



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:05 PM UTC

PDB ID : 3U30 / pdb_00003u30
Title : Crystal structure of a linear-specific Ubiquitin fab bound to linear ubiquitin
Authors : Matsumoto, M.L.; Dong, K.C.; Yu, C.; Phu, L.; Gao, X.; Hannoush, R.N.;
Hymowitz, S.G.; Kirkpatrick, D.S.; Dixit, V.M.; Kelley, R.F.
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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

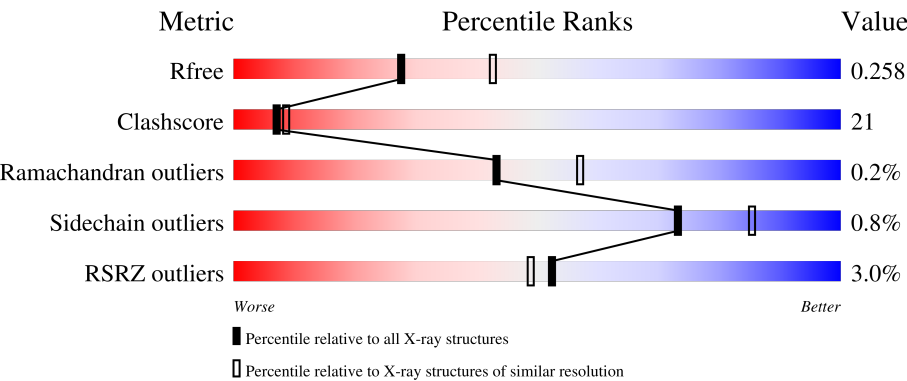
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

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}	180053	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (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 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	172	4% 68% 18% • 13%
1	D	172	3% 51% 33% • 16%
2	B	214	4% 71% 26% ••
2	E	214	4% 63% 34% ••
3	C	227	2% 63% 32% ••

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Mol	Chain	Length	Quality of chain
3	F	227	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: a green segment representing 70%, a yellow segment representing 25%, and a small grey segment at the end. Below the bar, the percentages '70%' and '25%' are labeled under their respective segments. To the right of the bar, there are two small black dots.

2 Entry composition [i](#)

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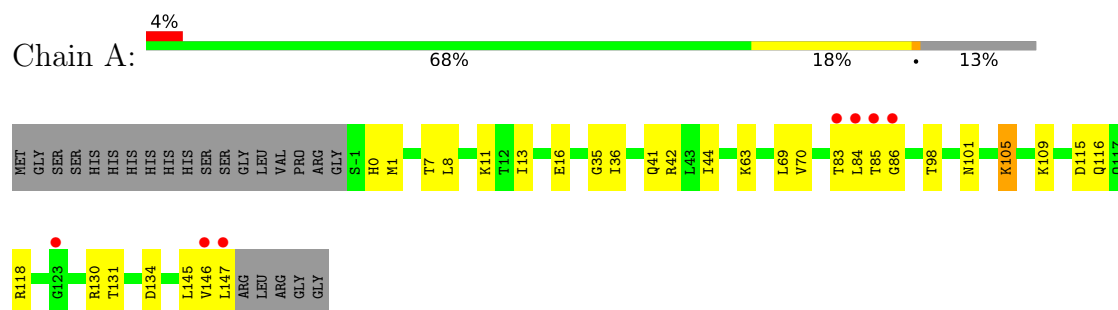
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	22	Total 22	O 22	0	0
4	E	42	Total 42	O 42	0	0
4	F	33	Total 33	O 33	0	0

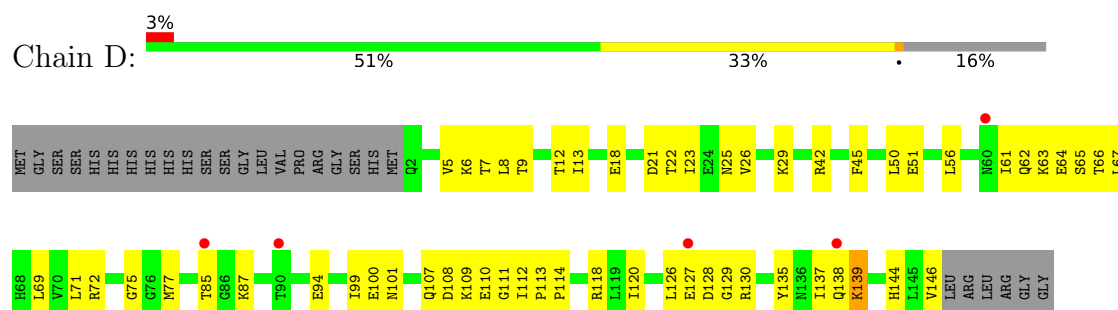
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, 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.

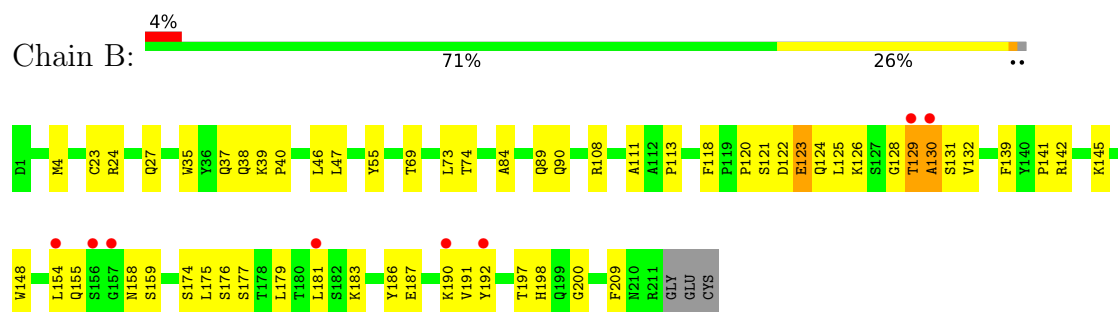
- Molecule 1: linear di-ubiquitin



- Molecule 1: linear di-ubiquitin

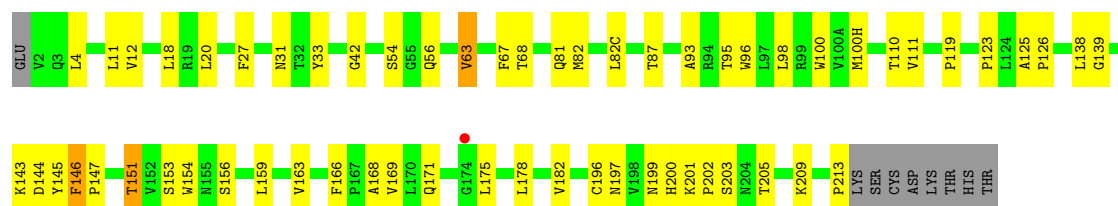


- Molecule 2: Light chain Fab



- Molecule 2: Light chain Fab





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	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) 94.8 (49.33-2.43)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 2.42Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.229 , 0.258 0.229 , 0.258	Depositor DCC
R_{free} test set	1957 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.350	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	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 ²)	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.*

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

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	0.26	0/1194	0.68	0/1609
1	D	0.28	0/1161	0.70	2/1565 (0.1%)
2	B	0.27	0/1656	0.79	2/2251 (0.1%)
2	E	0.38	1/1656 (0.1%)	0.76	3/2251 (0.1%)
3	C	0.26	0/1657	0.77	3/2264 (0.1%)
3	F	0.28	0/1657	0.72	2/2264 (0.1%)
All	All	0.29	1/8981 (0.0%)	0.74	12/12204 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	40	PRO	N-CD	-8.89	1.35	1.47

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	148	GLU	C-N-CD	-12.62	92.84	120.60
2	B	130	ALA	N-CA-CB	-12.29	91.96	110.40
2	B	130	ALA	N-CA-C	12.02	144.66	111.00
2	E	182	SER	N-CA-C	6.73	121.88	112.72
3	C	146	PHE	CB-CA-C	6.45	119.12	109.63
3	C	150	VAL	N-CA-C	6.14	118.23	108.89
3	F	144	ASP	N-CA-C	6.10	118.05	110.91
2	E	191	VAL	CA-C-N	-5.85	114.86	123.05
2	E	191	VAL	C-N-CA	-5.85	114.86	123.05
3	F	146	PHE	CB-CA-C	5.77	118.12	109.63
1	D	113	PRO	CA-C-N	5.09	124.26	118.97
1	D	113	PRO	C-N-CA	5.09	124.26	118.97

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	73	0
2	B	1620	0	1581	65	0
2	E	1620	0	1581	93	0
3	C	1618	0	1592	69	0
3	F	1618	0	1592	57	0
4	A	22	0	0	6	0
4	B	30	0	0	3	0
4	C	45	0	0	19	0
4	D	22	0	0	21	0
4	E	42	0	0	8	0
4	F	33	0	0	9	0
All	All	8998	0	8761	367	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 21.

All (367) 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)
3:C:148:GLU:HG3	3:C:149:PRO:CB	1.54	1.37
2:B:129:THR:CA	2:B:130:ALA:HB3	1.45	1.34
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:101:ASN:CG	4:D:153:HOH:O	2.09	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:O	2:E:209:PHE:CD1	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:D:107:GLN:C	4:D:160:HOH:O	2.24	0.79
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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
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:37:GLN:HB2	2:B:47:LEU:HD11	1.73	0.69
2:B:198:HIS:CD2	2:B:200:GLY:H	2.11	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
1:A:101:ASN:CG	4:A:164:HOH:O	2.36	0.67
3:C:94:ARG:NH1	3:C:101:ASP:OD1	2.27	0.67
1:D:108:ASP:C	4:D:160:HOH:O	2.38	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
3:F:18:LEU:HB3	3:F:82:MET:HE3	1.76	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
2:E:52:LYS:NZ	4:E:237:HOH:O	1.81	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:23:ILE:HG23	1:D:50:LEU:HD22	1.80	0.63
1:D:118:ARG:HH11	1:D:146:VAL:HG11	1.62	0.63
2:B:124:GLN:HE22	2:B:131:SER:HB2	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:9:THR:O	1:D:9:THR:HG22	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
1:D:77:MET:HE3	1:D:139:LYS:HA	1.83	0.61
2:E:6:GLN:HE21	2:E:99:GLY:HA3	1.66	0.61
2:E:198:HIS:CD2	2:E:200:GLY:H	2.19	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:130:ALA:HB1	2:B:181:LEU:C	2.27	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
1:D:127:GLU:CD	1:D:129:GLY:H	2.12	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:B:129:THR:CA	2:B:130:ALA:CB	2.31	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:B:123:GLU:H	2:B:123:GLU:CD	2.13	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:39:LYS:HB3	2:E:40:PRO:HD2	1.88	0.56
2:E:188:LYS:C	2:E:189:HIS:CG	2.83	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
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
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:7:THR:HG22	1:D:69:LEU:HD23	1.89	0.53
1:D:107:GLN:O	1:D:111:GLY:HA2	2.09	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:9:THR:O	1:D:9:THR:CG2	2.57	0.53
2:E:78:LEU:CD2	2:E:106:ILE:HD13	2.36	0.53
3:C:64:LYS:C	3:C:64:LYS:HD3	2.33	0.52
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:209:PHE:CD1	2:E:209:PHE:C	2.87	0.52
2:E:210:ASN:O	2:E:211:ARG:CG	2.57	0.52
3:F:123:PRO:HD3	3:F:209:LYS:HZ1	1.74	0.52
1:D:5:VAL:HG22	1:D:13:ILE:HB	1.91	0.52
1:D:107:GLN:HA	1:D:112:ILE:H	1.74	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
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
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
3:C:82:MET:HE1	3:C:109:VAL:HG21	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:190:LYS:HZ2	2:B:191:VAL:HG23	1.75	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
2:E:180:THR:C	2:E:181:LEU:HD12	2.35	0.51
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
2:E:190:LYS:HG2	2:E:211:ARG:C	2.36	0.50
3:C:12:VAL:HG11	3:C:82(C):LEU:HD13	1.94	0.50
3:C:100:TRP:N	4:C:241:HOH:O	2.01	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:56:LEU:O	1:D:61:ILE:HD12	2.12	0.49
1:D:110:GLU:HA	4:D:165:HOH:O	2.13	0.49
2:B:129:THR:HG22	2:B:130:ALA:C	2.38	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:113:PRO:HB3	2:B:139:PHE:HB3	1.96	0.48
2:E:4:MET:HE3	2:E:23:CYS:SG	2.53	0.48
2:B:35:TRP:CE3	2:B:73:LEU:HD12	2.49	0.48
3:C:39:GLN:HG3	3:C:44:GLY:O	2.14	0.48
3:C:178:LEU:HD12	3:C:178:LEU:C	2.39	0.48
3:C:184:VAL:HB	3:C:185:PRO:HD2	1.95	0.48
2:E:209:PHE:O	2:E:209:PHE:HD1	1.91	0.48
2:B:4:MET:HE3	2:B:23:CYS:SG	2.54	0.47
1:D:5:VAL:HG12	1:D:67:LEU:HB2	1.95	0.47
1:D:109:LYS:N	4:D:160:HOH:O	2.47	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:HD23	2:B:129:THR:C	2.38	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
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
2:E:125:LEU:C	2:E:127:SER:H	2.23	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
3:C:96:TRP:C	3:C:96:TRP:CD1	2.89	0.47
1:D:107:GLN:O	4:D:160:HOH:O	2.18	0.47
1:D:135:TYR:HB2	1:D:137:ILE:HD13	1.96	0.47
2:B:154:LEU:HD12	2:B:154:LEU:H	1.78	0.47
2:B:198:HIS:HD2	2:B:200:GLY:H	1.59	0.47
3:C:96:TRP:CD1	3:C:97:LEU:HD13	2.50	0.47
3:F:110:THR:HG23	4:F:253:HOH:O	2.15	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:N	1:A:13:ILE:HD12	2.30	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
3:F:178:LEU:C	3:F:178:LEU:HD12	2.40	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:73:LEU:HD23	2:B:74:THR:N	2.31	0.45
2:B:175:LEU:HD23	2:B:176:SER:N	2.30	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
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: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
2:E:151:ASP:OD1	2:E:191:VAL:HB	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
1:D:6:LYS:CG	1:D:12:THR:HG22	2.47	0.44
2:E:23:CYS:HB2	2:E:35:TRP:CH2	2.52	0.44
2:E:164:THR:HG23	3:F:166:PHE:CD1	2.52	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: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:20:LEU:HG	3:F:82:MET:HE2	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
2:E:175:LEU:C	2:E:175:LEU:HD23	2.43	0.43
1:A:1:MET:HG3	1:A:63:LYS:HA	2.01	0.43
1:A:146:VAL:HA	1:A:147:LEU:HA	1.62	0.43
2:B:148:TRP:HB2	2:B:155:GLN:HB2	2.00	0.43
2:E:140:TYR:CD2	2:E:141:PRO:HA	2.54	0.43
3:C:39:GLN:C	3:C:88:ALA:HB1	2.44	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
3:C:61:ASP:C	3:C:63:VAL:H	2.27	0.43
3:C:94:ARG:HD2	3:C:101:ASP:OD1	2.19	0.43
3:C:199:ASN:HD21	3:C:201:LYS:HZ1	1.64	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: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
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:8:LEU:HD13	3:F:96:TRP:CE3	2.54	0.42
1:D:45:PHE:HB2	1:D:67:LEU:HD22	2.02	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:0:HIS:CE1	1:A:16:GLU:OE2	2.72	0.42
1:A:109:LYS:HB3	1:A:109:LYS:HE2	1.83	0.42
3:C:74:SER:HB3	4:C:265:HOH:O	2.20	0.42
2:B:27:GLN:CD	4:B:232:HOH:O	2.53	0.41
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:18:ARG:HD3	4:E:250:HOH:O	2.20	0.41
2:E:131:SER:HA	2:E:179:LEU:O	2.18	0.41
1:D:87:LYS:NZ	1:D:110:GLU:CD	2.74	0.41
3:F:12:VAL:HG11	3:F:82(C):LEU:CD1	2.49	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:89:GLN:HG2	2:B:90:GLN:N	2.35	0.41
2:B:124:GLN:CG	2:B:129:THR:OG1	2.59	0.41
2:B:154:LEU:HD12	2:B:154:LEU:N	2.35	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: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
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

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
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%)	24	35
3	C	216/227 (95%)	205 (95%)	10 (5%)	1 (0%)	24	35
3	F	216/227 (95%)	202 (94%)	14 (6%)	0	100	100
All	All	1140/1226 (93%)	1070 (94%)	68 (6%)	2 (0%)	43	57

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%)	76	87
1	D	131/153 (86%)	130 (99%)	1 (1%)	73	85
2	B	186/188 (99%)	184 (99%)	2 (1%)	65	81
2	E	186/188 (99%)	185 (100%)	1 (0%)	81	90
3	C	182/191 (95%)	182 (100%)	0	100	100
3	F	182/191 (95%)	179 (98%)	3 (2%)	55	74
All	All	1002/1064 (94%)	994 (99%)	8 (1%)	73	85

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105	LYS
2	B	123	GLU
2	B	129	THR
1	D	139	LYS

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Mol	Chain	Res	Type
2	E	19	VAL
3	F	63	VAL
3	F	98	LEU
3	F	151	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	101	ASN
1	A	116	GLN
2	B	27	GLN
2	B	124	GLN
2	B	137	ASN
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	164	HIS
3	C	197	ASN
3	C	200	HIS
1	D	41	GLN
1	D	60	ASN
1	D	68	HIS
2	E	6	GLN
2	E	189	HIS
2	E	198	HIS
2	E	199	GLN
3	F	3	GLN
3	F	164	HIS
3	F	197	ASN
3	F	200	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	149/172 (86%)	0.18	7 (4%)	36 32	15, 34, 53, 67	0
1	D	145/172 (84%)	0.68	5 (3%)	48 44	22, 49, 64, 79	0
2	B	211/214 (98%)	0.32	8 (3%)	44 40	17, 39, 69, 75	0
2	E	211/214 (98%)	0.33	8 (3%)	44 40	22, 42, 75, 87	0
3	C	218/227 (96%)	0.24	5 (2%)	61 57	18, 42, 65, 84	0
3	F	218/227 (96%)	0.23	1 (0%)	87 86	20, 41, 62, 75	0
All	All	1152/1226 (93%)	0.32	34 (2%)	52 49	15, 41, 66, 87	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	84	LEU	3.8
1	A	85	THR	3.3
2	E	189	HIS	3.2
2	B	129	THR	3.1
1	A	86	GLY	3.0
3	C	130	SER	2.5
2	B	190	LYS	2.5
1	A	146	VAL	2.4
2	E	181	LEU	2.4
2	E	186	TYR	2.4
3	C	137	ALA	2.3
3	C	189	LEU	2.3
2	B	130	ALA	2.3
2	E	153	ALA	2.3
3	F	174	GLY	2.3
2	B	181	LEU	2.2
3	C	149	PRO	2.2
1	D	85	THR	2.2
1	D	90	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	147	LEU	2.2
2	E	191	VAL	2.2
1	D	127	GLU	2.2
2	B	157	GLY	2.2
2	E	128	GLY	2.2
1	D	138	GLN	2.1
1	A	123	GLY	2.1
2	B	192	TYR	2.1
2	E	183	LYS	2.1
3	C	114	ALA	2.1
2	B	154	LEU	2.1
1	A	83	THR	2.1
2	E	187	GLU	2.1
1	D	60	ASN	2.0
2	B	156	SER	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.