



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 31, 2021 – 04:03 AM EDT

PDB ID : 2E2L
Title : Helicobacter pylori formamidase AmiF contains a fine-tuned cysteine-glutamate-lysine catalytic triad
Authors : Wang, W.C.; Hung, C.L.
Deposited on : 2006-11-14
Resolution : 2.29 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.23.2
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) : 2.23.2

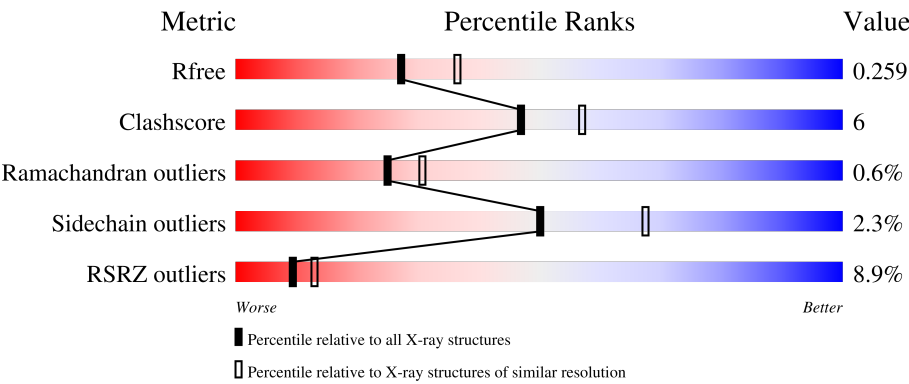
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.29 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (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 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	334	<div><div>7%</div><div>81%13%5%</div></div>
1	B	334	<div><div>9%</div><div>79%15%5%</div></div>
1	C	334	<div><div>7%</div><div>79%14%5%</div></div>
1	D	334	<div><div>7%</div><div>82%12%5%</div></div>
1	E	334	<div><div>10%</div><div>83%10%5%</div></div>

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Mol	Chain	Length	Quality of chain
1	F	334	

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	ARF	A	1335	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 15751 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.

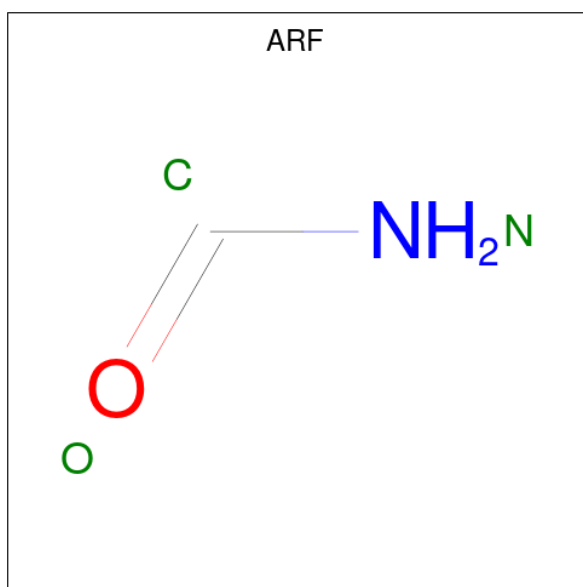
- Molecule 1 is a protein called Formamidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	317	Total	C	N	O	S	0	0	0
			2503	1610	421	461	11			
1	B	317	Total	C	N	O	S	0	0	0
			2503	1610	421	461	11			
1	C	317	Total	C	N	O	S	0	0	0
			2503	1610	421	461	11			
1	D	317	Total	C	N	O	S	0	0	0
			2503	1610	421	461	11			
1	E	316	Total	C	N	O	S	0	0	0
			2492	1604	417	460	11			
1	F	317	Total	C	N	O	S	0	0	0
			2503	1610	421	461	11			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	166	SER	CYS	engineered mutation	UNP O25836
B	166	SER	CYS	engineered mutation	UNP O25836
C	166	SER	CYS	engineered mutation	UNP O25836
D	166	SER	CYS	engineered mutation	UNP O25836
E	166	SER	CYS	engineered mutation	UNP O25836
F	166	SER	CYS	engineered mutation	UNP O25836

- Molecule 2 is FORMAMIDE (three-letter code: ARF) (formula: CH₃NO).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			3	1	1	1		
2	C	1	Total	C	N	O	0	0
			3	1	1	1		
2	F	1	Total	C	N	O	0	0
			3	1	1	1		

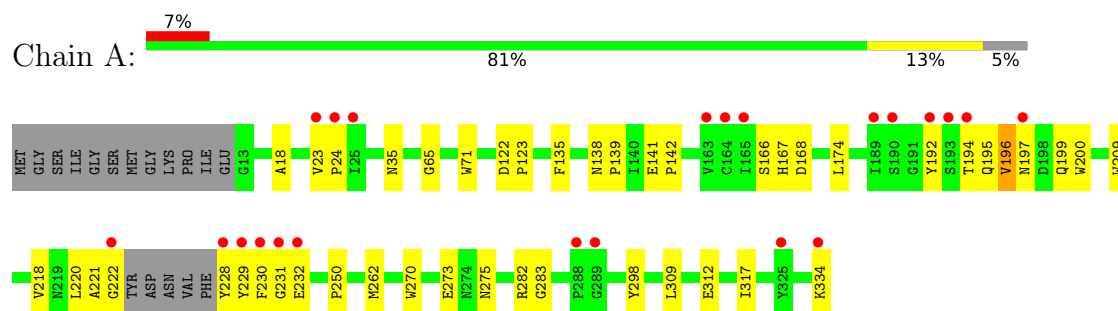
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	133	Total	O	0	0
			133	133		
3	B	131	Total	O	0	0
			131	131		
3	C	143	Total	O	0	0
			143	143		
3	D	102	Total	O	0	0
			102	102		
3	E	114	Total	O	0	0
			114	114		
3	F	112	Total	O	0	0
			112	112		

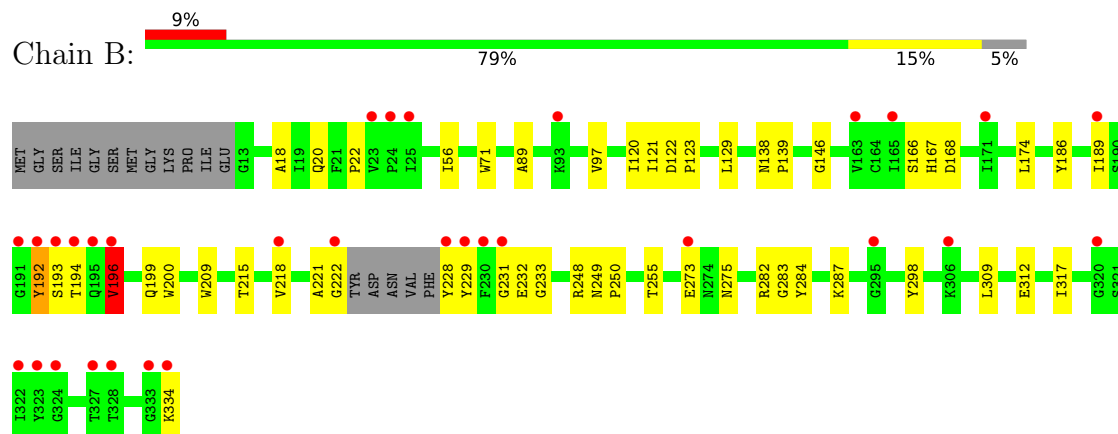
3 Residue-property plots

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.

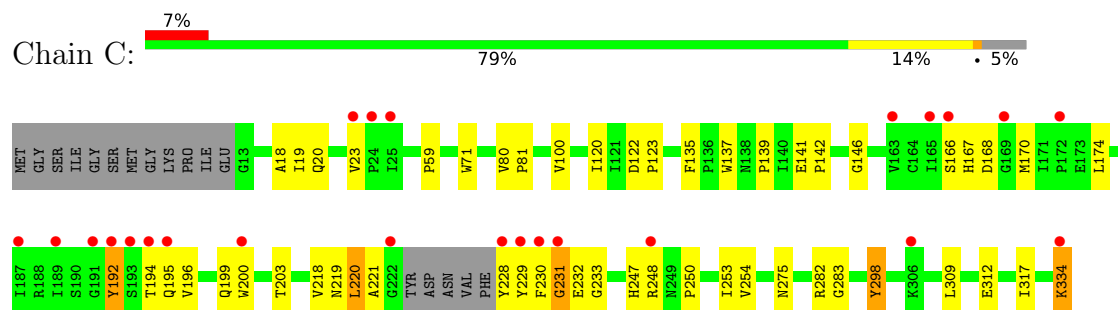
• Molecule 1: Formamidase



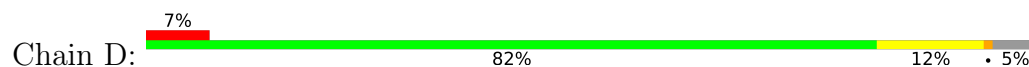
• Molecule 1: Formamidase

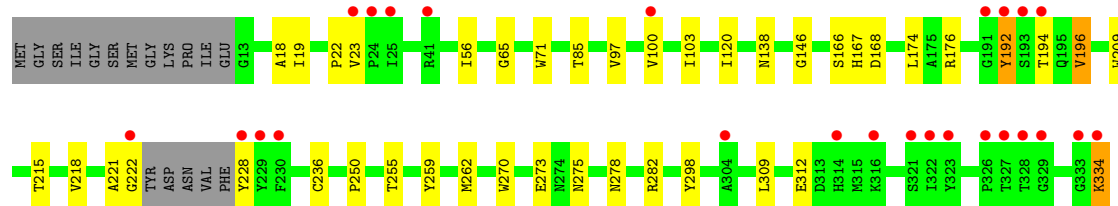


• Molecule 1: Formamidase

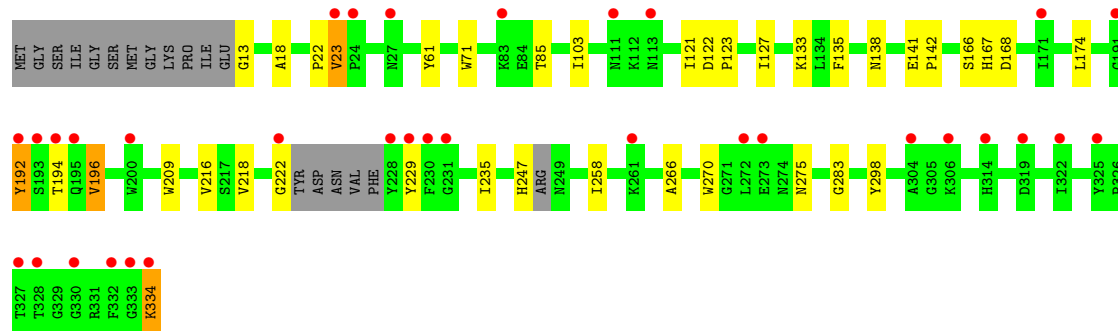
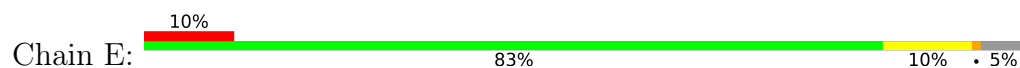


• Molecule 1: Formamidase

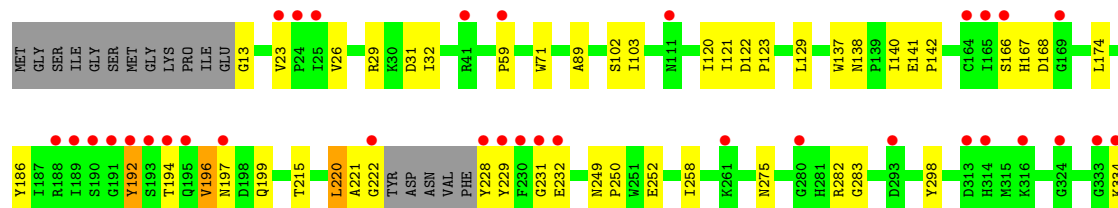
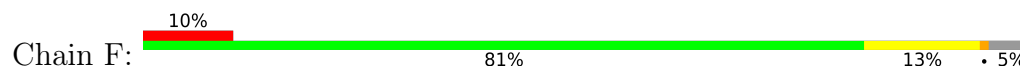




● Molecule 1: Formamidase



● Molecule 1: Formamidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.09Å 151.79Å 89.08Å 90.00° 114.99° 90.00°	Depositor
Resolution (Å)	30.00 – 2.29 29.62 – 2.29	Depositor EDS
% Data completeness (in resolution range)	90.9 (30.00-2.29) 90.9 (29.62-2.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.23 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.253 , 0.285 0.252 , 0.259	Depositor DCC
R_{free} test set	4091 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	35.7	Xtriage
Anisotropy	0.612	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 52.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.016 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	15751	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ARF

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	0.33	0/2574	0.50	0/3498
1	B	0.33	0/2574	0.48	0/3498
1	C	0.33	0/2574	0.50	0/3498
1	D	0.33	0/2574	0.49	0/3498
1	E	0.33	0/2562	0.48	0/3481
1	F	0.33	0/2574	0.48	0/3498
All	All	0.33	0/15432	0.49	0/20971

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	2503	0	2442	43	0
1	B	2503	0	2442	47	0
1	C	2503	0	2442	48	0
1	D	2503	0	2442	41	0
1	E	2492	0	2428	24	0
1	F	2503	0	2442	30	0
2	A	3	0	3	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	3	0	3	1	0
2	F	3	0	3	0	0
3	A	133	0	0	3	0
3	B	131	0	0	1	0
3	C	143	0	0	1	0
3	D	102	0	0	2	0
3	E	114	0	0	1	0
3	F	112	0	0	0	0
All	All	15751	0	14647	185	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 (185) 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:170:MET:CE	1:D:209:TRP:HE1	1.66	1.08
1:C:170:MET:HE2	1:D:209:TRP:HE1	1.24	1.00
1:A:192:TYR:HD2	2:A:1335:ARF:HN2	1.27	0.81
1:C:170:MET:HE2	1:D:209:TRP:NE1	1.96	0.80
1:C:170:MET:CE	1:D:209:TRP:NE1	2.47	0.77
1:C:282:ARG:NH2	1:D:146:GLY:O	2.20	0.74
1:C:219:ASN:HB3	1:C:231:GLY:HA2	1.68	0.73
1:A:275:ASN:HD22	1:B:167:HIS:HE1	1.37	0.71
1:C:334:LYS:HD2	1:C:334:LYS:H	1.58	0.69
1:C:275:ASN:ND2	1:D:167:HIS:HE1	1.91	0.68
1:C:275:ASN:HD22	1:D:167:HIS:HE1	1.38	0.67
1:A:283:GLY:HA2	1:B:138:ASN:HD21	1.58	0.67
1:D:23:VAL:HG21	1:D:65:GLY:HA2	1.77	0.66
1:A:309:LEU:HB2	1:A:312:GLU:HG3	1.79	0.65
1:B:200:TRP:HZ3	1:B:232:GLU:O	1.78	0.65
1:C:170:MET:HE3	1:C:203:THR:HG21	1.78	0.65
1:E:138:ASN:HD21	1:F:283:GLY:HA2	1.61	0.65
1:A:282:ARG:NH2	1:B:146:GLY:O	2.24	0.65
1:F:196:VAL:HG12	1:F:199:GLN:HB2	1.80	0.64
1:B:192:TYR:HD2	1:B:192:TYR:H	1.45	0.62
1:E:85:THR:HG22	1:E:103:ILE:HD13	1.81	0.61
1:E:192:TYR:N	1:E:192:TYR:HD2	1.99	0.60
1:A:230:PHE:HB2	3:A:1461:HOH:O	2.00	0.60
1:A:168:ASP:HB3	1:A:174:LEU:HD23	1.83	0.60
1:B:192:TYR:O	1:B:194:THR:N	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:GLY:HA3	1:A:228:TYR:HB3	1.83	0.59
1:A:229:TYR:CD2	1:D:250:PRO:O	2.56	0.58
1:C:200:TRP:HZ3	1:C:232:GLU:O	1.86	0.58
1:A:275:ASN:ND2	1:B:167:HIS:HE1	1.99	0.58
1:A:192:TYR:HD2	2:A:1335:ARF:N	1.99	0.58
1:F:59:PRO:HB2	1:F:220:LEU:HD13	1.85	0.58
1:A:23:VAL:N	1:A:222:GLY:O	2.36	0.58
1:A:275:ASN:HD22	1:B:167:HIS:CE1	2.20	0.57
1:C:167:HIS:HE1	1:D:275:ASN:ND2	2.03	0.57
1:E:192:TYR:N	1:E:192:TYR:CD2	2.71	0.57
1:E:266:ALA:O	1:E:270:TRP:HB2	2.05	0.56
1:A:135:PHE:CE1	1:B:282:ARG:HG2	2.40	0.56
1:C:146:GLY:O	1:D:282:ARG:NH2	2.33	0.56
1:B:309:LEU:HB2	1:B:312:GLU:HG3	1.87	0.56
1:A:195:GLN:HG3	1:D:255:THR:OG1	2.06	0.56
1:A:250:PRO:O	1:E:229:TYR:CD2	2.59	0.55
1:F:194:THR:C	1:F:196:VAL:H	2.10	0.55
1:B:209:TRP:HH2	1:B:273:GLU:HB3	1.72	0.55
1:C:199:GLN:HG2	1:D:209:TRP:CD2	2.42	0.55
1:D:194:THR:C	1:D:196:VAL:H	2.09	0.54
1:B:18:ALA:HB1	1:B:218:VAL:HB	1.89	0.54
1:C:283:GLY:HA2	1:D:138:ASN:HD21	1.73	0.54
1:E:283:GLY:HA2	1:F:138:ASN:HD21	1.71	0.54
1:B:250:PRO:O	1:C:229:TYR:CD2	2.61	0.54
1:E:247:HIS:C	3:E:447:HOH:O	2.45	0.54
1:C:250:PRO:O	1:F:229:TYR:CD2	2.61	0.54
1:A:275:ASN:ND2	1:B:167:HIS:CE1	2.75	0.53
1:D:192:TYR:N	1:D:192:TYR:CD2	2.75	0.53
1:A:270:TRP:CE3	1:A:273:GLU:HB2	2.44	0.53
1:B:20:GLN:NE2	1:B:233:GLY:H	2.07	0.53
1:A:221:ALA:HA	1:A:231:GLY:CA	2.39	0.52
1:F:231:GLY:O	1:F:232:GLU:HG2	2.10	0.52
1:C:167:HIS:CE1	1:D:275:ASN:ND2	2.77	0.52
1:C:194:THR:C	1:C:196:VAL:H	2.12	0.51
1:D:18:ALA:HB1	1:D:218:VAL:HB	1.93	0.51
1:A:138:ASN:HD21	1:B:283:GLY:HA2	1.75	0.51
1:B:20:GLN:HE21	1:B:233:GLY:H	1.59	0.51
1:A:192:TYR:HA	3:A:1464:HOH:O	2.11	0.51
1:F:222:GLY:HA2	1:F:228:TYR:HA	1.92	0.50
1:F:168:ASP:HB3	1:F:174:LEU:HD23	1.94	0.50
1:C:18:ALA:HB1	1:C:218:VAL:HB	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:275:ASN:ND2	1:D:167:HIS:CE1	2.76	0.50
1:D:221:ALA:HB3	1:D:250:PRO:HA	1.93	0.50
1:D:19:ILE:HD11	1:D:56:ILE:HG23	1.92	0.50
1:D:22:PRO:HA	1:D:222:GLY:H	1.76	0.50
1:A:220:LEU:O	1:A:231:GLY:HA2	2.12	0.49
1:A:24:PRO:HA	1:A:228:TYR:CE1	2.48	0.49
1:E:122:ASP:HB2	1:E:123:PRO:CD	2.43	0.49
1:C:192:TYR:H	1:C:192:TYR:HD2	1.60	0.49
1:D:282:ARG:HG3	3:D:387:HOH:O	2.13	0.49
1:F:221:ALA:HA	1:F:231:GLY:N	2.28	0.49
1:D:22:PRO:HB3	1:D:250:PRO:HB3	1.94	0.48
1:B:196:VAL:HG12	1:B:199:GLN:HB2	1.96	0.48
1:C:309:LEU:HB2	1:C:312:GLU:HG3	1.95	0.48
1:C:135:PHE:CE1	1:D:282:ARG:HG2	2.49	0.48
1:B:222:GLY:HA2	1:B:228:TYR:HA	1.96	0.48
1:C:100:VAL:HG22	1:C:120:ILE:HG12	1.95	0.48
1:D:309:LEU:HB2	1:D:312:GLU:HG3	1.94	0.48
1:A:228:TYR:O	1:A:250:PRO:HG3	2.14	0.48
1:D:270:TRP:CE3	1:D:273:GLU:HB2	2.49	0.48
1:B:228:TYR:O	1:B:250:PRO:HG3	2.14	0.47
1:B:168:ASP:HB3	1:B:174:LEU:HD23	1.97	0.47
1:B:209:TRP:CH2	1:B:273:GLU:HB3	2.49	0.47
1:A:18:ALA:HB1	1:A:218:VAL:HB	1.95	0.47
1:E:135:PHE:CE1	1:F:282:ARG:HG2	2.49	0.47
1:F:102:SER:O	1:F:103:ILE:HG13	2.14	0.47
1:B:249:ASN:HB3	1:C:229:TYR:CG	2.49	0.47
1:C:192:TYR:CD2	1:C:192:TYR:N	2.81	0.47
1:A:122:ASP:HB2	1:A:123:PRO:CD	2.45	0.47
1:C:122:ASP:HB2	1:C:123:PRO:CD	2.45	0.47
1:D:23:VAL:CG2	1:D:65:GLY:HA2	2.42	0.47
1:E:216:VAL:HG13	1:E:235:ILE:HG12	1.96	0.46
1:A:167:HIS:CE1	1:B:275:ASN:HD22	2.33	0.46
1:E:275:ASN:ND2	1:F:167:HIS:CE1	2.84	0.46
1:C:221:ALA:HA	1:C:231:GLY:H	1.80	0.46
1:E:334:LYS:H	1:E:334:LYS:HD2	1.80	0.46
1:B:194:THR:C	1:B:196:VAL:H	2.19	0.46
1:D:215:THR:OG1	1:D:236:CYS:HB2	2.16	0.46
1:F:221:ALA:HB3	1:F:250:PRO:HA	1.98	0.46
1:C:221:ALA:HB3	1:C:250:PRO:HA	1.98	0.46
1:E:18:ALA:HB1	1:E:218:VAL:HB	1.98	0.46
1:E:168:ASP:HB3	1:E:174:LEU:HD23	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:284:TYR:HA	1:B:287:LYS:O	2.16	0.45
1:B:89:ALA:HA	1:B:121:ILE:HG21	1.99	0.45
1:B:282:ARG:HG3	3:B:369:HOH:O	2.16	0.45
1:F:89:ALA:HA	1:F:121:ILE:HG21	1.99	0.45
1:D:209:TRP:HH2	1:D:273:GLU:HB3	1.82	0.45
1:E:192:TYR:HD2	1:E:192:TYR:H	1.64	0.45
1:A:35:ASN:ND2	3:A:1376:HOH:O	2.50	0.44
1:A:270:TRP:HE3	1:A:273:GLU:HB2	1.82	0.44
1:C:221:ALA:O	1:C:228:TYR:C	2.56	0.44
1:F:122:ASP:HB2	1:F:123:PRO:CD	2.47	0.44
1:D:100:VAL:HG22	1:D:120:ILE:HG12	2.00	0.44
1:A:200:TRP:CZ3	1:A:232:GLU:HB3	2.53	0.44
1:B:255:THR:OG1	1:C:195:GLN:HG3	2.18	0.44
1:D:194:THR:C	1:D:196:VAL:N	2.70	0.44
1:F:29:ARG:HA	1:F:32:ILE:HD12	1.98	0.44
1:A:222:GLY:CA	1:A:228:TYR:HB3	2.45	0.44
1:B:186:TYR:HB3	1:B:215:THR:HG22	2.00	0.44
1:B:229:TYR:CG	1:F:249:ASN:HB3	2.52	0.43
1:C:167:HIS:HE1	1:D:275:ASN:HD22	1.64	0.43
1:F:120:ILE:HD12	1:F:129:LEU:HD13	1.99	0.43
1:D:85:THR:HG22	1:D:103:ILE:HD13	2.00	0.43
1:F:13:GLY:HA3	1:F:258:ILE:O	2.19	0.43
1:A:167:HIS:CE1	1:B:275:ASN:ND2	2.86	0.43
1:B:248:ARG:NH1	1:F:252:GLU:OE2	2.51	0.43
1:E:121:ILE:HG12	1:E:127:ILE:HG12	2.01	0.43
1:B:221:ALA:HA	1:B:231:GLY:H	1.84	0.43
1:F:141:GLU:HA	1:F:142:PRO:HD3	1.85	0.43
1:B:229:TYR:CD2	1:F:250:PRO:O	2.72	0.43
1:C:233:GLY:HA3	1:C:254:VAL:HG21	2.00	0.43
1:D:168:ASP:HB3	1:D:174:LEU:HD23	2.00	0.43
1:A:221:ALA:HA	1:A:231:GLY:HA3	1.99	0.43
1:B:122:ASP:HB2	1:B:123:PRO:CD	2.48	0.43
1:B:189:ILE:HG13	1:B:189:ILE:O	2.19	0.42
1:A:209:TRP:CD2	1:B:199:GLN:HG2	2.54	0.42
1:E:209:TRP:CD2	1:F:199:GLN:HG2	2.54	0.42
1:A:139:PRO:HG3	1:A:317:ILE:HG21	2.01	0.42
1:F:137:TRP:O	1:F:140:ILE:HG22	2.19	0.42
1:C:59:PRO:HB2	1:C:220:LEU:HD13	2.02	0.42
1:F:186:TYR:HB3	1:F:215:THR:HG22	2.02	0.42
1:C:141:GLU:HA	1:C:142:PRO:HD3	1.82	0.42
1:E:167:HIS:CE1	1:F:275:ASN:ND2	2.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:192:TYR:O	1:F:194:THR:N	2.44	0.42
1:C:168:ASP:HB3	1:C:174:LEU:HD23	2.02	0.42
1:B:192:TYR:CD2	1:B:192:TYR:N	2.85	0.42
1:D:56:ILE:HD12	1:D:97:VAL:HG21	2.02	0.42
1:A:23:VAL:HG21	1:A:65:GLY:HA2	2.01	0.42
1:C:192:TYR:HA	3:C:2472:HOH:O	2.19	0.41
1:C:139:PRO:HG3	1:C:317:ILE:HG21	2.01	0.41
1:E:13:GLY:HA3	1:E:258:ILE:O	2.20	0.41
1:E:141:GLU:HA	1:E:142:PRO:HD3	1.86	0.41
1:E:194:THR:C	1:E:196:VAL:H	2.22	0.41
1:B:139:PRO:HG3	1:B:317:ILE:HG21	2.02	0.41
1:C:194:THR:HB	1:C:230:PHE:CD1	2.55	0.41
1:D:334:LYS:HD2	1:D:334:LYS:H	1.85	0.41
1:A:194:THR:C	1:A:196:VAL:H	2.24	0.41
1:A:199:GLN:HG2	1:B:209:TRP:CD2	2.55	0.41
1:B:249:ASN:ND2	1:C:248:ARG:HB2	2.36	0.41
1:C:19:ILE:HG13	1:C:253:ILE:HD12	2.02	0.41
1:C:80:VAL:HA	1:C:81:PRO:HA	1.91	0.41
1:D:278:ASN:ND2	3:D:339:HOH:O	2.54	0.41
1:C:20:GLN:NE2	1:C:233:GLY:H	2.18	0.41
1:C:170:MET:CE	1:C:203:THR:HG21	2.50	0.41
1:A:141:GLU:HA	1:A:142:PRO:HD3	1.85	0.41
1:A:283:GLY:CA	1:B:138:ASN:HD21	2.30	0.41
1:B:120:ILE:HD12	1:B:129:LEU:HD13	2.03	0.41
1:C:137:TRP:CE3	2:C:2335:ARF:H	2.56	0.41
1:A:221:ALA:HB3	1:A:250:PRO:HA	2.03	0.41
1:E:23:VAL:HG21	1:E:61:TYR:O	2.21	0.41
1:C:231:GLY:O	1:C:247:HIS:O	2.39	0.40
1:D:259:TYR:HB3	1:D:262:MET:HE3	2.03	0.40
1:E:22:PRO:HA	1:E:222:GLY:HA3	2.03	0.40
1:F:26:VAL:HG13	1:F:31:ASP:HB2	2.02	0.40
1:C:298:TYR:CE1	1:D:176:ARG:HG2	2.56	0.40
1:F:23:VAL:N	1:F:222:GLY:O	2.54	0.40
1:A:138:ASN:HD21	1:B:283:GLY:CA	2.34	0.40
1:A:167:HIS:HE1	1:B:275:ASN:ND2	2.20	0.40
1:D:23:VAL:HG12	1:D:222:GLY:C	2.42	0.40
1:B:56:ILE:HD12	1:B:97:VAL:HG21	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	313/334 (94%)	300 (96%)	12 (4%)	1 (0%)	41	50
1	B	313/334 (94%)	298 (95%)	11 (4%)	4 (1%)	12	12
1	C	313/334 (94%)	298 (95%)	13 (4%)	2 (1%)	25	31
1	D	313/334 (94%)	302 (96%)	10 (3%)	1 (0%)	41	50
1	E	310/334 (93%)	299 (96%)	9 (3%)	2 (1%)	25	31
1	F	313/334 (94%)	299 (96%)	13 (4%)	1 (0%)	41	50
All	All	1875/2004 (94%)	1796 (96%)	68 (4%)	11 (1%)	25	31

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	166	SER
1	B	193	SER
1	C	231	GLY
1	E	166	SER
1	F	166	SER
1	B	166	SER
1	C	166	SER
1	D	166	SER
1	E	133	LYS
1	B	22	PRO
1	B	196	VAL

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	263/277 (95%)	257 (98%)	6 (2%)	50	67
1	B	263/277 (95%)	258 (98%)	5 (2%)	57	73
1	C	263/277 (95%)	257 (98%)	6 (2%)	50	67
1	D	263/277 (95%)	257 (98%)	6 (2%)	50	67
1	E	262/277 (95%)	256 (98%)	6 (2%)	50	67
1	F	263/277 (95%)	256 (97%)	7 (3%)	44	61
All	All	1577/1662 (95%)	1541 (98%)	36 (2%)	50	67

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

Mol	Chain	Res	Type
1	A	71	TRP
1	A	196	VAL
1	A	197	ASN
1	A	262	MET
1	A	298	TYR
1	A	334	LYS
1	B	71	TRP
1	B	192	TYR
1	B	196	VAL
1	B	298	TYR
1	B	334	LYS
1	C	23	VAL
1	C	71	TRP
1	C	192	TYR
1	C	220	LEU
1	C	298	TYR
1	C	334	LYS
1	D	71	TRP
1	D	192	TYR
1	D	196	VAL
1	D	228	TYR
1	D	298	TYR
1	D	334	LYS
1	E	23	VAL
1	E	71	TRP
1	E	192	TYR
1	E	196	VAL
1	E	298	TYR
1	E	334	LYS
1	F	71	TRP

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Mol	Chain	Res	Type
1	F	192	TYR
1	F	196	VAL
1	F	197	ASN
1	F	220	LEU
1	F	298	TYR
1	F	334	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	138	ASN
1	A	167	HIS
1	A	210	HIS
1	A	275	ASN
1	A	278	ASN
1	A	281	HIS
1	B	20	GLN
1	B	138	ASN
1	B	167	HIS
1	B	197	ASN
1	B	210	HIS
1	B	281	HIS
1	B	292	HIS
1	C	20	GLN
1	C	138	ASN
1	C	167	HIS
1	C	210	HIS
1	C	249	ASN
1	C	275	ASN
1	C	278	ASN
1	D	138	ASN
1	D	167	HIS
1	D	197	ASN
1	D	210	HIS
1	D	249	ASN
1	D	275	ASN
1	D	278	ASN
1	D	281	HIS
1	D	292	HIS
1	E	138	ASN
1	E	167	HIS
1	E	197	ASN

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Mol	Chain	Res	Type
1	E	210	HIS
1	E	275	ASN
1	F	20	GLN
1	F	138	ASN
1	F	167	HIS
1	F	210	HIS
1	F	275	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 ⓘ

3 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	ARF	A	1335	-	2,2,2	2.31	1 (50%)	1,1,1	0.77	0
2	ARF	F	3335	-	2,2,2	2.38	1 (50%)	1,1,1	0.74	0
2	ARF	C	2335	-	2,2,2	2.31	1 (50%)	1,1,1	0.78	0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	3335	ARF	O-C	3.33	1.33	1.22
2	C	2335	ARF	O-C	3.23	1.33	1.22
2	A	1335	ARF	O-C	3.22	1.33	1.22

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 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1335	ARF	2	0
2	C	2335	ARF	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, 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	317/334 (94%)	0.44	22 (6%) 16 22	24, 34, 42, 49	0
1	B	317/334 (94%)	0.49	31 (9%) 7 10	24, 35, 49, 52	0
1	C	317/334 (94%)	0.40	24 (7%) 13 18	25, 34, 44, 47	0
1	D	317/334 (94%)	0.45	25 (7%) 12 17	27, 36, 51, 56	0
1	E	316/334 (94%)	0.60	33 (10%) 6 9	28, 39, 55, 59	0
1	F	317/334 (94%)	0.67	34 (10%) 6 8	28, 39, 48, 51	0
All	All	1901/2004 (94%)	0.51	169 (8%) 9 13	24, 36, 48, 59	0

All (169) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	193	SER	12.3
1	A	229	TYR	11.6
1	F	228	TYR	11.4
1	F	229	TYR	10.7
1	B	192	TYR	10.5
1	A	228	TYR	10.4
1	B	228	TYR	9.8
1	B	229	TYR	8.3
1	D	229	TYR	8.3
1	C	229	TYR	8.3
1	E	192	TYR	8.0
1	E	229	TYR	7.7
1	C	228	TYR	7.5
1	F	192	TYR	7.5
1	B	193	SER	7.5
1	A	192	TYR	7.3
1	C	193	SER	7.2
1	C	222	GLY	7.0
1	E	228	TYR	7.0

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Mol	Chain	Res	Type	RSRZ
1	D	228	TYR	6.5
1	C	192	TYR	6.5
1	F	24	PRO	6.4
1	F	193	SER	6.4
1	B	24	PRO	6.0
1	E	327	THR	5.9
1	D	192	TYR	5.9
1	C	25	ILE	5.6
1	A	193	SER	5.6
1	D	24	PRO	5.6
1	C	194	THR	5.4
1	A	24	PRO	5.2
1	F	191	GLY	5.2
1	E	191	GLY	5.0
1	C	230	PHE	5.0
1	F	194	THR	4.9
1	C	24	PRO	4.7
1	A	25	ILE	4.5
1	A	194	THR	4.5
1	B	191	GLY	4.4
1	A	230	PHE	4.4
1	D	25	ILE	4.3
1	D	193	SER	4.3
1	E	222	GLY	4.2
1	E	23	VAL	4.1
1	F	189	ILE	4.1
1	B	222	GLY	4.1
1	F	324	GLY	4.1
1	E	230	PHE	4.0
1	D	334	LYS	4.0
1	E	24	PRO	4.0
1	A	334	LYS	4.0
1	D	328	THR	3.9
1	B	25	ILE	3.9
1	E	328	THR	3.9
1	D	194	THR	3.9
1	A	231	GLY	3.8
1	D	222	GLY	3.8
1	D	333	GLY	3.7
1	D	230	PHE	3.7
1	B	23	VAL	3.6
1	C	23	VAL	3.6

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Mol	Chain	Res	Type	RSRZ
1	E	314	HIS	3.6
1	A	222	GLY	3.4
1	F	41	ARG	3.4
1	D	323	TYR	3.4
1	B	333	GLY	3.4
1	F	23	VAL	3.4
1	D	316	LYS	3.4
1	C	191	GLY	3.3
1	B	194	THR	3.3
1	F	230	PHE	3.3
1	F	261	LYS	3.3
1	B	327	THR	3.3
1	C	200	TRP	3.3
1	D	321	SER	3.2
1	D	191	GLY	3.2
1	D	314	HIS	3.2
1	B	231	GLY	3.2
1	C	189	ILE	3.1
1	D	327	THR	3.1
1	A	189	ILE	3.1
1	E	333	GLY	3.1
1	B	230	PHE	3.1
1	E	334	LYS	3.0
1	A	163	VAL	3.0
1	E	325	TYR	3.0
1	F	334	LYS	3.0
1	B	328	THR	3.0
1	C	187	ILE	3.0
1	D	329	GLY	2.9
1	F	333	GLY	2.9
1	F	190	SER	2.8
1	F	231	GLY	2.8
1	A	325	TYR	2.7
1	F	164	CYS	2.7
1	A	190	SER	2.7
1	C	195	GLN	2.7
1	B	324	GLY	2.7
1	B	195	GLN	2.6
1	E	194	THR	2.6
1	B	306	LYS	2.6
1	F	25	ILE	2.6
1	F	188	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	165	ILE	2.5
1	A	23	VAL	2.5
1	E	111	ASN	2.5
1	E	306	LYS	2.5
1	C	165	ILE	2.5
1	E	319	ASP	2.5
1	B	334	LYS	2.4
1	D	322	ILE	2.4
1	E	261	LYS	2.4
1	C	163	VAL	2.4
1	C	172	PRO	2.4
1	F	195	GLN	2.4
1	B	171	ILE	2.4
1	E	322	ILE	2.4
1	F	165	ILE	2.4
1	A	197	ASN	2.4
1	B	322	ILE	2.4
1	D	326	PRO	2.4
1	B	320	GLY	2.3
1	D	304	ALA	2.3
1	B	93	LYS	2.3
1	F	169	GLY	2.3
1	F	197	ASN	2.3
1	F	293	ASP	2.3
1	F	280	GLY	2.3
1	C	306	LYS	2.3
1	E	304	ALA	2.3
1	C	334	LYS	2.3
1	F	316	LYS	2.3
1	E	195	GLN	2.3
1	F	314	HIS	2.3
1	E	273	GLU	2.3
1	B	189	ILE	2.3
1	E	272	LEU	2.2
1	E	231	GLY	2.2
1	B	163	VAL	2.2
1	E	200	TRP	2.2
1	A	165	ILE	2.2
1	F	232	GLU	2.2
1	B	323	TYR	2.2
1	C	169	GLY	2.2
1	F	166	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	231	GLY	2.2
1	E	171	ILE	2.2
1	B	295	GLY	2.2
1	E	330	GLY	2.2
1	F	313	ASP	2.2
1	E	332	PHE	2.2
1	A	232	GLU	2.2
1	E	27	ASN	2.1
1	D	23	VAL	2.1
1	E	113	ASN	2.1
1	A	164	CYS	2.1
1	B	218	VAL	2.1
1	D	41	ARG	2.1
1	D	100	VAL	2.1
1	A	288	PRO	2.1
1	F	222	GLY	2.1
1	B	273	GLU	2.1
1	F	111	ASN	2.1
1	B	196	VAL	2.1
1	C	166	SER	2.1
1	F	59	PRO	2.0
1	C	248	ARG	2.0
1	E	83	LYS	2.0
1	A	289	GLY	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, 95th 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(\AA^2)	Q<0.9
2	ARF	A	1335	3/3	0.83	0.26	36,36,40,40	0
2	ARF	C	2335	3/3	0.88	0.27	41,41,44,44	0
2	ARF	F	3335	3/3	0.90	0.28	40,40,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.