%)	57 71
1	D	199/237 (84%)	189 (95%)	10 (5%)	24 30
All	All	832/948 (88%)	804 (97%)	28 (3%)	37 47

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	ASN
1	A	95	LEU
1	A	125	ARG
1	A	132	ASP
1	A	146	ILE
1	A	162	GLN
1	A	211	GLU
1	A	229	LEU

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Mol	Chain	Res	Type
1	A	238	LEU
1	B	85	MET
1	B	121	VAL
1	B	162	GLN
1	B	211	GLU
1	B	311	GLU
1	C	59	SER
1	C	125	ARG
1	C	162	GLN
1	C	188	ILE
1	D	125	ARG
1	D	134	LEU
1	D	149	ARG
1	D	162	GLN
1	D	199	TYR
1	D	201	GLU
1	D	210	SER
1	D	220	TRP
1	D	229	LEU
1	D	274	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	108	GLN
1	A	162	GLN
1	B	124	GLN
1	B	162	GLN
1	B	299	GLN
1	C	162	GLN
1	D	162	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	ACT	B	1	-	1,3,3	0.61	0	0,3,3	0.00	-
2	ACT	D	1	-	1,3,3	1.80	0	0,3,3	0.00	-
2	ACT	A	350	-	1,3,3	0.61	0	0,3,3	0.00	-
2	ACT	A	1	-	1,3,3	1.24	0	0,3,3	0.00	-
2	ACT	C	1	-	1,3,3	0.53	0	0,3,3	0.00	-

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.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1	ACT	2	0
2	A	350	ACT	3	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^{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	278/303 (91%)	-0.02	3 (1%) 80 79	24, 34, 50, 58	0
1	B	291/303 (96%)	-0.26	1 (0%) 94 93	21, 30, 43, 52	0
1	C	285/303 (94%)	0.14	9 (3%) 47 45	26, 42, 54, 61	0
1	D	277/303 (91%)	0.71	29 (10%) 6 5	35, 49, 61, 66	0
All	All	1131/1212 (93%)	0.14	42 (3%) 41 39	21, 39, 56, 66	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	208	LEU	4.7
1	D	89	GLY	4.6
1	D	140	GLY	4.1
1	D	206	ALA	4.0
1	D	178	SER	3.8
1	D	210	SER	3.8
1	D	209	ARG	3.4
1	D	242	ASP	3.3
1	C	203	LEU	3.3
1	C	317	ALA	3.2
1	D	241	PRO	3.2
1	D	176	GLY	3.1
1	D	240	LYS	3.0
1	D	313	LEU	3.0
1	C	239	GLY	3.0
1	D	272	ASP	3.0
1	C	134	LEU	2.9
1	D	155	GLY	2.9
1	D	243	GLY	2.9
1	D	173	ILE	2.8
1	D	169	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	213	ALA	2.6
1	C	55	ALA	2.6
1	A	117	ARG	2.5
1	C	315	GLY	2.5
1	D	203	LEU	2.4
1	D	324	THR	2.4
1	D	180	ILE	2.4
1	A	112	LEU	2.4
1	C	140	GLY	2.3
1	B	55	ALA	2.3
1	A	103	PRO	2.3
1	D	200	LEU	2.2
1	C	240	LYS	2.2
1	D	175	LEU	2.2
1	C	152	GLY	2.2
1	D	172	LEU	2.2
1	D	304	VAL	2.1
1	D	118	VAL	2.1
1	D	181	ALA	2.0
1	D	338	PRO	2.0
1	D	231	THR	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	ACT	A	350	4/4	0.91	0.14	58,58,59,59	0
2	ACT	C	1	4/4	0.91	0.23	43,43,44,44	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	ACT	B	1	4/4	0.95	0.18	39,39,40,40	0
2	ACT	A	1	4/4	0.97	0.19	39,39,40,40	0
2	ACT	D	1	4/4	0.97	0.16	52,52,52,53	0

6.5 Other polymers (i)

There are no such residues in this entry.