



# Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 10:21 am BST

PDB ID : 3U30  
Title : Crystal structure of a linear-specific Ubiquitin fab bound to linear ubiquitin  
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Deposited on : 2011-10-04  
Resolution : 2.43 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](mailto: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.

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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	:	2.11
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)	:	2.11

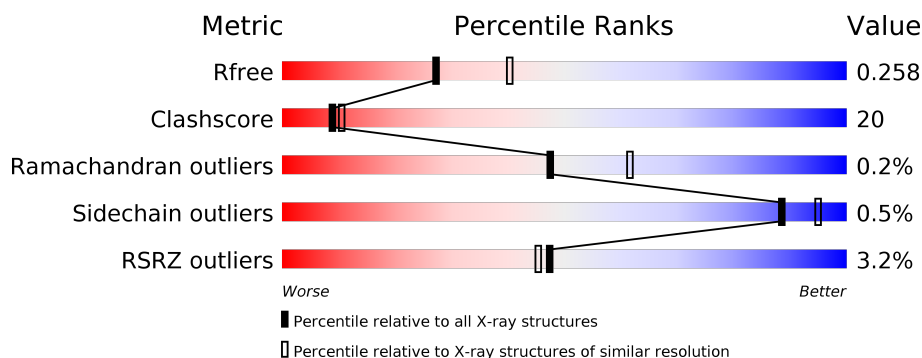
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.43 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	4647 (2.44-2.40)
Clashscore	141614	5161 (2.44-2.40)
Ramachandran outliers	138981	5073 (2.44-2.40)
Sidechain outliers	138945	5074 (2.44-2.40)
RSRZ outliers	127900	4543 (2.44-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	172	<div> <div>%</div> <div> <div></div> <div>68%</div> <div>18%</div> <div>•</div> <div>13%</div> </div> </div>
1	D	172	<div> <div>%</div> <div> <div></div> <div>51%</div> <div>33%</div> <div>16%</div> </div> </div>
2	B	214	<div> <div>5%</div> <div> <div></div> <div>72%</div> <div>26%</div> <div>••</div> </div> </div>
2	E	214	<div> <div>5%</div> <div> <div></div> <div>63%</div> <div>35%</div> <div>••</div> </div> </div>
3	C	227	<div> <div>5%</div> <div> <div></div> <div>65%</div> <div>30%</div> <div>••</div> </div> </div>
3	F	227	<div> <div>%</div> <div> <div></div> <div>70%</div> <div>25%</div> <div>•</div> </div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8998 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 linear di-ubiquitin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	149	Total	C	N	O	S	0	0	0
			1180	743	203	232	2			
1	D	145	Total	C	N	O	S	0	0	0
			1148	723	197	227	1			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP P0CG48
A	-18	GLY	-	EXPRESSION TAG	UNP P0CG48
A	-17	SER	-	EXPRESSION TAG	UNP P0CG48
A	-16	SER	-	EXPRESSION TAG	UNP P0CG48
A	-15	HIS	-	EXPRESSION TAG	UNP P0CG48
A	-14	HIS	-	EXPRESSION TAG	UNP P0CG48
A	-13	HIS	-	EXPRESSION TAG	UNP P0CG48
A	-12	HIS	-	EXPRESSION TAG	UNP P0CG48
A	-11	HIS	-	EXPRESSION TAG	UNP P0CG48
A	-10	HIS	-	EXPRESSION TAG	UNP P0CG48
A	-9	SER	-	EXPRESSION TAG	UNP P0CG48
A	-8	SER	-	EXPRESSION TAG	UNP P0CG48
A	-7	GLY	-	EXPRESSION TAG	UNP P0CG48
A	-6	LEU	-	EXPRESSION TAG	UNP P0CG48
A	-5	VAL	-	EXPRESSION TAG	UNP P0CG48
A	-4	PRO	-	EXPRESSION TAG	UNP P0CG48
A	-3	ARG	-	EXPRESSION TAG	UNP P0CG48
A	-2	GLY	-	EXPRESSION TAG	UNP P0CG48
A	-1	SER	-	EXPRESSION TAG	UNP P0CG48
A	0	HIS	-	EXPRESSION TAG	UNP P0CG48
D	-19	MET	-	EXPRESSION TAG	UNP P0CG48
D	-18	GLY	-	EXPRESSION TAG	UNP P0CG48
D	-17	SER	-	EXPRESSION TAG	UNP P0CG48
D	-16	SER	-	EXPRESSION TAG	UNP P0CG48
D	-15	HIS	-	EXPRESSION TAG	UNP P0CG48

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-14	HIS	-	EXPRESSION TAG	UNP P0CG48
D	-13	HIS	-	EXPRESSION TAG	UNP P0CG48
D	-12	HIS	-	EXPRESSION TAG	UNP P0CG48
D	-11	HIS	-	EXPRESSION TAG	UNP P0CG48
D	-10	HIS	-	EXPRESSION TAG	UNP P0CG48
D	-9	SER	-	EXPRESSION TAG	UNP P0CG48
D	-8	SER	-	EXPRESSION TAG	UNP P0CG48
D	-7	GLY	-	EXPRESSION TAG	UNP P0CG48
D	-6	LEU	-	EXPRESSION TAG	UNP P0CG48
D	-5	VAL	-	EXPRESSION TAG	UNP P0CG48
D	-4	PRO	-	EXPRESSION TAG	UNP P0CG48
D	-3	ARG	-	EXPRESSION TAG	UNP P0CG48
D	-2	GLY	-	EXPRESSION TAG	UNP P0CG48
D	-1	SER	-	EXPRESSION TAG	UNP P0CG48
D	0	HIS	-	EXPRESSION TAG	UNP P0CG48

- Molecule 2 is a protein called Light chain Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	211	Total	C	N	O	S	0	0	0
			1620	1017	269	329	5			
2	E	211	Total	C	N	O	S	0	0	0
			1620	1017	269	329	5			

- Molecule 3 is a protein called Heavy chain Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	218	Total	C	N	O	S	0	0	0
			1618	1022	271	319	6			
3	F	218	Total	C	N	O	S	0	0	0
			1618	1022	271	319	6			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	22	Total	O	0	0
			22	22		
4	B	30	Total	O	0	0
			30	30		
4	C	45	Total	O	0	0
			45	45		

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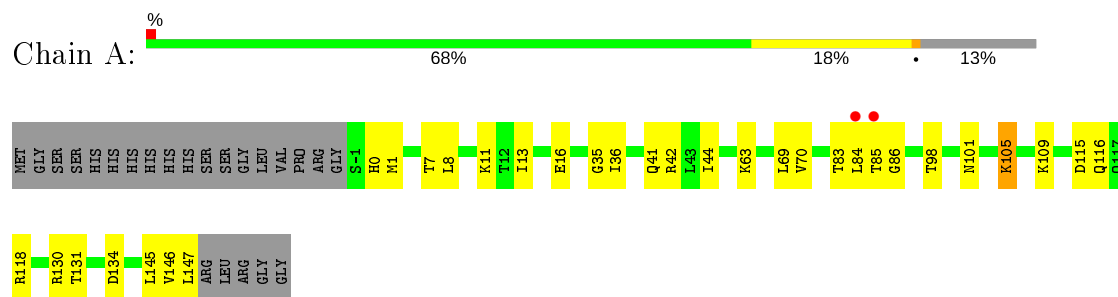
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	22	Total 22	O 22	0	0
4	E	42	Total 42	O 42	0	0
4	F	33	Total 33	O 33	0	0

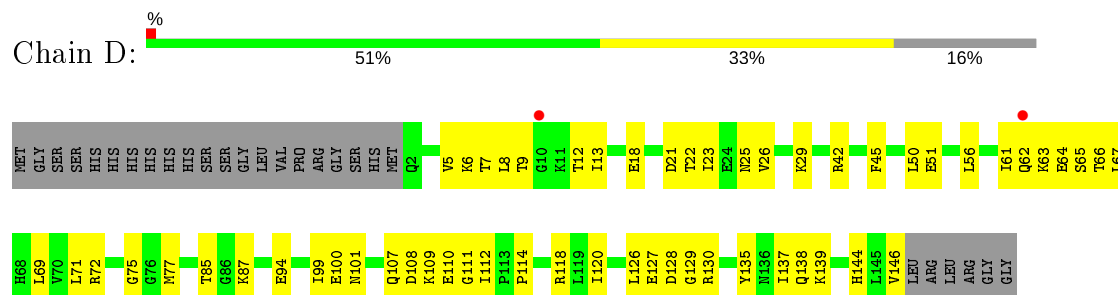
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

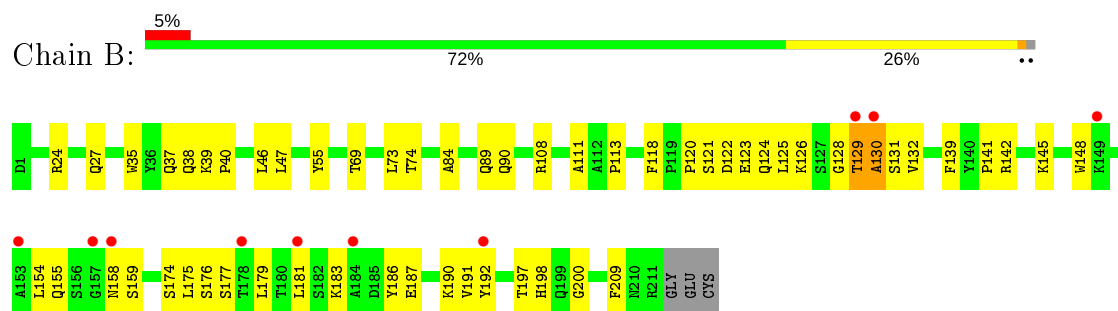
#### • Molecule 1: linear di-ubiquitin



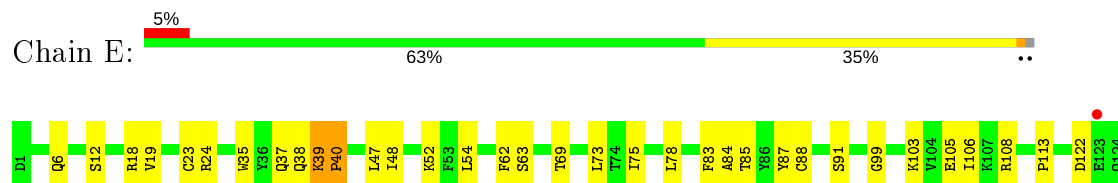
#### • Molecule 1: linear di-ubiquitin

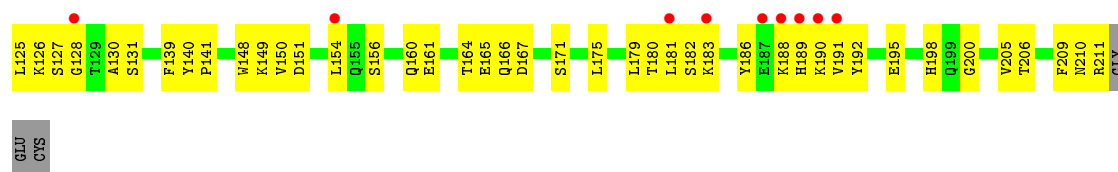


#### • Molecule 2: Light chain Fab



#### • Molecule 2: Light chain Fab





• Molecule 3: Heavy chain Fab



• Molecule 3: Heavy chain Fab



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	53.32Å 59.77Å 96.03Å 87.08° 76.77° 71.70°	Depositor
Resolution (Å)	49.33 – 2.43 49.33 – 2.43	Depositor EDS
% Data completeness (in resolution range)	89.6 (49.33-2.43) 94.9 (49.33-2.43)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.09 (at 2.42Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, $R_{free}$	0.229 , 0.258 0.228 , 0.258	Depositor DCC
$R_{free}$ test set	1956 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.9	Xtriage
Anisotropy	0.350	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 46.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	8998	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.87% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

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	0.22	0/1194	0.40	0/1609
1	D	0.23	0/1161	0.42	0/1565
2	B	0.23	0/1656	0.57	2/2251 (0.1%)
2	E	0.34	1/1656 (0.1%)	0.44	0/2251
3	C	0.23	0/1657	0.51	1/2264 (0.0%)
3	F	0.24	0/1657	0.45	0/2264
All	All	0.25	1/8981 (0.0%)	0.48	3/12204 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	40	PRO	N-CD	-8.96	1.35	1.47

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	130	ALA	N-CA-CB	-12.96	91.96	110.10
3	C	148	GLU	C-N-CD	-12.62	92.84	120.60
2	B	130	ALA	N-CA-C	12.47	144.66	111.00

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	1180	0	1225	27	0
1	D	1148	0	1190	72	0
2	B	1620	0	1581	63	0
2	E	1620	0	1581	91	0
3	C	1618	0	1592	66	0
3	F	1618	0	1592	55	0
4	A	22	0	0	6	0
4	B	30	0	0	2	0
4	C	45	0	0	19	0
4	D	22	0	0	22	0
4	E	42	0	0	8	0
4	F	33	0	0	9	0
All	All	8998	0	8761	357	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 20.

All (357) 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)
2:B:129:THR:CA	2:B:130:ALA:HB3	1.45	1.43
3:C:148:GLU:HG3	3:C:149:PRO:CB	1.54	1.37
3:C:148:GLU:CG	3:C:149:PRO:HB3	1.60	1.28
2:B:129:THR:HA	2:B:130:ALA:CB	1.62	1.28
3:F:110:THR:HA	4:F:246:HOH:O	1.31	1.22
1:D:130:ARG:NH1	4:D:166:HOH:O	1.81	1.11
2:E:52:LYS:NZ	4:E:237:HOH:O	1.81	1.11
2:B:125:LEU:HA	2:B:129:THR:O	1.50	1.10
3:F:111:VAL:N	4:F:246:HOH:O	1.83	1.08
1:D:130:ARG:NE	4:D:181:HOH:O	1.90	1.04
1:D:101:ASN:ND2	4:D:153:HOH:O	1.92	1.02
3:C:177:SER:HB2	4:C:264:HOH:O	1.58	1.00
1:D:87:LYS:NZ	1:D:110:GLU:OE2	1.96	0.97
1:D:101:ASN:HB2	4:D:153:HOH:O	1.65	0.95
1:D:61:ILE:O	1:D:62:GLN:HG2	1.66	0.94
3:F:126:PRO:HG3	3:F:138:LEU:HB3	1.49	0.93
1:D:101:ASN:CB	4:D:153:HOH:O	2.14	0.93
2:B:27:GLN:OE1	4:B:232:HOH:O	1.84	0.93
2:E:195:GLU:O	4:E:245:HOH:O	1.88	0.92
2:E:151:ASP:OD1	2:E:191:VAL:HG23	1.70	0.91
2:E:188:LYS:O	2:E:189:HIS:CD2	2.25	0.90
1:D:101:ASN:CG	4:D:153:HOH:O	2.09	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:42:ARG:HB2	1:D:72:ARG:HG3	1.55	0.88
2:B:130:ALA:HB1	2:B:181:LEU:O	1.73	0.88
2:E:128:GLY:HA2	2:E:183:LYS:HD3	1.52	0.88
1:D:62:GLN:CG	1:D:63:LYS:H	1.85	0.88
1:D:23:ILE:HD12	1:D:50:LEU:HD21	1.54	0.88
3:C:2:VAL:N	4:C:263:HOH:O	2.06	0.86
1:A:85:THR:OG1	1:A:86:GLY:HA2	1.75	0.86
3:F:82:MET:HB3	3:F:82(C):LEU:HD21	1.57	0.86
3:C:30:SER:O	4:C:232:HOH:O	1.93	0.86
1:D:130:ARG:CZ	4:D:181:HOH:O	2.21	0.85
2:E:190:LYS:HB3	2:E:211:ARG:H	1.40	0.85
2:E:209:PHE:CD1	2:E:209:PHE:O	2.30	0.83
3:F:42:GLY:N	4:F:247:HOH:O	2.03	0.83
2:E:150:VAL:HG22	2:E:192:TYR:CD1	2.13	0.83
3:C:72:ASP:OD1	4:C:265:HOH:O	1.95	0.83
1:A:41:GLN:O	4:A:153:HOH:O	1.98	0.82
2:E:38:GLN:OE1	4:E:239:HOH:O	1.97	0.82
1:D:128:ASP:O	4:D:169:HOH:O	1.97	0.81
3:C:171:GLN:NE2	4:C:264:HOH:O	2.05	0.81
2:E:210:ASN:O	2:E:211:ARG:HB2	1.79	0.80
3:F:87:THR:OG1	4:F:246:HOH:O	1.99	0.80
2:B:155:GLN:HB3	2:B:158:ASN:HD21	1.46	0.80
1:A:101:ASN:OD1	4:A:164:HOH:O	1.99	0.79
2:E:151:ASP:OD1	2:E:191:VAL:CG2	2.30	0.79
2:B:124:GLN:HG2	2:B:129:THR:OG1	1.81	0.79
2:E:150:VAL:HG22	2:E:192:TYR:HD1	1.47	0.79
3:F:42:GLY:CA	4:F:247:HOH:O	2.31	0.78
2:E:78:LEU:HD21	2:E:106:ILE:HD13	1.63	0.78
3:F:42:GLY:HA2	4:F:247:HOH:O	1.84	0.77
1:A:35:GLY:O	4:A:168:HOH:O	2.02	0.77
2:B:108:ARG:HH12	2:B:111:ALA:HB2	1.49	0.77
2:E:210:ASN:O	2:E:211:ARG:CB	2.33	0.77
3:C:47:TRP:O	4:C:243:HOH:O	2.04	0.76
3:C:148:GLU:HG3	3:C:149:PRO:CA	2.17	0.75
1:D:107:GLN:C	4:D:160:HOH:O	2.24	0.75
3:C:156:SER:H	3:C:197:ASN:HD21	1.34	0.75
3:F:156:SER:H	3:F:197:ASN:HD21	1.36	0.74
2:E:186:TYR:HA	2:E:192:TYR:OH	1.88	0.73
2:B:129:THR:CA	2:B:130:ALA:CB	2.31	0.73
1:D:77:MET:HB3	3:F:56:GLN:NE2	2.03	0.72
2:E:160:GLN:HB3	3:F:169:VAL:HG11	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:130:ARG:NH2	4:D:181:HOH:O	2.20	0.72
3:C:119:PRO:HB3	3:C:145:TYR:HB3	1.72	0.72
1:A:101:ASN:ND2	4:A:164:HOH:O	2.17	0.72
2:E:18:ARG:NE	4:E:250:HOH:O	2.23	0.70
2:E:190:LYS:HB3	2:E:211:ARG:N	2.07	0.70
3:C:185:PRO:HG2	3:C:188:SER:HB3	1.74	0.70
2:E:131:SER:O	4:E:240:HOH:O	2.10	0.70
3:C:148:GLU:HG3	3:C:149:PRO:HB3	0.74	0.69
2:B:198:HIS:CD2	2:B:200:GLY:H	2.11	0.69
2:B:37:GLN:HB2	2:B:47:LEU:HD11	1.73	0.69
1:D:111:GLY:N	4:D:160:HOH:O	2.24	0.68
1:D:6:LYS:HG2	1:D:12:THR:HG22	1.76	0.68
1:D:75:GLY:O	1:D:94:GLU:HG2	1.93	0.67
1:D:62:GLN:CG	1:D:63:LYS:N	2.56	0.67
3:C:94:ARG:NH1	3:C:101:ASP:OD1	2.27	0.67
3:F:110:THR:CA	4:F:246:HOH:O	2.00	0.66
3:C:193:THR:HG23	3:C:210:LYS:HE3	1.77	0.66
2:B:124:GLN:HE22	2:B:131:SER:CB	2.07	0.66
2:E:19:VAL:HG22	2:E:75:ILE:HB	1.78	0.66
2:B:158:ASN:ND2	2:B:179:LEU:HD11	2.12	0.65
3:C:213:PRO:O	4:C:252:HOH:O	2.13	0.65
2:B:145:LYS:HB3	2:B:197:THR:OG1	1.97	0.65
3:C:29:PHE:O	3:C:32:THR:HG22	1.97	0.65
1:D:22:THR:N	4:D:161:HOH:O	1.95	0.65
3:F:11:LEU:HB2	3:F:147:PRO:HG3	1.79	0.65
1:D:77:MET:HB3	3:F:56:GLN:HE21	1.60	0.65
1:A:101:ASN:CG	4:A:164:HOH:O	2.36	0.64
3:C:200:HIS:HD2	3:C:203:SER:OG	1.81	0.63
3:C:123:PRO:HD3	3:C:209:LYS:HE2	1.81	0.63
1:D:100:GLU:HB3	4:D:169:HOH:O	1.99	0.63
1:D:118:ARG:HH11	1:D:146:VAL:HG11	1.62	0.63
1:D:23:ILE:HG23	1:D:50:LEU:HD22	1.80	0.63
2:B:124:GLN:HE22	2:B:131:SER:HB2	1.64	0.62
1:D:108:ASP:C	4:D:160:HOH:O	2.38	0.62
1:D:9:THR:HG22	1:D:9:THR:O	1.98	0.62
3:C:99:ARG:NH2	4:C:257:HOH:O	2.32	0.62
3:C:144:ASP:OD1	4:C:261:HOH:O	2.16	0.62
2:E:37:GLN:HB2	2:E:47:LEU:HD11	1.82	0.62
3:F:147:PRO:O	3:F:200:HIS:HE1	1.82	0.62
2:E:151:ASP:HA	2:E:191:VAL:HB	1.80	0.61
2:E:198:HIS:CD2	2:E:200:GLY:H	2.19	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:6:GLN:HE21	2:E:99:GLY:HA3	1.66	0.61
3:C:156:SER:H	3:C:197:ASN:ND2	1.99	0.61
1:A:83:THR:O	1:A:86:GLY:HA3	2.00	0.61
2:E:188:LYS:O	2:E:189:HIS:CG	2.54	0.60
1:D:62:GLN:HG3	1:D:63:LYS:H	1.65	0.60
2:B:125:LEU:HD23	2:B:129:THR:O	2.00	0.60
1:D:62:GLN:HG2	1:D:63:LYS:H	1.65	0.60
3:F:168:ALA:HA	3:F:178:LEU:HB3	1.81	0.60
2:B:38:GLN:O	2:B:84:ALA:HB1	2.01	0.60
1:D:61:ILE:O	1:D:62:GLN:CG	2.46	0.60
3:F:31:ASN:ND2	4:F:239:HOH:O	2.34	0.60
2:B:120:PRO:HG2	2:B:186:TYR:CE1	2.36	0.59
2:E:190:LYS:HD3	2:E:211:ARG:HA	1.84	0.59
1:D:107:GLN:OE1	1:D:114:PRO:HD3	2.03	0.59
3:F:119:PRO:HB3	3:F:145:TYR:HB3	1.84	0.59
1:D:23:ILE:CD1	1:D:50:LEU:HD21	2.28	0.58
2:E:6:GLN:HE22	2:E:87:TYR:HA	1.68	0.58
2:E:150:VAL:HA	2:E:191:VAL:O	2.03	0.58
3:F:63:VAL:HG13	3:F:67:PHE:HB2	1.86	0.58
1:D:100:GLU:CB	4:D:169:HOH:O	2.52	0.58
2:E:182:SER:O	2:E:183:LYS:HG2	2.04	0.58
3:F:156:SER:H	3:F:197:ASN:ND2	2.01	0.58
2:E:151:ASP:OD2	2:E:189:HIS:HB3	2.04	0.57
2:E:83:PHE:CE2	2:E:165:GLU:HB2	2.39	0.57
1:D:23:ILE:HG13	1:D:51:GLU:O	2.05	0.57
2:E:210:ASN:O	2:E:211:ARG:HG3	2.05	0.57
1:D:62:GLN:HG2	1:D:63:LYS:N	2.20	0.57
2:E:148:TRP:HB2	4:E:231:HOH:O	2.04	0.57
2:E:151:ASP:H	2:E:191:VAL:HB	1.70	0.56
2:E:39:LYS:HB3	2:E:40:PRO:HD2	1.88	0.56
2:B:130:ALA:HB1	2:B:181:LEU:C	2.27	0.56
2:B:129:THR:HA	2:B:130:ALA:HB3	0.66	0.56
2:B:142:ARG:NH1	4:B:236:HOH:O	2.30	0.55
3:F:126:PRO:HD2	3:F:213:PRO:HA	1.88	0.55
2:E:180:THR:HG21	3:F:143:LYS:NZ	2.21	0.55
3:C:199:ASN:HD21	3:C:201:LYS:NZ	2.04	0.55
3:F:151:THR:OG1	3:F:199:ASN:HB3	2.06	0.55
3:F:200:HIS:HD2	3:F:203:SER:OG	1.89	0.55
3:F:18:LEU:HB3	3:F:82:MET:HE3	1.89	0.55
1:A:0:HIS:HE1	1:A:16:GLU:OE2	1.90	0.55
3:C:151:THR:OG1	3:C:199:ASN:HB3	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:108:ARG:NH1	2:B:111:ALA:HB2	2.20	0.54
3:C:144:ASP:N	4:C:264:HOH:O	1.93	0.54
3:C:124:LEU:HD12	3:C:139:GLY:HA3	1.90	0.54
3:C:93:ALA:HB1	3:C:100(E):MET:HB3	1.90	0.54
2:B:128:GLY:O	2:B:183:LYS:N	2.35	0.54
2:E:83:PHE:HE2	2:E:165:GLU:HB2	1.72	0.53
1:D:107:GLN:O	1:D:111:GLY:HA2	2.09	0.53
1:D:7:THR:HG22	1:D:69:LEU:HD23	1.89	0.53
2:E:151:ASP:CA	2:E:191:VAL:HB	2.37	0.53
2:B:124:GLN:HG3	3:C:122:PHE:CE2	2.43	0.53
2:E:161:GLU:OE1	2:E:175:LEU:HD11	2.09	0.53
1:D:127:GLU:CD	1:D:129:GLY:H	2.12	0.53
1:D:107:GLN:O	4:D:160:HOH:O	2.18	0.53
1:D:9:THR:CG2	1:D:9:THR:O	2.57	0.53
2:E:78:LEU:CD2	2:E:106:ILE:HD13	2.36	0.53
3:C:168:ALA:HB2	3:C:178:LEU:HD23	1.91	0.52
2:E:122:ASP:O	2:E:126:LYS:HG2	2.08	0.52
2:E:210:ASN:O	2:E:211:ARG:CG	2.57	0.52
1:D:107:GLN:HA	1:D:112:ILE:H	1.74	0.52
1:D:5:VAL:HG22	1:D:13:ILE:HB	1.91	0.52
2:B:148:TRP:HE1	2:B:177:SER:HB3	1.72	0.52
2:E:78:LEU:HD21	2:E:106:ILE:CD1	2.37	0.52
2:E:188:LYS:C	2:E:189:HIS:CG	2.83	0.52
3:F:93:ALA:HB1	3:F:100(H):MET:HB3	1.92	0.52
1:A:7:THR:OG1	1:A:11:LYS:HB3	2.10	0.52
1:A:8:LEU:HD11	4:C:245:HOH:O	2.10	0.52
2:B:123:GLU:H	2:B:123:GLU:CD	2.13	0.52
3:C:185:PRO:HG2	3:C:188:SER:CB	2.40	0.51
2:B:122:ASP:O	2:B:126:LYS:HG3	2.10	0.51
3:C:31:ASN:ND2	4:C:227:HOH:O	2.43	0.51
2:E:40:PRO:HG3	2:E:165:GLU:HG3	1.92	0.51
2:B:158:ASN:HD22	2:B:179:LEU:HD11	1.74	0.51
2:E:151:ASP:N	2:E:191:VAL:HB	2.26	0.51
1:A:118:ARG:HB3	1:A:146:VAL:O	2.11	0.51
1:A:146:VAL:HA	1:A:147:LEU:HD23	1.93	0.51
2:B:46:LEU:HD23	2:B:55:TYR:CG	2.45	0.51
3:F:159:LEU:HD21	3:F:182:VAL:HG11	1.92	0.51
2:E:150:VAL:HG13	2:E:192:TYR:CE1	2.46	0.51
2:E:190:LYS:HB3	2:E:211:ARG:CA	2.41	0.51
1:A:130:ARG:NH1	1:A:134:ASP:HB3	2.27	0.51
3:C:12:VAL:O	3:C:111:VAL:HA	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:ASP:OD1	1:A:116:GLN:HG3	2.11	0.50
2:B:124:GLN:HE22	2:B:131:SER:HB3	1.73	0.50
3:C:3:GLN:OE1	4:C:260:HOH:O	2.19	0.50
3:F:68:THR:OG1	3:F:81:GLN:HB3	2.11	0.50
2:E:150:VAL:CG1	2:E:192:TYR:CE1	2.95	0.50
3:C:100:TRP:N	4:C:241:HOH:O	2.01	0.50
3:C:12:VAL:HG11	3:C:82(C):LEU:HD13	1.94	0.50
1:A:36:ILE:HD12	1:A:69:LEU:HD21	1.94	0.50
2:E:151:ASP:OD1	2:E:191:VAL:CB	2.60	0.50
3:C:209:LYS:HD2	2:E:156:SER:O	2.12	0.49
3:F:123:PRO:HD3	3:F:209:LYS:NZ	2.26	0.49
1:D:110:GLU:HA	4:D:165:HOH:O	2.13	0.49
1:D:56:LEU:O	1:D:61:ILE:HD12	2.12	0.49
3:C:40:ALA:HB3	3:C:43:LYS:HB2	1.95	0.49
1:D:61:ILE:HG23	1:D:65:SER:HB2	1.95	0.49
1:A:98:THR:HA	1:A:131:THR:HA	1.94	0.49
3:F:126:PRO:HG3	3:F:138:LEU:CB	2.33	0.49
3:F:12:VAL:HG11	3:F:82(C):LEU:HD12	1.94	0.49
2:E:12:SER:HA	2:E:105:GLU:O	2.12	0.49
2:E:85:THR:OG1	2:E:103:LYS:HD2	2.13	0.49
1:D:25:ASN:O	1:D:29:LYS:HG3	2.13	0.48
3:C:11:LEU:HB2	3:C:147:PRO:HG3	1.96	0.48
2:E:78:LEU:CD2	2:E:106:ILE:CD1	2.91	0.48
2:E:190:LYS:CD	2:E:211:ARG:HA	2.44	0.48
2:E:150:VAL:CG2	2:E:192:TYR:CD1	2.92	0.48
1:D:66:THR:HG23	4:D:227:HOH:O	2.14	0.48
2:B:113:PRO:HB3	2:B:139:PHE:HB3	1.96	0.48
3:C:64:LYS:C	3:C:64:LYS:HD3	2.33	0.48
2:E:209:PHE:CD1	2:E:209:PHE:C	2.87	0.48
2:B:35:TRP:CE3	2:B:73:LEU:HD12	2.49	0.48
3:C:184:VAL:HB	3:C:185:PRO:HD2	1.95	0.48
3:C:39:GLN:HG3	3:C:44:GLY:O	2.14	0.48
2:E:209:PHE:HD1	2:E:209:PHE:O	1.91	0.48
1:D:109:LYS:N	4:D:160:HOH:O	2.47	0.47
1:D:5:VAL:HG12	1:D:67:LEU:HB2	1.95	0.47
2:E:78:LEU:HG	2:E:106:ILE:CD1	2.44	0.47
2:E:180:THR:HG21	3:F:143:LYS:HZ3	1.79	0.47
1:A:8:LEU:HD21	1:A:70:VAL:HG12	1.96	0.47
2:B:125:LEU:CD2	2:B:130:ALA:H	2.27	0.47
1:D:85:THR:OG1	1:D:87:LYS:HE2	2.14	0.47
3:C:83:ARG:HD3	3:C:85:GLU:OE2	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:190:LYS:CG	2:E:211:ARG:HA	2.44	0.47
2:B:121:SER:HB3	2:B:124:GLN:HB2	1.97	0.47
3:F:201:LYS:N	3:F:202:PRO:CD	2.77	0.47
1:A:105:LYS:NZ	4:A:164:HOH:O	2.24	0.47
1:D:135:TYR:HB2	1:D:137:ILE:HD13	1.96	0.47
2:B:154:LEU:H	2:B:154:LEU:HD12	1.78	0.47
2:B:198:HIS:HD2	2:B:200:GLY:H	1.59	0.47
3:F:110:THR:HG23	4:F:253:HOH:O	2.15	0.47
3:C:96:TRP:CD1	3:C:97:LEU:HD13	2.50	0.47
1:D:127:GLU:OE1	1:D:130:ARG:HG2	2.14	0.47
2:E:164:THR:HG23	3:F:166:PHE:CE1	2.50	0.47
1:A:13:ILE:HD12	1:A:13:ILE:N	2.30	0.47
2:E:180:THR:C	2:E:181:LEU:HD12	2.35	0.47
2:E:18:ARG:CD	4:E:250:HOH:O	2.63	0.46
2:B:174:SER:O	3:C:166:PHE:HE1	1.97	0.46
2:E:39:LYS:HB3	2:E:40:PRO:CD	2.45	0.46
2:B:46:LEU:HD23	2:B:55:TYR:CD1	2.51	0.46
2:B:39:LYS:HB3	2:B:40:PRO:HD2	1.97	0.46
2:E:140:TYR:CG	2:E:141:PRO:HA	2.51	0.46
2:B:125:LEU:HD23	2:B:130:ALA:N	2.30	0.46
2:B:183:LYS:HE2	2:B:187:GLU:OE2	2.16	0.46
3:C:100:TRP:HA	4:C:241:HOH:O	2.13	0.46
1:A:130:ARG:HH11	1:A:134:ASP:HB3	1.80	0.45
2:B:175:LEU:HD23	2:B:176:SER:N	2.30	0.45
2:B:73:LEU:HD23	2:B:74:THR:N	2.31	0.45
2:E:24:ARG:HA	2:E:69:THR:O	2.16	0.45
1:D:63:LYS:HG3	1:D:64:GLU:HG2	1.98	0.45
2:E:190:LYS:HG2	2:E:211:ARG:C	2.36	0.45
3:C:30:SER:HA	4:C:226:HOH:O	2.15	0.45
3:C:82:MET:HB3	3:C:82(C):LEU:HD21	1.99	0.45
1:D:26:VAL:HG21	1:D:56:LEU:HD21	1.97	0.45
1:D:42:ARG:HB2	1:D:72:ARG:CG	2.36	0.45
2:E:151:ASP:OD1	2:E:191:VAL:HB	2.16	0.45
2:E:48:ILE:CD1	2:E:54:LEU:HD23	2.47	0.45
2:E:63:SER:O	2:E:73:LEU:HD12	2.16	0.45
3:F:4:LEU:HD21	3:F:27:PHE:HZ	1.81	0.45
3:C:148:GLU:HA	3:C:149:PRO:HA	1.49	0.45
2:E:113:PRO:HB3	2:E:139:PHE:HB3	1.99	0.45
2:E:38:GLN:O	2:E:84:ALA:HB1	2.16	0.45
3:F:125:ALA:HA	3:F:126:PRO:HD3	1.83	0.45
3:F:171:GLN:HB2	3:F:175:LEU:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:6:LYS:CG	1:D:12:THR:HG22	2.47	0.44
2:E:164:THR:HG23	3:F:166:PHE:CD1	2.52	0.44
2:E:23:CYS:HB2	2:E:35:TRP:CH2	2.52	0.44
2:B:129:THR:HG22	2:B:130:ALA:C	2.38	0.44
2:B:118:PHE:CD1	3:C:124:LEU:HB3	2.53	0.44
1:D:120:ILE:HG22	1:D:144:HIS:O	2.18	0.44
2:B:158:ASN:HD22	2:B:179:LEU:CD1	2.30	0.44
2:B:159:SER:HB3	2:B:179:LEU:HD12	1.98	0.44
3:C:143:LYS:HG3	4:C:264:HOH:O	2.18	0.44
3:C:199:ASN:HD21	3:C:201:LYS:HZ1	1.63	0.44
3:C:74:SER:CB	4:C:265:HOH:O	2.65	0.44
2:E:149:LYS:NZ	2:E:154:LEU:HD21	2.32	0.44
1:D:108:ASP:N	4:D:160:HOH:O	2.46	0.44
1:D:110:GLU:CA	4:D:165:HOH:O	2.64	0.44
3:C:131:THR:HG22	3:C:132:SER:N	2.32	0.44
2:E:6:GLN:NE2	2:E:88:CYS:H	2.15	0.44
2:B:24:ARG:HA	2:B:69:THR:O	2.18	0.43
2:B:192:TYR:HB2	2:B:209:PHE:CE1	2.53	0.43
2:E:141:PRO:O	2:E:198:HIS:HE1	1.99	0.43
3:F:139:GLY:HA2	3:F:154:TRP:CH2	2.53	0.43
1:A:0:HIS:NE2	1:A:16:GLU:HB3	2.34	0.43
3:C:96:TRP:CD1	3:C:97:LEU:CD1	3.01	0.43
1:D:118:ARG:HD3	1:D:146:VAL:HB	2.00	0.43
2:E:180:THR:O	2:E:181:LEU:HD12	2.18	0.43
2:B:190:LYS:NZ	2:B:191:VAL:HG23	2.33	0.43
1:A:146:VAL:HA	1:A:147:LEU:HA	1.62	0.43
1:A:1:MET:HG3	1:A:63:LYS:HA	2.01	0.43
2:B:148:TRP:HB2	2:B:155:GLN:HB2	2.00	0.43
3:C:178:LEU:C	3:C:178:LEU:HD12	2.39	0.43
2:E:140:TYR:CD2	2:E:141:PRO:HA	2.54	0.43
1:D:138:GLN:HA	3:F:54:SER:O	2.19	0.43
2:E:91:SER:HB2	3:F:100:TRP:HB2	2.01	0.43
3:F:163:VAL:HG22	3:F:182:VAL:HG22	2.01	0.43
2:B:125:LEU:HD23	2:B:129:THR:C	2.38	0.43
3:C:94:ARG:HD2	3:C:101:ASP:OD1	2.19	0.43
1:D:99:ILE:HG13	1:D:126:LEU:HB3	2.00	0.43
2:E:108:ARG:HG3	2:E:140:TYR:CD1	2.54	0.43
3:F:119:PRO:CB	3:F:145:TYR:HB3	2.49	0.43
2:E:125:LEU:C	2:E:127:SER:H	2.23	0.42
2:E:54:LEU:HD21	2:E:62:PHE:O	2.19	0.42
2:B:129:THR:HG22	2:B:130:ALA:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:96:TRP:C	3:C:96:TRP:CD1	2.89	0.42
1:D:120:ILE:CG2	1:D:144:HIS:HB2	2.49	0.42
1:D:22:THR:OG1	1:D:25:ASN:HB2	2.19	0.42
3:F:153:SER:OG	3:F:197:ASN:HB2	2.19	0.42
2:B:148:TRP:HZ2	2:B:177:SER:O	2.01	0.42
1:D:139:LYS:HD3	3:F:56:GLN:HG2	2.02	0.42
3:F:146:PHE:HA	3:F:147:PRO:HA	1.73	0.42
1:D:23:ILE:HD12	1:D:50:LEU:CD2	2.38	0.42
2:E:113:PRO:HD3	2:E:198:HIS:CD2	2.53	0.42
2:B:124:GLN:HE21	2:B:129:THR:HG21	1.85	0.42
1:D:18:GLU:O	1:D:56:LEU:HD12	2.20	0.42
1:D:45:PHE:HB2	1:D:67:LEU:HD22	2.02	0.42
1:D:8:LEU:HD13	3:F:96:TRP:CE3	2.54	0.42
1:A:84:LEU:HD13	1:A:145:LEU:O	2.20	0.42
2:B:141:PRO:O	2:B:198:HIS:HE1	2.02	0.42
3:F:200:HIS:HB3	3:F:205:THR:HB	2.00	0.42
1:A:109:LYS:HB3	1:A:109:LYS:HE2	1.83	0.42
1:A:0:HIS:CE1	1:A:16:GLU:OE2	2.72	0.42
3:C:74:SER:HB3	4:C:265:HOH:O	2.20	0.42
2:B:130:ALA:O	2:B:181:LEU:N	2.53	0.41
2:B:132:VAL:HG13	2:B:179:LEU:HB3	2.02	0.41
3:C:20:LEU:HD12	3:C:80:LEU:HD23	2.01	0.41
2:E:131:SER:HA	2:E:179:LEU:O	2.18	0.41
2:E:18:ARG:HD3	4:E:250:HOH:O	2.20	0.41
3:F:12:VAL:HG11	3:F:82(C):LEU:CD1	2.49	0.41
3:F:178:LEU:HD12	3:F:178:LEU:C	2.40	0.41
3:C:166:PHE:HA	3:C:167:PRO:HD3	1.95	0.41
1:D:71:LEU:HD22	3:F:96:TRP:HA	2.02	0.41
2:E:167:ASP:O	2:E:171:SER:HA	2.21	0.41
3:F:196:CYS:SG	3:F:209:LYS:HB3	2.61	0.41
2:B:190:LYS:HZ2	2:B:191:VAL:HG23	1.85	0.41
2:B:124:GLN:CG	2:B:129:THR:OG1	2.59	0.41
2:B:89:GLN:HG2	2:B:90:GLN:N	2.35	0.41
1:D:108:ASP:CA	4:D:160:HOH:O	2.68	0.41
3:C:2:VAL:HA	3:C:25:SER:O	2.21	0.41
2:E:130:ALA:O	2:E:180:THR:HA	2.21	0.41
2:B:154:LEU:HD12	2:B:154:LEU:N	2.35	0.41
2:E:166:GLN:HE21	2:E:171:SER:HB3	1.84	0.41
1:A:42:ARG:HG2	1:A:44:ILE:HD11	2.03	0.41
3:C:146:PHE:HA	3:C:147:PRO:HA	1.69	0.40
2:E:205:VAL:CG1	2:E:206:THR:N	2.83	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:139:GLY:HA2	3:F:154:TRP:CZ2	2.57	0.40
2:B:55:TYR:CZ	3:C:101:ASP:HB2	2.57	0.40
1:D:18:GLU:O	1:D:21:ASP:HB2	2.21	0.40
3:F:33:TYR:HB2	3:F:95:THR:HB	2.02	0.40
3:C:188:SER:O	3:C:192:GLN:HB3	2.21	0.40
3:C:59:TYR:CZ	3:C:69:ILE:HG22	2.56	0.40
3:C:201:LYS:N	3:C:202:PRO:CD	2.84	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	147/172 (86%)	142 (97%)	5 (3%)	0	100	100
1	D	143/172 (83%)	133 (93%)	10 (7%)	0	100	100
2	B	209/214 (98%)	196 (94%)	13 (6%)	0	100	100
2	E	209/214 (98%)	192 (92%)	16 (8%)	1 (0%)	29	40
3	C	216/227 (95%)	205 (95%)	10 (5%)	1 (0%)	29	40
3	F	216/227 (95%)	202 (94%)	14 (6%)	0	100	100
All	All	1140/1226 (93%)	1070 (94%)	68 (6%)	2 (0%)	47	61

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	149	PRO
2	E	39	LYS

### 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	135/153 (88%)	134 (99%)	1 (1%)	84	92
1	D	131/153 (86%)	131 (100%)	0	100	100
2	B	186/188 (99%)	185 (100%)	1 (0%)	88	95
2	E	186/188 (99%)	186 (100%)	0	100	100
3	C	182/191 (95%)	181 (100%)	1 (0%)	88	95
3	F	182/191 (95%)	180 (99%)	2 (1%)	73	86
All	All	1002/1064 (94%)	997 (100%)	5 (0%)	88	95

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105	LYS
2	B	129	THR
3	C	96	TRP
3	F	96	TRP
3	F	98	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	GLN
1	A	101	ASN
1	A	116	GLN
2	B	27	GLN
2	B	124	GLN
2	B	155	GLN
2	B	158	ASN
2	B	189	HIS
2	B	198	HIS
2	B	210	ASN
3	C	31	ASN
3	C	171	GLN

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Mol	Chain	Res	Type
3	C	197	ASN
3	C	199	ASN
3	C	200	HIS
1	D	41	GLN
1	D	68	HIS
2	E	6	GLN
2	E	189	HIS
2	E	198	HIS
3	F	3	GLN
3	F	56	GLN
3	F	197	ASN
3	F	200	HIS

### 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 carbohydrates 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

### 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	149/172 (86%)	-0.07	2 (1%) 77 75	15, 34, 53, 67	0
1	D	145/172 (84%)	0.25	2 (1%) 75 73	22, 49, 64, 79	0
2	B	211/214 (98%)	0.27	10 (4%) 31 29	17, 39, 69, 75	0
2	E	211/214 (98%)	0.15	10 (4%) 31 29	22, 42, 75, 87	0
3	C	218/227 (96%)	0.27	11 (5%) 28 26	18, 42, 65, 84	0
3	F	218/227 (96%)	0.08	2 (0%) 84 82	20, 41, 62, 75	0
All	All	1152/1226 (93%)	0.17	37 (3%) 47 45	15, 41, 66, 87	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	129	THR	5.1
1	A	84	LEU	4.6
2	E	128	GLY	4.4
3	C	130	SER	4.1
3	C	191	THR	3.4
2	E	181	LEU	3.3
2	B	158	ASN	3.3
2	B	153	ALA	3.3
2	B	181	LEU	3.1
2	B	130	ALA	3.1
2	E	154	LEU	3.0
3	C	190	GLY	3.0
2	E	187	GLU	2.7
2	B	157	GLY	2.6
2	B	184	ALA	2.6
3	C	189	LEU	2.6
3	C	136	ALA	2.6
1	D	10	GLY	2.6
3	F	190	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
2	E	183	LYS	2.4
1	A	85	THR	2.4
2	E	190	LYS	2.4
3	C	206	LYS	2.4
3	F	171	GLN	2.3
1	D	62	GLN	2.3
2	B	192	TYR	2.3
2	E	191	VAL	2.2
2	E	123	GLU	2.2
2	E	188	LYS	2.2
2	E	189	HIS	2.1
2	B	178	THR	2.1
2	B	149	LYS	2.1
3	C	185	PRO	2.0
3	C	115	SER	2.0
3	C	118	GLY	2.0
3	C	186	SER	2.0
3	C	208	ASP	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 carbohydrates 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.