



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:24 PM GMT

PDB ID : 1K47
Title : Crystal Structure of the Streptococcus pneumoniae Phosphomevalonate Kinase (PMK)
Authors : Romanowski, M.J.; Bonanno, J.B.; Burley, S.K.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2001-10-05
Resolution : 2.42 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
<http://wwpdb.org/validation/2016/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	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

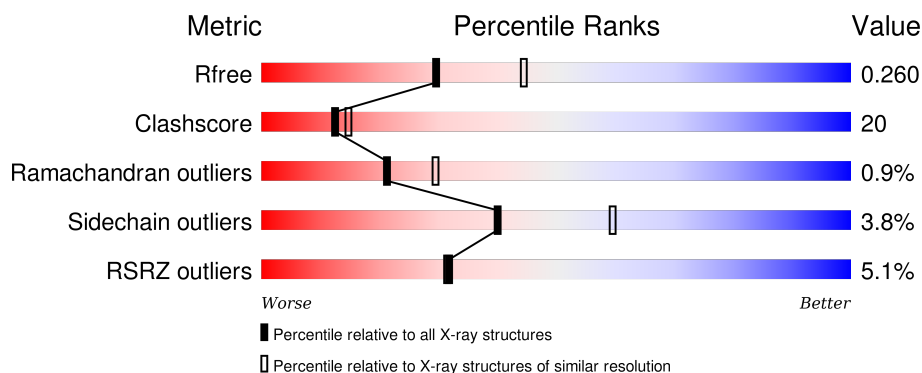
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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}	91344	3386 (2.44-2.40)
Clashscore	102246	3897 (2.44-2.40)
Ramachandran outliers	100387	3837 (2.44-2.40)
Sidechain outliers	100360	3838 (2.44-2.40)
RSRZ outliers	91569	3396 (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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	337	<div> <div>5%</div> <div>64%</div> <div>31%</div> <div>• •</div> </div>
1	B	337	<div> <div>5%</div> <div>65%</div> <div>27%</div> <div>• •</div> </div>
1	C	337	<div> <div>5%</div> <div>66%</div> <div>27%</div> <div>• •</div> </div>
1	D	337	<div> <div>6%</div> <div>63%</div> <div>30%</div> <div>• •</div> </div>
1	E	337	<div> <div>4%</div> <div>61%</div> <div>33%</div> <div>• •</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	337	<div><div></div><div>4%</div><div>67%</div><div>26%</div><div></div><div></div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 15779 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 phosphomevalonate kinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	Se	0	0	0
			2500	1595	413	480	4	8			
1	B	322	Total	C	N	O	S	Se	0	0	0
			2500	1595	413	480	4	8			
1	C	322	Total	C	N	O	S	Se	0	0	0
			2500	1595	413	480	4	8			
1	D	322	Total	C	N	O	S	Se	0	0	0
			2500	1595	413	480	4	8			
1	E	322	Total	C	N	O	S	Se	0	0	0
			2500	1595	413	480	4	8			
1	F	322	Total	C	N	O	S	Se	0	0	0
			2500	1595	413	480	4	8			

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1001	HIS	-	CLONING ARTIFACT	UNP Q8DR49
A	1002	MSE	-	CLONING ARTIFACT	UNP Q8DR49
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
A	34	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
A	50	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
A	73	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
A	95	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
A	148	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
A	216	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
B	1001	HIS	-	CLONING ARTIFACT	UNP Q8DR49
B	1002	MSE	-	CLONING ARTIFACT	UNP Q8DR49
B	1	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
B	34	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
B	50	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
B	73	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
B	95	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
B	148	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49

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Chain	Residue	Modelled	Actual	Comment	Reference
B	216	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
C	1001	HIS	-	CLONING ARTIFACT	UNP Q8DR49
C	1002	MSE	-	CLONING ARTIFACT	UNP Q8DR49
C	1	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
C	34	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
C	50	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
C	73	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
C	95	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
C	148	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
C	216	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
D	1001	HIS	-	CLONING ARTIFACT	UNP Q8DR49
D	1002	MSE	-	CLONING ARTIFACT	UNP Q8DR49
D	1	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
D	34	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
D	50	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
D	73	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
D	95	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
D	148	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
D	216	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
E	1001	HIS	-	CLONING ARTIFACT	UNP Q8DR49
E	1002	MSE	-	CLONING ARTIFACT	UNP Q8DR49
E	1	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
E	34	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
E	50	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
E	73	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
E	95	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
E	148	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
E	216	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
F	1001	HIS	-	CLONING ARTIFACT	UNP Q8DR49
F	1002	MSE	-	CLONING ARTIFACT	UNP Q8DR49
F	1	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
F	34	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
F	50	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
F	73	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
F	95	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
F	148	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49
F	216	MSE	MET	MODIFIED RESIDUE	UNP Q8DR49

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	128	Total O 128 128	0	0

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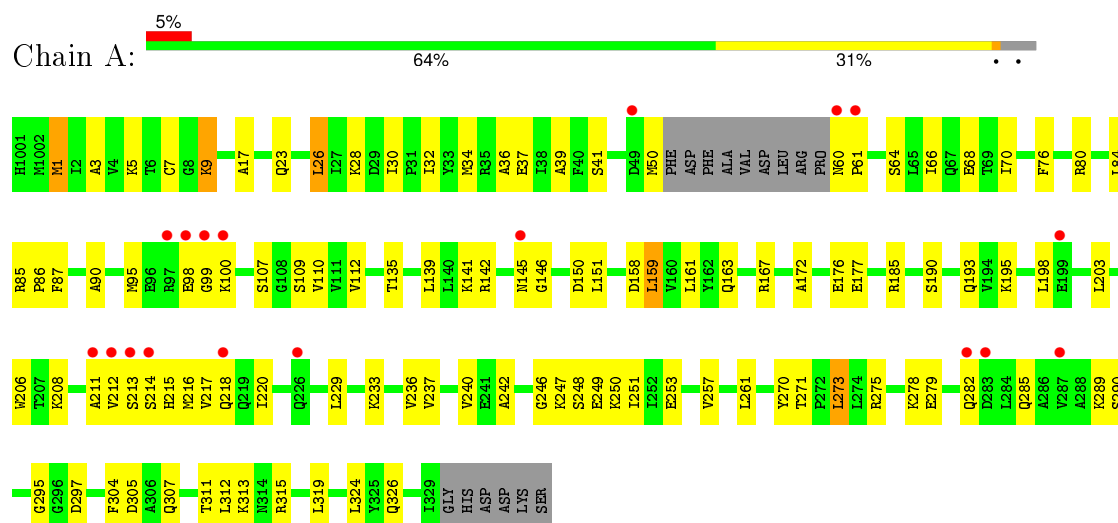
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	133	Total 133	O 133	0	0
2	C	122	Total 122	O 122	0	0
2	D	136	Total 136	O 136	0	0
2	E	131	Total 131	O 131	0	0
2	F	129	Total 129	O 129	0	0

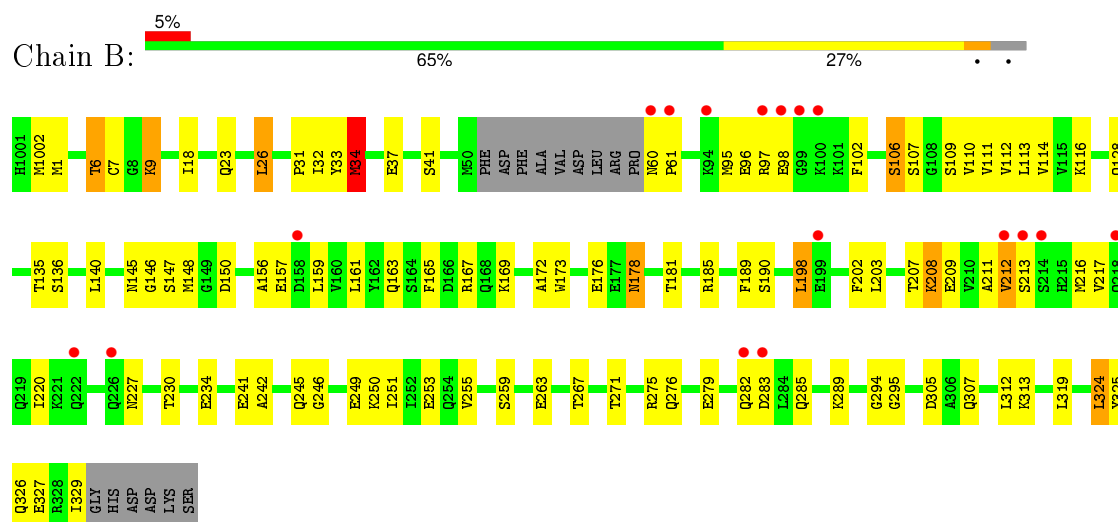
3 Residue-property plots [i](#)

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

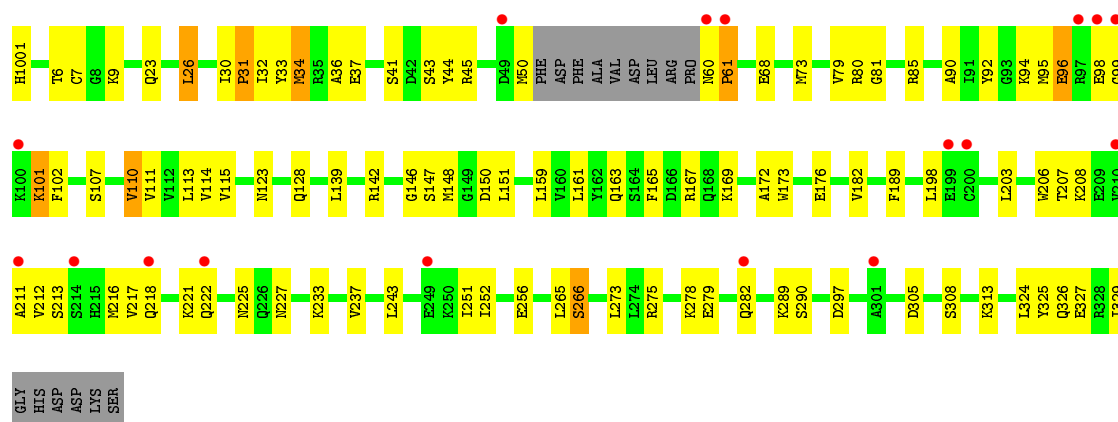


- Molecule 1: phosphomevalonate kinase



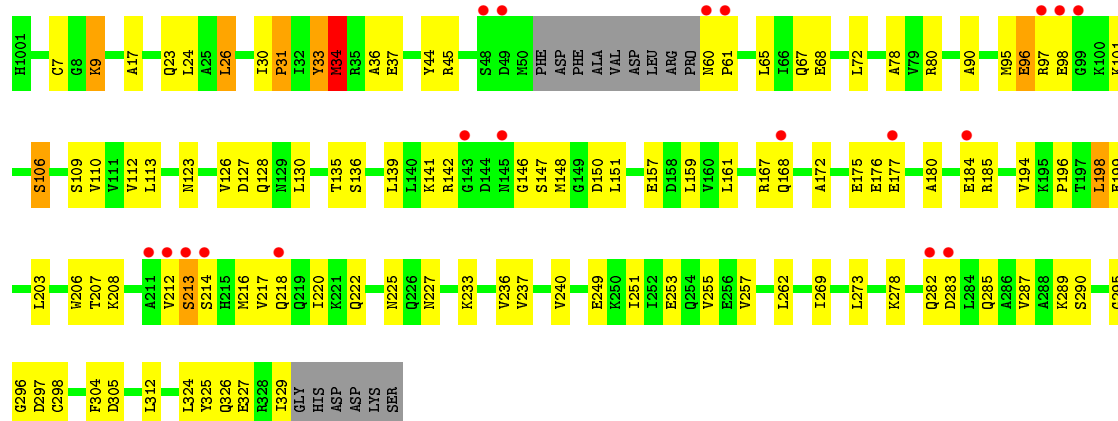
- Molecule 1: phosphomevalonate kinase





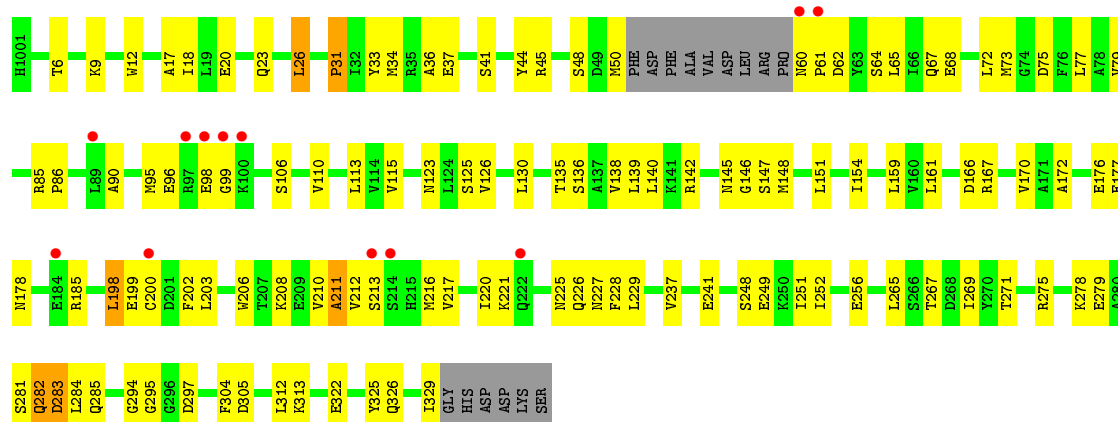
- Molecule 1: phosphomevalonate kinase

Chain D: 6% 63% 30%



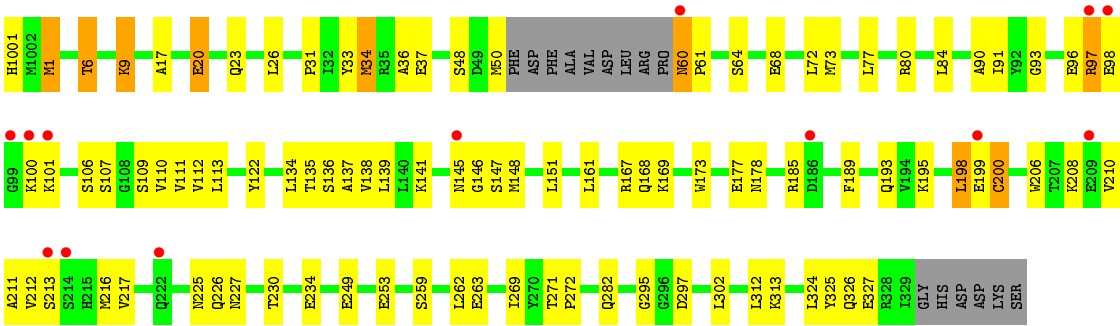
- Molecule 1: phosphomevalonate kinase

Chain E: 4% 61% 33%



- Molecule 1: phosphomevalonate kinase

Chain F: 4% 67% 26%



4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	231.61Å 231.61Å 88.59Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.42 20.11 – 2.42	Depositor EDS
% Data completeness (in resolution range)	95.0 (20.00-2.42) 97.7 (20.11-2.42)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 2.41Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.215 , 0.257 0.218 , 0.260	Depositor DCC
R_{free} test set	9706 reflections (9.61%)	DCC
Wilson B-factor (Å ²)	35.6	Xtriage
Anisotropy	0.074	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 41.5	EDS
Estimated twinning fraction	0.017 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 198848 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	15779	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.37	0/2529	0.62	2/3401 (0.1%)
1	B	0.38	0/2529	0.63	2/3401 (0.1%)
1	C	0.38	0/2529	0.63	1/3401 (0.0%)
1	D	0.36	0/2529	0.61	2/3401 (0.1%)
1	E	0.38	0/2529	0.63	1/3401 (0.0%)
1	F	0.38	0/2529	0.64	2/3401 (0.1%)
All	All	0.37	0/15174	0.63	10/20406 (0.0%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	33	TYR	N-CA-C	6.23	127.83	111.00
1	C	33	TYR	N-CA-C	6.04	127.30	111.00
1	A	159	LEU	N-CA-C	-5.61	95.86	111.00
1	B	33	TYR	N-CA-C	5.60	126.12	111.00
1	B	34	MSE	N-CA-C	-5.30	96.69	111.00
1	F	34	MSE	N-CA-C	-5.29	96.73	111.00
1	F	33	TYR	N-CA-C	5.10	124.77	111.00
1	D	34	MSE	N-CA-C	-5.08	97.28	111.00
1	A	32	ILE	N-CA-C	-5.03	97.42	111.00
1	D	33	TYR	N-CA-C	5.01	124.52	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	2500	0	2539	93	0
1	B	2500	0	2539	99	0
1	C	2500	0	2539	90	0
1	D	2500	0	2539	104	0
1	E	2500	0	2539	118	0
1	F	2500	0	2539	95	0
2	A	128	0	0	16	0
2	B	133	0	0	18	0
2	C	122	0	0	16	0
2	D	136	0	0	18	0
2	E	131	0	0	19	0
2	F	129	0	0	16	0
All	All	15779	0	15234	590	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 (590) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:225:ASN:HD21	1:D:227:ASN:HB2	1.05	1.13
1:A:213:SER:H	1:A:216:MSE:HE3	1.20	1.06
1:E:60:ASN:HB2	1:E:61:PRO:HD3	1.38	1.05
1:F:148:MSE:HE1	1:F:151:LEU:HD12	1.34	1.04
1:F:148:MSE:CE	1:F:151:LEU:HD12	1.88	1.03
1:D:34:MSE:HE1	1:D:110:VAL:HA	1.35	1.01
1:B:212:VAL:HG13	1:B:216:MSE:HB2	1.44	0.99
1:C:60:ASN:HB2	1:C:61:PRO:HD3	1.45	0.98
1:C:213:SER:H	1:C:216:MSE:HE3	1.27	0.96
1:D:60:ASN:HB2	1:D:61:PRO:HD3	1.50	0.94
1:A:212:VAL:HG13	1:A:216:MSE:HB2	1.50	0.92
1:E:212:VAL:HG13	1:E:216:MSE:HB2	1.50	0.92
1:A:60:ASN:HB2	1:A:61:PRO:HD3	1.51	0.92
1:E:95:MSE:HE1	1:E:106:SER:HA	1.53	0.90
1:D:225:ASN:ND2	1:D:227:ASN:HB2	1.87	0.88
1:F:193:GLN:HE21	1:F:195:LYS:HG2	1.37	0.88
1:F:177:GLU:HG3	1:F:185:ARG:HH21	1.38	0.88
1:D:212:VAL:HG13	1:D:216:MSE:HB2	1.57	0.86
1:D:136:SER:HA	2:D:1103:HOH:O	1.74	0.85
1:C:212:VAL:HG22	1:C:216:MSE:HB2	1.58	0.85
1:F:313:LYS:NZ	1:F:326:GLN:HE22	1.73	0.85
1:A:275:ARG:HG3	2:A:1112:HOH:O	1.75	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:6:THR:HB	2:B:1069:HOH:O	1.74	0.84
1:B:60:ASN:HB2	1:B:61:PRO:HD3	1.58	0.83
1:D:177:GLU:HG3	1:D:185:ARG:NH2	1.92	0.82
1:D:213:SER:H	1:D:216:MSE:CE	1.92	0.82
1:A:146:GLY:O	1:A:167:ARG:HD3	1.79	0.82
1:B:113:LEU:HB2	2:B:1069:HOH:O	1.79	0.81
1:D:34:MSE:HE3	1:D:113:LEU:HD22	1.62	0.81
1:A:98:GLU:HB2	2:A:1081:HOH:O	1.79	0.81
1:C:278:LYS:HD3	2:C:1057:HOH:O	1.80	0.81
1:C:113:LEU:HD13	2:C:1065:HOH:O	1.81	0.80
1:C:218:GLN:O	1:C:222:GLN:HG3	1.82	0.80
1:E:267:THR:HB	2:E:1078:HOH:O	1.80	0.80
1:F:34:MSE:HB3	2:F:1062:HOH:O	1.81	0.80
1:D:34:MSE:CE	1:D:110:VAL:HA	2.11	0.79
1:B:249:GLU:O	1:B:253:GLU:HG3	1.83	0.79
1:E:213:SER:H	1:E:216:MSE:HE3	1.45	0.78
1:C:146:GLY:O	1:C:167:ARG:HD3	1.83	0.78
1:B:185:ARG:HD3	2:B:1047:HOH:O	1.84	0.78
1:C:213:SER:H	1:C:216:MSE:CE	1.97	0.77
1:E:6:THR:HB	2:E:1064:HOH:O	1.86	0.76
1:F:177:GLU:HG3	1:F:185:ARG:NH2	2.00	0.76
1:E:185:ARG:HD3	2:E:1042:HOH:O	1.84	0.76
1:E:34:MSE:HE3	1:E:113:LEU:HD22	1.68	0.75
1:C:92:TYR:HA	1:C:96:GLU:OE1	1.86	0.75
1:C:34:MSE:HE1	1:C:110:VAL:HG22	1.67	0.74
1:B:207:THR:O	1:B:209:GLU:N	2.20	0.74
2:C:1045:HOH:O	1:E:45:ARG:HD3	1.87	0.73
1:C:211:ALA:HB2	1:C:297:ASP:OD2	1.89	0.72
1:E:34:MSE:CE	1:E:110:VAL:HA	2.19	0.72
1:A:177:GLU:HG3	1:A:185:ARG:NH2	2.05	0.72
1:A:279:GLU:HB3	1:A:282:GLN:HE21	1.55	0.72
1:E:34:MSE:HG2	1:E:95:MSE:CB	2.20	0.71
1:E:282:GLN:O	1:E:284:LEU:N	2.22	0.71
1:D:26:LEU:HD22	1:D:159:LEU:HD11	1.71	0.71
1:B:212:VAL:HG13	1:B:216:MSE:CB	2.20	0.71
1:B:172:ALA:O	1:B:176:GLU:HG3	1.91	0.70
1:E:67:GLN:HG3	2:F:1092:HOH:O	1.91	0.70
1:D:68:GLU:OE1	1:D:142:ARG:HG3	1.90	0.70
2:A:1052:HOH:O	1:D:45:ARG:HD3	1.91	0.70
1:D:34:MSE:HG2	1:D:95:MSE:HB3	1.73	0.70
1:E:34:MSE:HG2	1:E:95:MSE:HB3	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:126:VAL:HB	1:D:130:LEU:HD23	1.73	0.69
1:D:206:TRP:CZ3	1:D:208:LYS:HG2	2.27	0.69
1:D:146:GLY:O	1:D:167:ARG:HD3	1.92	0.69
1:F:20:GLU:HB2	2:F:1090:HOH:O	1.90	0.69
1:D:225:ASN:ND2	1:D:227:ASN:H	1.91	0.69
1:F:169:LYS:HE2	1:F:173:TRP:CH2	2.27	0.69
1:C:213:SER:N	1:C:216:MSE:HE3	2.05	0.69
1:D:285:GLN:HB3	1:D:305:ASP:OD1	1.93	0.69
1:D:212:VAL:HG23	1:D:295:GLY:HA3	1.74	0.69
1:F:100:LYS:HG2	1:F:106:SER:HB3	1.74	0.69
1:A:213:SER:O	1:A:217:VAL:HG23	1.93	0.69
1:C:32:ILE:HG22	1:C:95:MSE:HG2	1.75	0.69
1:F:193:GLN:HE21	1:F:195:LYS:CG	2.07	0.68
1:F:193:GLN:NE2	1:F:195:LYS:HE2	2.08	0.68
1:B:242:ALA:HB2	1:B:250:LYS:HG3	1.75	0.68
1:B:208:LYS:HE3	2:B:1008:HOH:O	1.94	0.68
1:B:106:SER:O	1:B:110:VAL:HG23	1.94	0.68
1:F:34:MSE:SE	1:F:110:VAL:HG22	2.44	0.68
1:C:213:SER:OG	1:C:216:MSE:HG3	1.93	0.67
1:E:211:ALA:HB2	1:E:297:ASP:OD2	1.94	0.67
1:F:212:VAL:HG13	1:F:216:MSE:HB2	1.76	0.67
2:C:1022:HOH:O	1:E:37:GLU:HG3	1.93	0.67
1:D:213:SER:OG	1:D:216:MSE:HG3	1.95	0.67
1:D:218:GLN:O	1:D:222:GLN:HG3	1.95	0.67
1:E:139:LEU:HD12	1:E:151:LEU:HD11	1.77	0.67
1:F:213:SER:OG	1:F:216:MSE:HE3	1.95	0.67
1:B:34:MSE:HE3	1:B:113:LEU:HD22	1.77	0.66
1:B:169:LYS:HE2	1:B:173:TRP:CZ2	2.30	0.66
1:D:213:SER:H	1:D:216:MSE:HE3	1.58	0.66
1:E:279:GLU:O	1:E:282:GLN:HG3	1.95	0.66
1:E:248:SER:HA	1:E:251:ILE:HD12	1.77	0.66
1:F:60:ASN:HB2	1:F:61:PRO:HD3	1.77	0.66
1:F:230:THR:HG22	1:F:234:GLU:OE2	1.95	0.66
1:D:213:SER:H	1:D:216:MSE:HE2	1.58	0.66
1:D:212:VAL:CG1	1:D:216:MSE:HB2	2.24	0.66
1:E:198:LEU:HB3	2:E:1129:HOH:O	1.96	0.66
2:C:1068:HOH:O	1:E:50:MSE:HE2	1.96	0.66
1:C:233:LYS:O	1:C:237:VAL:HG23	1.96	0.66
1:D:24:LEU:HD22	1:D:161:LEU:HD21	1.78	0.66
1:E:313:LYS:NZ	1:E:326:GLN:HE22	1.94	0.65
1:A:216:MSE:O	1:A:220:ILE:HG13	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:185:ARG:HD3	2:D:1041:HOH:O	1.96	0.65
1:B:26:LEU:HD22	1:B:159:LEU:HD11	1.78	0.65
1:E:123:ASN:HB3	2:E:1061:HOH:O	1.97	0.65
1:D:253:GLU:O	1:D:257:VAL:HG23	1.97	0.65
1:E:60:ASN:HB2	1:E:61:PRO:CD	2.21	0.64
1:F:113:LEU:HD13	2:F:1062:HOH:O	1.96	0.64
1:C:289:LYS:HG2	1:C:290:SER:O	1.98	0.64
1:E:212:VAL:HG23	1:E:295:GLY:HA3	1.79	0.64
1:E:95:MSE:CE	1:E:106:SER:HA	2.25	0.64
1:D:95:MSE:HE1	1:D:106:SER:HA	1.80	0.64
1:E:98:GLU:HB2	2:E:1085:HOH:O	1.97	0.64
1:F:145:ASN:HB3	2:F:1060:HOH:O	1.96	0.64
1:D:67:GLN:HE21	1:E:210:VAL:HG23	1.62	0.64
1:F:168:GLN:HG2	2:F:1077:HOH:O	1.97	0.64
1:F:198:LEU:HD22	1:F:199:GLU:O	1.97	0.64
1:D:198:LEU:HD23	1:D:199:GLU:H	1.63	0.64
1:D:9:LYS:HE3	1:D:150:ASP:HB3	1.80	0.63
1:A:279:GLU:HB3	1:A:282:GLN:NE2	2.13	0.63
1:A:212:VAL:HG12	1:A:217:VAL:HG23	1.79	0.63
1:F:148:MSE:HE1	1:F:151:LEU:CD1	2.18	0.63
1:C:282:GLN:HA	2:C:1083:HOH:O	1.99	0.63
1:C:9:LYS:HE3	1:C:150:ASP:HB3	1.81	0.63
1:D:249:GLU:HG3	2:D:1051:HOH:O	1.99	0.62
1:D:216:MSE:O	1:D:220:ILE:HG13	1.98	0.62
1:F:325:TYR:CE2	1:F:327:GLU:HB2	2.33	0.62
1:D:278:LYS:HE3	2:D:1112:HOH:O	2.00	0.62
1:B:34:MSE:CE	1:B:110:VAL:HA	2.29	0.62
1:E:279:GLU:HB3	2:E:1090:HOH:O	1.99	0.62
1:B:140:LEU:HG	1:B:148:MSE:HE2	1.81	0.62
1:E:249:GLU:HG2	2:E:1055:HOH:O	1.98	0.62
1:E:206:TRP:CZ3	1:E:208:LYS:HG2	2.35	0.62
1:F:169:LYS:HE2	1:F:173:TRP:CZ2	2.35	0.62
2:C:1113:HOH:O	1:E:48:SER:HB2	1.99	0.62
1:B:213:SER:H	1:B:216:MSE:HE3	1.64	0.61
1:E:213:SER:O	1:E:217:VAL:HG23	2.00	0.61
1:B:216:MSE:O	1:B:220:ILE:HG13	1.98	0.61
1:A:107:SER:HB2	2:A:1101:HOH:O	1.99	0.61
1:B:136:SER:CB	1:B:148:MSE:HE3	2.31	0.61
1:E:65:LEU:HD13	1:E:142:ARG:NH1	2.15	0.61
1:F:211:ALA:HB2	1:F:297:ASP:OD2	2.00	0.61
1:B:34:MSE:HG2	1:B:95:MSE:HB2	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:106:SER:O	1:F:110:VAL:HG23	2.00	0.61
1:E:146:GLY:O	1:E:167:ARG:HD3	2.00	0.61
1:A:145:ASN:HB3	2:A:1079:HOH:O	2.00	0.61
1:C:6:THR:HB	2:C:1065:HOH:O	2.01	0.60
1:B:212:VAL:HG12	1:B:217:VAL:HG23	1.81	0.60
1:D:98:GLU:HB2	2:D:1078:HOH:O	1.99	0.60
1:E:136:SER:HB3	1:E:148:MSE:HE3	1.83	0.60
1:F:212:VAL:HG12	1:F:217:VAL:HG23	1.82	0.60
1:A:313:LYS:NZ	1:A:326:GLN:HE22	2.00	0.60
1:F:313:LYS:HZ1	1:F:326:GLN:HE22	1.50	0.60
1:E:283:ASP:N	2:E:1073:HOH:O	2.33	0.60
1:A:257:VAL:O	1:A:261:LEU:HG	2.00	0.60
1:C:34:MSE:HE1	1:C:110:VAL:HG13	1.84	0.60
2:A:1007:HOH:O	1:C:43:SER:HB2	2.02	0.59
1:A:139:LEU:HD12	1:A:151:LEU:HD11	1.84	0.59
1:C:107:SER:O	1:C:110:VAL:HG23	2.01	0.59
2:B:1102:HOH:O	1:F:48:SER:HB2	2.01	0.59
1:C:98:GLU:HB2	2:C:1076:HOH:O	2.02	0.59
1:A:198:LEU:HD21	1:A:246:GLY:O	2.02	0.59
1:B:34:MSE:HG3	1:B:95:MSE:HE3	1.84	0.59
1:B:136:SER:HB3	1:B:148:MSE:HE3	1.83	0.59
1:F:64:SER:O	1:F:68:GLU:HG3	2.03	0.59
1:B:212:VAL:CG1	1:B:216:MSE:HB2	2.26	0.59
1:D:34:MSE:HG2	1:D:95:MSE:CB	2.33	0.59
1:E:113:LEU:HB2	2:E:1064:HOH:O	2.02	0.59
1:F:6:THR:HG21	1:F:109:SER:O	2.03	0.59
1:A:34:MSE:CE	1:A:109:SER:HB2	2.33	0.59
1:A:64:SER:O	1:A:68:GLU:HG3	2.02	0.59
1:E:199:GLU:O	1:E:200:CYS:SG	2.61	0.58
1:A:198:LEU:HG	1:A:246:GLY:HA2	1.84	0.58
1:A:17:ALA:O	1:A:23:GLN:HG3	2.02	0.58
1:E:271:THR:HG23	1:E:295:GLY:HA2	1.85	0.58
1:B:140:LEU:HD21	1:B:167:ARG:HG2	1.85	0.58
1:A:34:MSE:SE	1:A:110:VAL:HG22	2.54	0.58
1:E:75:ASP:O	1:E:79:VAL:HG23	2.04	0.58
1:E:126:VAL:HB	1:E:130:LEU:HD23	1.84	0.58
1:E:206:TRP:CE2	1:E:208:LYS:HA	2.39	0.58
1:C:147:SER:O	1:C:148:MSE:HB2	2.03	0.58
1:C:85:ARG:HD3	2:C:1072:HOH:O	2.04	0.57
1:B:34:MSE:CE	1:B:110:VAL:HG22	2.34	0.57
1:B:95:MSE:HA	2:B:1134:HOH:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:325:TYR:CE2	1:B:327:GLU:HB2	2.39	0.57
1:A:36:ALA:HA	1:A:90:ALA:O	2.04	0.57
1:F:206:TRP:CE2	1:F:208:LYS:HA	2.39	0.57
1:A:60:ASN:HB2	1:A:61:PRO:CD	2.32	0.57
1:E:34:MSE:HE2	1:E:110:VAL:HA	1.84	0.57
1:B:140:LEU:HG	1:B:148:MSE:CE	2.34	0.57
1:D:206:TRP:HB2	1:D:324:LEU:HG	1.86	0.57
1:D:207:THR:HB	1:D:273:LEU:HD23	1.86	0.57
1:C:80:ARG:NE	1:C:80:ARG:HA	2.19	0.57
1:C:34:MSE:CE	1:C:110:VAL:HG22	2.35	0.57
1:C:213:SER:O	1:C:217:VAL:HG23	2.03	0.57
1:E:177:GLU:OE2	1:E:185:ARG:NH2	2.38	0.57
1:B:207:THR:O	1:B:207:THR:OG1	2.21	0.57
1:B:146:GLY:O	1:B:167:ARG:HD3	2.05	0.57
1:C:123:ASN:HB3	2:C:1061:HOH:O	2.04	0.57
1:B:34:MSE:HE1	1:B:110:VAL:HG22	1.85	0.57
1:F:213:SER:O	1:F:217:VAL:HG23	2.03	0.57
1:D:34:MSE:HE3	1:D:113:LEU:CD2	2.34	0.56
1:E:212:VAL:HG13	1:E:216:MSE:CB	2.31	0.56
1:A:28:LYS:HD3	1:A:159:LEU:HD13	1.86	0.56
1:C:139:LEU:HD12	1:C:151:LEU:HD11	1.87	0.56
1:B:61:PRO:HD2	1:F:50:MSE:SE	2.56	0.56
1:D:206:TRP:CZ2	1:D:208:LYS:HA	2.41	0.56
1:E:269:ILE:HG23	1:E:294:GLY:HA3	1.87	0.56
1:F:147:SER:O	1:F:148:MSE:HB2	2.05	0.56
1:A:23:GLN:NE2	1:A:229:LEU:HB3	2.21	0.56
1:B:313:LYS:NZ	1:B:326:GLN:HE22	2.03	0.56
1:A:7:CYS:HB2	1:A:30:ILE:O	2.05	0.56
1:E:285:GLN:HB2	1:E:305:ASP:OD1	2.06	0.56
1:B:41:SER:HB2	1:F:37:GLU:OE2	2.06	0.56
1:B:116:LYS:HD2	2:B:1052:HOH:O	2.05	0.56
1:C:275:ARG:O	1:C:279:GLU:HG3	2.05	0.55
1:C:26:LEU:HD22	1:C:159:LEU:HD11	1.87	0.55
1:D:225:ASN:HD21	1:D:227:ASN:CB	1.98	0.55
1:E:216:MSE:O	1:E:220:ILE:HG13	2.06	0.55
1:C:98:GLU:HB3	1:C:101:LYS:HD2	1.89	0.55
1:F:97:ARG:HB3	1:F:97:ARG:NH1	2.22	0.55
1:B:169:LYS:HE2	1:B:173:TRP:CH2	2.42	0.55
1:B:34:MSE:HE3	1:B:113:LEU:CD2	2.37	0.55
1:E:203:LEU:HD23	1:E:326:GLN:HG3	1.89	0.55
1:D:7:CYS:SG	1:D:157:GLU:HG2	2.47	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:249:GLU:CG	2:E:1055:HOH:O	2.53	0.55
1:B:9:LYS:C	1:B:9:LYS:HD2	2.28	0.55
1:D:161:LEU:C	1:D:161:LEU:HD23	2.28	0.54
1:F:72:LEU:HD23	1:F:138:VAL:HG23	1.89	0.54
1:E:203:LEU:CD2	1:E:326:GLN:HG3	2.38	0.54
1:F:1001:HIS:N	2:F:1066:HOH:O	2.40	0.54
1:B:282:GLN:O	1:B:283:ASP:HB2	2.06	0.54
1:B:241:GLU:O	1:B:245:GLN:HG3	2.07	0.54
1:C:313:LYS:NZ	1:C:326:GLN:HE22	2.05	0.54
1:B:163:GLN:HB3	1:B:190:SER:HB3	1.90	0.54
1:E:213:SER:H	1:E:216:MSE:CE	2.16	0.54
1:F:206:TRP:CZ3	1:F:208:LYS:HG2	2.43	0.54
1:E:172:ALA:O	1:E:176:GLU:HG3	2.07	0.54
1:D:225:ASN:HD22	1:D:227:ASN:H	1.55	0.54
1:C:50:MSE:SE	1:E:61:PRO:HD2	2.57	0.54
1:E:106:SER:O	1:E:110:VAL:HG23	2.07	0.54
1:B:249:GLU:HG3	2:B:1056:HOH:O	2.08	0.54
1:A:214:SER:HB2	1:A:218:GLN:HE22	1.71	0.54
1:A:248:SER:HA	1:A:251:ILE:HD12	1.89	0.54
1:B:32:ILE:HG22	1:B:95:MSE:HG2	1.88	0.54
1:F:31:PRO:HD2	1:F:325:TYR:CZ	2.43	0.54
1:A:206:TRP:CE2	1:A:208:LYS:HA	2.43	0.54
1:A:275:ARG:HD2	2:A:1062:HOH:O	2.07	0.54
1:B:98:GLU:HB2	2:B:1076:HOH:O	2.07	0.54
1:E:44:TYR:O	1:E:45:ARG:HG3	2.08	0.54
1:D:147:SER:O	1:D:148:MSE:HB2	2.07	0.54
1:A:247:LYS:HB2	1:A:250:LYS:HB3	1.89	0.53
1:D:251:ILE:O	1:D:255:VAL:HG23	2.07	0.53
1:A:1:MSE:HB3	1:A:39:ALA:HB2	1.89	0.53
1:F:112:VAL:HG22	1:F:135:THR:HG21	1.91	0.53
1:D:329:ILE:O	1:D:329:ILE:HG22	2.08	0.53
1:E:200:CYS:SG	1:E:304:PHE:CD2	3.00	0.53
1:F:17:ALA:O	1:F:23:GLN:HG3	2.08	0.53
1:A:215:HIS:HB2	2:A:1095:HOH:O	2.07	0.53
1:A:211:ALA:HB2	1:A:297:ASP:OD2	2.08	0.53
1:C:169:LYS:HE2	1:C:173:TRP:CZ2	2.43	0.53
1:E:212:VAL:HA	1:E:216:MSE:HE3	1.90	0.53
1:F:313:LYS:CE	1:F:326:GLN:HE22	2.21	0.53
1:F:282:GLN:HB3	2:F:1080:HOH:O	2.09	0.53
1:B:7:CYS:SG	1:B:157:GLU:HG2	2.49	0.53
1:B:34:MSE:HG2	1:B:95:MSE:CB	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:34:MSE:HG3	1:C:95:MSE:HE3	1.91	0.53
1:E:281:SER:O	1:E:282:GLN:O	2.26	0.53
1:E:96:GLU:O	1:E:96:GLU:HG2	2.08	0.52
1:B:60:ASN:CB	1:B:61:PRO:HD3	2.35	0.52
1:E:64:SER:O	1:E:68:GLU:HG3	2.08	0.52
1:D:206:TRP:HB2	1:D:324:LEU:CG	2.38	0.52
1:E:199:GLU:HA	1:E:329:ILE:O	2.09	0.52
1:E:206:TRP:CZ2	1:E:208:LYS:HA	2.44	0.52
1:B:275:ARG:O	1:B:279:GLU:HG3	2.09	0.52
1:C:172:ALA:O	1:C:176:GLU:HG3	2.09	0.52
1:F:212:VAL:HG23	1:F:295:GLY:HA3	1.90	0.52
1:A:242:ALA:HB2	1:A:250:LYS:HG3	1.91	0.52
1:D:123:ASN:HB3	2:D:1060:HOH:O	2.08	0.52
1:C:252:ILE:HG22	1:C:256:GLU:OE2	2.08	0.52
1:A:158:ASP:HB3	1:A:193:GLN:HG3	1.90	0.52
1:B:276:GLN:OE1	1:B:319:LEU:HD22	2.09	0.52
1:D:177:GLU:HG3	1:D:185:ARG:HH22	1.74	0.52
1:A:26:LEU:HD22	1:A:159:LEU:HD11	1.91	0.52
1:F:210:VAL:HG23	2:F:1092:HOH:O	2.07	0.52
1:C:94:LYS:HD3	2:C:1110:HOH:O	2.10	0.52
1:D:289:LYS:HG2	1:D:290:SER:O	2.10	0.52
1:F:100:LYS:HG2	1:F:106:SER:CB	2.40	0.51
1:D:198:LEU:HD23	1:D:199:GLU:N	2.25	0.51
1:E:167:ARG:HD2	2:E:1062:HOH:O	2.09	0.51
1:C:45:ARG:NH1	2:C:1046:HOH:O	2.43	0.51
1:C:146:GLY:O	1:C:167:ARG:CD	2.57	0.51
1:D:34:MSE:CG	1:D:95:MSE:HB3	2.40	0.51
1:F:313:LYS:HE2	1:F:326:GLN:NE2	2.25	0.51
1:A:253:GLU:O	1:A:257:VAL:HG23	2.10	0.51
1:C:60:ASN:HB2	1:C:61:PRO:CD	2.29	0.51
1:E:6:THR:HG23	1:E:154:ILE:HG23	1.91	0.51
1:B:60:ASN:HB2	1:B:61:PRO:CD	2.37	0.51
1:D:225:ASN:ND2	1:D:227:ASN:N	2.59	0.51
1:A:203:LEU:HD23	1:A:326:GLN:HG3	1.93	0.51
1:B:230:THR:O	1:B:234:GLU:HG3	2.11	0.51
1:E:73:MSE:O	1:E:77:LEU:HG	2.11	0.51
1:A:236:VAL:O	1:A:240:VAL:HG23	2.11	0.51
1:C:165:PHE:HB3	1:C:189:PHE:HB3	1.92	0.51
1:B:34:MSE:HE2	1:B:110:VAL:HA	1.91	0.51
1:A:172:ALA:O	1:A:176:GLU:HG3	2.10	0.51
1:D:212:VAL:HA	1:D:216:MSE:HE3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:LYS:O	1:A:237:VAL:HG23	2.11	0.51
1:A:198:LEU:HD11	1:A:246:GLY:O	2.11	0.51
1:B:178:ASN:ND2	1:B:181:THR:H	2.09	0.51
1:C:207:THR:HB	1:C:273:LEU:HD23	1.93	0.51
1:B:227:ASN:ND2	2:B:1094:HOH:O	2.43	0.51
1:F:226:GLN:HB3	2:F:1104:HOH:O	2.11	0.50
1:C:34:MSE:HG2	1:C:95:MSE:HB3	1.92	0.50
1:F:212:VAL:HG13	1:F:216:MSE:CB	2.41	0.50
1:E:26:LEU:HD22	1:E:159:LEU:HD11	1.94	0.50
2:A:1025:HOH:O	1:D:37:GLU:HG3	2.10	0.50
1:D:262:LEU:O	1:D:269:ILE:HD12	2.11	0.50
1:E:36:ALA:HA	1:E:90:ALA:O	2.11	0.50
1:F:262:LEU:HD22	1:F:269:ILE:HD13	1.93	0.50
1:F:206:TRP:HB2	1:F:324:LEU:HG	1.92	0.50
1:E:267:THR:O	1:E:267:THR:HG22	2.10	0.50
1:E:136:SER:CB	1:E:148:MSE:HE3	2.41	0.50
1:F:325:TYR:CZ	1:F:327:GLU:HB2	2.46	0.50
1:A:34:MSE:HE1	1:A:110:VAL:HG23	1.94	0.50
1:B:9:LYS:HE3	1:B:150:ASP:HB3	1.93	0.50
1:B:279:GLU:HB3	1:B:282:GLN:NE2	2.27	0.50
1:D:236:VAL:O	1:D:240:VAL:HG23	2.11	0.50
1:E:61:PRO:HG3	1:E:96:GLU:HG3	1.93	0.50
1:E:115:VAL:HG21	1:E:135:THR:HG22	1.94	0.50
1:F:80:ARG:NH2	2:F:1074:HOH:O	2.45	0.50
1:F:93:GLY:O	1:F:97:ARG:HG3	2.11	0.50
1:C:325:TYR:CE2	1:C:327:GLU:HB2	2.47	0.50
1:F:178:ASN:HB2	2:F:1030:HOH:O	2.10	0.50
1:E:217:VAL:HG12	1:E:221:LYS:HE3	1.93	0.50
1:E:281:SER:C	1:E:282:GLN:O	2.50	0.50
1:E:237:VAL:O	1:E:241:GLU:HG3	2.12	0.50
1:A:213:SER:N	1:A:216:MSE:HE3	2.06	0.49
1:E:34:MSE:HE2	1:E:110:VAL:CA	2.42	0.49
1:B:26:LEU:HD22	1:B:159:LEU:CD1	2.42	0.49
1:D:9:LYS:HD2	1:D:9:LYS:C	2.32	0.49
1:D:34:MSE:HE2	1:D:109:SER:C	2.32	0.49
1:B:112:VAL:HG22	1:B:135:THR:HG21	1.94	0.49
1:E:34:MSE:HE3	1:E:113:LEU:CD2	2.40	0.49
1:C:31:PRO:HD2	1:C:325:TYR:CZ	2.47	0.49
1:B:259:SER:O	1:B:263:GLU:HG3	2.13	0.49
1:A:279:GLU:O	1:A:282:GLN:HG3	2.12	0.49
1:C:31:PRO:HG2	1:C:325:TYR:OH	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:18:ILE:HG13	1:E:18:ILE:O	2.13	0.49
1:F:193:GLN:NE2	1:F:195:LYS:HG2	2.19	0.49
1:A:249:GLU:O	1:A:253:GLU:HG3	2.13	0.48
1:C:225:ASN:ND2	1:C:227:ASN:HB2	2.28	0.48
1:D:180:ALA:O	1:D:184:GLU:HG3	2.13	0.48
1:B:213:SER:O	1:B:217:VAL:HG23	2.12	0.48
1:E:213:SER:N	1:E:216:MSE:HE3	2.23	0.48
1:B:213:SER:CB	1:B:216:MSE:HE3	2.43	0.48
1:C:173:TRP:HB3	1:C:182:VAL:HG22	1.95	0.48
1:A:206:TRP:CZ2	1:A:208:LYS:HA	2.49	0.48
1:D:175:GLU:HG3	1:D:175:GLU:O	2.13	0.48
1:B:37:GLU:HG3	2:F:1003:HOH:O	2.13	0.48
1:E:198:LEU:HD11	1:E:248:SER:HB3	1.95	0.48
1:D:298:CYS:SG	1:D:324:LEU:HD12	2.54	0.48
1:C:7:CYS:HB2	1:C:30:ILE:O	2.13	0.48
1:F:36:ALA:HA	1:F:90:ALA:O	2.13	0.48
1:F:137:ALA:O	1:F:141:LYS:HG3	2.14	0.48
1:D:233:LYS:O	1:D:237:VAL:HG23	2.14	0.48
1:C:206:TRP:CE2	1:C:208:LYS:HA	2.48	0.48
1:A:275:ARG:O	1:A:279:GLU:HG3	2.14	0.48
1:E:12:TRP:HH2	1:E:329:ILE:HD13	1.79	0.48
1:B:203:LEU:HD23	1:B:326:GLN:HG3	1.95	0.48
1:C:313:LYS:HZ3	1:C:326:GLN:HE22	1.62	0.47
1:D:325:TYR:CZ	1:D:327:GLU:HB2	2.49	0.47
1:A:311:THR:O	1:A:315:ARG:HG3	2.14	0.47
1:C:217:VAL:HG12	1:C:221:LYS:HE3	1.96	0.47
1:C:203:LEU:HD23	1:C:326:GLN:HG3	1.95	0.47
1:D:112:VAL:HG22	1:D:135:THR:HG21	1.95	0.47
1:D:17:ALA:O	1:D:23:GLN:HG3	2.14	0.47
1:C:68:GLU:OE1	1:C:142:ARG:HG3	2.15	0.47
1:B:34:MSE:HB2	2:B:1069:HOH:O	2.13	0.47
1:B:102:PHE:CG	1:B:324:LEU:HD11	2.49	0.47
1:A:37:GLU:HG3	2:D:1004:HOH:O	2.15	0.47
1:E:213:SER:OG	1:E:216:MSE:HG3	2.15	0.47
1:F:185:ARG:HD3	2:F:1042:HOH:O	2.14	0.47
2:B:1027:HOH:O	1:F:37:GLU:HG3	2.14	0.47
1:B:267:THR:HG23	1:B:275:ARG:HH22	1.80	0.47
1:D:213:SER:O	1:D:217:VAL:HG23	2.14	0.47
1:C:252:ILE:O	1:C:256:GLU:HG3	2.15	0.47
1:F:73:MSE:O	1:F:77:LEU:HG	2.15	0.47
1:A:66:ILE:O	1:A:70:ILE:HG13	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:329:ILE:C	2:E:1060:HOH:O	2.53	0.47
1:C:34:MSE:HE1	1:C:110:VAL:CG2	2.41	0.46
1:A:193:GLN:NE2	1:A:195:LYS:HE2	2.29	0.46
1:A:278:LYS:HE3	2:A:1044:HOH:O	2.15	0.46
1:E:34:MSE:CE	1:E:110:VAL:HG22	2.46	0.46
1:C:206:TRP:CH2	1:C:208:LYS:HG2	2.50	0.46
1:E:147:SER:O	1:E:148:MSE:HB2	2.15	0.46
1:D:285:GLN:HG3	1:D:304:PHE:HB2	1.97	0.46
1:D:161:LEU:HD22	2:D:1021:HOH:O	2.14	0.46
1:E:228:PHE:CZ	1:E:265:LEU:HD22	2.50	0.46
1:B:97:ARG:HD3	2:B:1083:HOH:O	2.15	0.46
1:D:139:LEU:HB2	2:D:1103:HOH:O	2.16	0.46
1:C:289:LYS:HG2	1:C:290:SER:N	2.31	0.46
1:C:94:LYS:HE3	2:C:1098:HOH:O	2.15	0.46
1:A:198:LEU:HG	1:A:246:GLY:CA	2.46	0.46
1:D:127:ASP:HB2	2:D:1122:HOH:O	2.15	0.46
1:A:289:LYS:HG2	1:A:290:SER:O	2.14	0.46
1:D:106:SER:O	1:D:110:VAL:HG23	2.16	0.46
1:B:271:THR:HG23	1:B:295:GLY:HA2	1.98	0.46
1:A:34:MSE:HG3	1:A:95:MSE:HB2	1.98	0.46
1:A:9:LYS:HE3	1:A:150:ASP:HB3	1.98	0.46
1:C:289:LYS:CG	1:C:290:SER:N	2.79	0.46
1:D:98:GLU:HB3	1:D:101:LYS:HD2	1.98	0.46
1:D:30:ILE:HA	1:D:31:PRO:HD3	1.80	0.46
1:E:17:ALA:O	1:E:23:GLN:HG3	2.16	0.46
1:F:148:MSE:HE3	1:F:151:LEU:HD12	1.89	0.45
1:D:139:LEU:HD12	1:D:151:LEU:HD11	1.99	0.45
1:F:230:THR:O	1:F:234:GLU:HG3	2.16	0.45
1:C:329:ILE:O	1:C:329:ILE:HG22	2.14	0.45
1:F:6:THR:CG2	1:F:109:SER:OG	2.64	0.45
1:C:34:MSE:HE3	1:C:110:VAL:N	2.31	0.45
1:A:34:MSE:HE3	1:A:109:SER:CB	2.47	0.45
1:D:203:LEU:CD2	1:D:326:GLN:HG3	2.47	0.45
1:B:107:SER:O	1:B:111:VAL:HG23	2.16	0.45
1:E:140:LEU:CD2	1:E:148:MSE:HE2	2.46	0.45
1:E:200:CYS:SG	1:E:304:PHE:CE2	3.08	0.45
1:F:9:LYS:C	1:F:9:LYS:HD2	2.37	0.45
1:D:282:GLN:O	1:D:283:ASP:HB2	2.16	0.45
1:F:113:LEU:HD22	2:F:1062:HOH:O	2.17	0.45
1:D:282:GLN:HA	2:D:1093:HOH:O	2.16	0.45
1:D:95:MSE:HE2	1:D:109:SER:CB	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:270:TYR:C	2:A:1112:HOH:O	2.55	0.45
1:D:126:VAL:CB	1:D:130:LEU:HD23	2.43	0.45
1:E:208:LYS:HG3	2:E:1131:HOH:O	2.17	0.45
1:C:36:ALA:HA	1:C:90:ALA:O	2.17	0.45
1:B:34:MSE:HE1	1:B:110:VAL:HA	1.99	0.45
1:C:41:SER:HB2	1:E:37:GLU:OE2	2.17	0.45
1:A:271:THR:HB	2:A:1039:HOH:O	2.17	0.45
1:C:265:LEU:O	1:C:266:SER:HB3	2.16	0.45
1:A:41:SER:HB2	1:D:37:GLU:OE2	2.16	0.44
1:A:163:GLN:HB3	1:A:190:SER:HB3	1.99	0.44
1:A:305:ASP:OD2	1:A:307:GLN:HB3	2.17	0.44
1:F:139:LEU:HD12	1:F:148:MSE:HE1	1.99	0.44
1:B:110:VAL:O	1:B:114:VAL:HG23	2.17	0.44
1:E:206:TRP:HE3	1:E:322:GLU:OE1	2.00	0.44
1:A:9:LYS:C	1:A:9:LYS:HD2	2.38	0.44
1:E:226:GLN:NE2	1:E:229:LEU:HD12	2.32	0.44
1:B:202:PHE:HB2	1:B:329:ILE:CD1	2.48	0.44
1:B:211:ALA:HB1	1:B:294:GLY:O	2.17	0.44
1:C:73:MSE:SE	1:C:115:VAL:HG13	2.67	0.44
1:F:1:MSE:HE2	1:F:37:GLU:HB3	2.00	0.44
1:D:214:SER:HB2	2:D:1100:HOH:O	2.16	0.44
1:B:34:MSE:HE3	1:B:34:MSE:HB3	1.90	0.44
1:D:44:TYR:O	1:D:45:ARG:HG3	2.16	0.44
1:A:3:ALA:HB2	1:A:37:GLU:OE1	2.17	0.44
1:F:249:GLU:O	1:F:253:GLU:HG3	2.18	0.44
1:C:305:ASP:OD1	1:C:308:SER:HB2	2.18	0.44
1:A:279:GLU:CB	1:A:282:GLN:HE21	2.27	0.44
1:B:98:GLU:CB	2:B:1076:HOH:O	2.65	0.44
1:B:165:PHE:HB3	1:B:189:PHE:HB3	1.99	0.44
1:E:275:ARG:O	1:E:279:GLU:HG3	2.18	0.44
1:A:85:ARG:CZ	2:A:1046:HOH:O	2.66	0.44
1:F:6:THR:HB	2:F:1062:HOH:O	2.18	0.44
1:E:139:LEU:HB2	1:E:148:MSE:HE1	1.99	0.44
1:C:225:ASN:HD22	1:C:227:ASN:H	1.66	0.44
1:D:141:LYS:NZ	2:D:1123:HOH:O	2.50	0.44
1:C:102:PHE:N	1:C:102:PHE:CD2	2.83	0.44
1:C:206:TRP:CZ3	1:C:208:LYS:HG2	2.53	0.43
1:E:278:LYS:HE3	2:E:1043:HOH:O	2.17	0.43
1:A:50:MSE:HE2	2:D:1070:HOH:O	2.17	0.43
1:E:34:MSE:HG2	1:E:95:MSE:HB2	1.97	0.43
1:B:167:ARG:NH1	2:B:1067:HOH:O	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:96:GLU:C	1:F:98:GLU:H	2.21	0.43
1:E:198:LEU:HD23	1:E:199:GLU:N	2.32	0.43
1:D:212:VAL:CG2	1:D:295:GLY:HA3	2.46	0.43
1:F:313:LYS:CE	1:F:326:GLN:NE2	2.81	0.43
1:A:193:GLN:HE21	1:A:195:LYS:HG2	1.84	0.43
1:A:315:ARG:O	1:A:319:LEU:HG	2.19	0.43
1:C:37:GLU:HG3	2:E:1004:HOH:O	2.19	0.43
1:F:189:PHE:C	1:F:189:PHE:CD1	2.91	0.43
1:D:96:GLU:O	1:D:98:GLU:N	2.52	0.43
1:F:96:GLU:O	1:F:98:GLU:N	2.52	0.43
1:D:33:TYR:O	1:D:95:MSE:HB2	2.19	0.43
1:B:178:ASN:HD22	1:B:178:ASN:C	2.21	0.43
1:B:128:GLN:HE21	1:B:156:ALA:HA	1.83	0.43
1:D:36:ALA:HA	1:D:90:ALA:O	2.19	0.43
1:B:1002:MSE:O	1:F:1001:HIS:HA	2.19	0.43
1:D:78:ALA:HA	2:D:1053:HOH:O	2.18	0.43
1:B:251:ILE:O	1:B:255:VAL:HG23	2.18	0.43
1:E:50:MSE:HG2	1:E:50:MSE:O	2.18	0.43
1:B:146:GLY:O	1:B:167:ARG:CD	2.66	0.43
1:A:212:VAL:CG1	1:A:216:MSE:HB2	2.36	0.43
1:E:34:MSE:HE1	1:E:110:VAL:HA	1.98	0.43
1:F:225:ASN:ND2	1:F:227:ASN:HB2	2.34	0.43
1:F:136:SER:CB	1:F:148:MSE:HE2	2.49	0.42
1:F:313:LYS:HZ3	1:F:326:GLN:HE22	1.60	0.42
1:B:26:LEU:CD2	1:B:159:LEU:HD11	2.47	0.42
1:B:26:LEU:HD23	1:B:26:LEU:HA	1.90	0.42
1:B:147:SER:O	1:B:148:MSE:HB2	2.19	0.42
1:A:76:PHE:O	1:A:80:ARG:HG2	2.19	0.42
1:B:6:THR:HG22	1:B:109:SER:HB3	2.00	0.42
1:C:6:THR:CB	2:C:1065:HOH:O	2.65	0.42
1:C:92:TYR:HA	1:C:96:GLU:CD	2.39	0.42
1:E:6:THR:HG23	1:E:154:ILE:CG2	2.49	0.42
1:E:313:LYS:HZ3	1:E:326:GLN:HE22	1.65	0.42
1:A:313:LYS:HZ3	1:A:326:GLN:HE22	1.67	0.42
1:E:20:GLU:HB3	1:E:23:GLN:CG	2.49	0.42
1:D:172:ALA:O	1:D:176:GLU:HG3	2.19	0.42
1:A:141:LYS:NZ	2:A:1091:HOH:O	2.51	0.42
1:E:282:GLN:O	1:E:283:ASP:C	2.58	0.42
1:F:60:ASN:CB	1:F:61:PRO:HD3	2.46	0.42
1:D:282:GLN:HB3	2:D:1137:HOH:O	2.18	0.42
1:E:72:LEU:HD23	1:E:138:VAL:HG23	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:60:ASN:HB2	1:D:61:PRO:CD	2.33	0.42
1:F:31:PRO:HG2	1:F:325:TYR:OH	2.19	0.42
1:B:178:ASN:HD22	1:B:181:THR:H	1.66	0.42
1:C:111:VAL:O	1:C:115:VAL:HG23	2.20	0.42
1:C:79:VAL:C	1:C:81:GLY:H	2.23	0.42
1:E:202:PHE:HB2	1:E:329:ILE:HD12	2.02	0.42
1:B:23:GLN:O	1:B:163:GLN:HA	2.20	0.42
1:F:200:CYS:SG	1:F:302:LEU:HD22	2.60	0.42
1:F:147:SER:O	1:F:148:MSE:CB	2.68	0.42
1:C:147:SER:O	1:C:148:MSE:CB	2.68	0.42
1:B:279:GLU:HB3	1:B:282:GLN:HE21	1.85	0.42
1:A:34:MSE:HE1	1:A:110:VAL:N	2.34	0.42
1:A:112:VAL:HG22	1:A:135:THR:HG21	2.01	0.42
1:A:247:LYS:HB2	1:A:250:LYS:CB	2.49	0.42
1:D:31:PRO:HD2	1:D:325:TYR:CZ	2.55	0.42
1:C:23:GLN:O	1:C:163:GLN:HA	2.20	0.42
1:E:34:MSE:HE2	1:E:110:VAL:N	2.35	0.42
1:F:206:TRP:CZ2	1:F:208:LYS:HA	2.54	0.42
1:A:206:TRP:CZ3	1:A:208:LYS:HG2	2.55	0.42
1:C:206:TRP:HB2	1:C:324:LEU:HG	2.02	0.42
1:C:1001:HIS:N	2:C:1074:HOH:O	2.53	0.42
1:D:255:VAL:HB	1:D:287:VAL:HG21	2.02	0.41
1:B:289:LYS:NZ	2:B:1131:HOH:O	2.43	0.41
1:A:84:LEU:HD11	1:A:87:PHE:HE2	1.85	0.41
1:F:259:SER:O	1:F:263:GLU:HG3	2.20	0.41
1:C:212:VAL:O	1:C:212:VAL:HG13	2.21	0.41
1:E:68:GLU:OE1	1:E:142:ARG:HG3	2.19	0.41
1:A:270:TYR:HB3	2:A:1112:HOH:O	2.19	0.41
1:A:193:GLN:HE22	1:A:195:LYS:HE2	1.84	0.41
1:B:189:PHE:CD1	1:B:189:PHE:C	2.93	0.41
1:B:198:LEU:N	2:B:1103:HOH:O	2.51	0.41
1:B:198:LEU:HB2	1:B:246:GLY:HA2	2.01	0.41
1:E:85:ARG:NH2	2:E:1020:HOH:O	2.53	0.41
1:A:212:VAL:HG23	1:A:295:GLY:HA3	2.01	0.41
1:F:213:SER:CB	1:F:216:MSE:HE3	2.50	0.41
1:E:50:MSE:C	2:E:1098:HOH:O	2.58	0.41
1:A:198:LEU:HD11	1:A:246:GLY:C	2.39	0.41
1:F:91:ILE:HG22	1:F:96:GLU:OE1	2.20	0.41
1:F:146:GLY:O	1:F:167:ARG:HD3	2.20	0.41
1:C:212:VAL:HG22	1:C:216:MSE:CB	2.39	0.41
1:C:110:VAL:O	1:C:114:VAL:HG23	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:37:GLU:HA	1:D:37:GLU:OE1	2.20	0.41
1:D:194:VAL:O	1:D:196:PRO:HD3	2.20	0.41
1:F:139:LEU:CD1	1:F:151:LEU:HD11	2.50	0.41
1:D:168:GLN:HB3	2:D:1098:HOH:O	2.19	0.41
1:F:134:LEU:HD23	1:F:134:LEU:HA	1.92	0.41
1:A:50:MSE:HE1	1:D:61:PRO:HD2	2.02	0.41
1:E:34:MSE:HE1	1:E:110:VAL:HG22	2.02	0.41
1:D:296:GLY:O	1:D:297:ASP:HB2	2.21	0.41
1:A:41:SER:O	1:A:86:PRO:HB2	2.21	0.41
1:E:62:ASP:N	2:E:1027:HOH:O	2.48	0.41
1:D:139:LEU:CG	2:D:1103:HOH:O	2.68	0.41
1:B:34:MSE:CB	2:B:1069:HOH:O	2.69	0.41
1:C:313:LYS:NZ	1:C:326:GLN:NE2	2.69	0.41
1:A:273:LEU:HD11	1:C:44:TYR:CE1	2.55	0.41
1:A:146:GLY:O	1:A:167:ARG:CD	2.61	0.40
1:D:26:LEU:HD22	1:D:159:LEU:CD1	2.47	0.40
1:B:18:ILE:HG13	1:B:147:SER:HB2	2.02	0.40
1:E:252:ILE:O	1:E:256:GLU:HB2	2.21	0.40
1:A:100:LYS:HE2	2:A:1101:HOH:O	2.21	0.40
1:B:203:LEU:CD2	1:B:326:GLN:HG3	2.50	0.40
1:C:243:LEU:HD23	1:C:251:ILE:HD13	2.03	0.40
1:E:41:SER:O	1:E:86:PRO:HB2	2.21	0.40
1:C:221:LYS:HE2	1:C:221:LYS:HB3	1.86	0.40
1:C:34:MSE:HG2	1:C:95:MSE:CB	2.51	0.40
1:A:150:ASP:N	1:A:150:ASP:OD1	2.54	0.40
1:F:225:ASN:HD21	1:F:227:ASN:HB2	1.85	0.40
1:E:225:ASN:ND2	1:E:227:ASN:HD22	2.19	0.40
1:E:31:PRO:HD2	1:E:325:TYR:CZ	2.56	0.40
1:B:212:VAL:HG23	1:B:295:GLY:HA3	2.03	0.40
1:D:65:LEU:HD13	1:D:142:ARG:NH1	2.37	0.40
1:E:166:ASP:O	1:E:170:VAL:HG23	2.21	0.40
1:A:304:PHE:O	1:A:305:ASP:HB3	2.22	0.40
1:F:271:THR:HB	1:F:272:PRO:HD2	2.04	0.40
1:F:107:SER:O	1:F:111:VAL:HG23	2.21	0.40
1:F:84:LEU:HA	1:F:122:TYR:OH	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	318/337 (94%)	299 (94%)	18 (6%)	1 (0%)	46	62
1	B	318/337 (94%)	304 (96%)	11 (4%)	3 (1%)	21	29
1	C	318/337 (94%)	302 (95%)	11 (4%)	5 (2%)	12	15
1	D	318/337 (94%)	300 (94%)	15 (5%)	3 (1%)	21	29
1	E	318/337 (94%)	303 (95%)	10 (3%)	5 (2%)	12	15
1	F	318/337 (94%)	306 (96%)	11 (4%)	1 (0%)	46	62
All	All	1908/2022 (94%)	1814 (95%)	76 (4%)	18 (1%)	21	29

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	99	GLY
1	B	208	LYS
1	E	282	GLN
1	E	283	ASP
1	C	99	GLY
1	D	97	ARG
1	F	97	ARG
1	B	212	VAL
1	C	101	LYS
1	D	213	SER
1	C	266	SER
1	E	211	ALA
1	B	31	PRO
1	C	61	PRO
1	D	31	PRO
1	C	31	PRO
1	E	31	PRO
1	E	99	GLY

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	271/276 (98%)	261 (96%)	10 (4%)	41	61
1	B	271/276 (98%)	255 (94%)	16 (6%)	24	37
1	C	271/276 (98%)	264 (97%)	7 (3%)	54	74
1	D	271/276 (98%)	261 (96%)	10 (4%)	41	61
1	E	271/276 (98%)	263 (97%)	8 (3%)	48	69
1	F	271/276 (98%)	260 (96%)	11 (4%)	37	56
All	All	1626/1656 (98%)	1564 (96%)	62 (4%)	40	59

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	5	LYS
1	A	9	LYS
1	A	26	LEU
1	A	142	ARG
1	A	161	LEU
1	A	273	LEU
1	A	285	GLN
1	A	312	LEU
1	A	324	LEU
1	B	1	MSE
1	B	6	THR
1	B	9	LYS
1	B	26	LEU
1	B	34	MSE
1	B	96	GLU
1	B	106	SER
1	B	145	ASN
1	B	161	LEU
1	B	178	ASN
1	B	198	LEU
1	B	285	GLN

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Mol	Chain	Res	Type
1	B	305	ASP
1	B	307	GLN
1	B	312	LEU
1	B	324	LEU
1	C	26	LEU
1	C	34	MSE
1	C	96	GLU
1	C	110	VAL
1	C	128	GLN
1	C	161	LEU
1	C	198	LEU
1	D	9	LYS
1	D	26	LEU
1	D	34	MSE
1	D	72	LEU
1	D	80	ARG
1	D	96	GLU
1	D	106	SER
1	D	128	GLN
1	D	198	LEU
1	D	312	LEU
1	E	9	LYS
1	E	26	LEU
1	E	125	SER
1	E	145	ASN
1	E	161	LEU
1	E	178	ASN
1	E	198	LEU
1	E	312	LEU
1	F	1	MSE
1	F	6	THR
1	F	9	LYS
1	F	20	GLU
1	F	26	LEU
1	F	60	ASN
1	F	101	LYS
1	F	161	LEU
1	F	198	LEU
1	F	200	CYS
1	F	312	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	GLN
1	A	193	GLN
1	A	218	GLN
1	A	226	GLN
1	A	254	GLN
1	A	282	GLN
1	A	285	GLN
1	A	326	GLN
1	B	128	GLN
1	B	178	ASN
1	B	282	GLN
1	B	285	GLN
1	B	326	GLN
1	C	128	GLN
1	C	223	ASN
1	C	225	ASN
1	C	326	GLN
1	D	60	ASN
1	D	67	GLN
1	D	193	GLN
1	D	225	ASN
1	D	282	GLN
1	E	128	GLN
1	E	193	GLN
1	E	226	GLN
1	E	227	ASN
1	E	285	GLN
1	E	326	GLN
1	F	193	GLN
1	F	219	GLN
1	F	225	ASN
1	F	326	GLN

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 [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, 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	314/337 (93%)	0.13	18 (5%)	27 27	21, 38, 61, 75	0
1	B	314/337 (93%)	0.09	17 (5%)	29 29	21, 37, 59, 75	0
1	C	314/337 (93%)	0.11	17 (5%)	29 29	22, 38, 58, 75	0
1	D	314/337 (93%)	0.24	19 (6%)	25 24	20, 41, 64, 73	0
1	E	314/337 (93%)	0.15	12 (3%)	44 44	20, 35, 59, 77	0
1	F	314/337 (93%)	0.12	13 (4%)	41 41	22, 36, 60, 78	0
All	All	1884/2022 (93%)	0.14	96 (5%)	32 31	20, 37, 61, 78	0

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	99	GLY	10.4
1	E	98	GLU	6.6
1	F	99	GLY	6.4
1	F	97	ARG	6.3
1	F	98	GLU	5.7
1	D	61	PRO	5.7
1	D	60	ASN	5.5
1	C	61	PRO	5.5
1	F	60	ASN	5.4
1	D	99	GLY	5.3
1	B	97	ARG	5.3
1	B	99	GLY	5.2
1	B	100	LYS	5.2
1	D	283	ASP	5.2
1	B	60	ASN	5.1
1	C	60	ASN	5.0
1	E	61	PRO	4.9
1	D	97	ARG	4.8
1	A	60	ASN	4.8

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Mol	Chain	Res	Type	RSRZ
1	A	211	ALA	4.7
1	B	98	GLU	4.6
1	C	98	GLU	4.6
1	A	98	GLU	4.5
1	D	212	VAL	4.5
1	A	99	GLY	4.4
1	D	98	GLU	4.3
1	A	218	GLN	4.2
1	E	214	SER	4.0
1	A	214	SER	3.9
1	A	213	SER	3.9
1	F	100	LYS	3.9
1	C	218	GLN	3.9
1	C	214	SER	3.8
1	C	210	VAL	3.7
1	E	97	ARG	3.7
1	C	99	GLY	3.6
1	C	100	LYS	3.6
1	A	212	VAL	3.5
1	A	100	LYS	3.4
1	D	218	GLN	3.3
1	D	48	SER	3.3
1	E	60	ASN	3.3
1	D	214	SER	3.3
1	B	214	SER	3.2
1	A	61	PRO	3.2
1	F	214	SER	3.1
1	C	282	GLN	3.1
1	E	100	LYS	3.1
1	F	145	ASN	3.0
1	F	199	GLU	3.0
1	D	282	GLN	3.0
1	D	143	GLY	3.0
1	C	97	ARG	3.0
1	B	282	GLN	3.0
1	D	213	SER	3.0
1	F	101	LYS	3.0
1	B	218	GLN	2.9
1	E	213	SER	2.9
1	A	283	ASP	2.8
1	D	211	ALA	2.8
1	B	61	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	49	ASP	2.8
1	C	200	CYS	2.8
1	A	97	ARG	2.7
1	B	212	VAL	2.7
1	D	145	ASN	2.6
1	B	283	ASP	2.5
1	C	222	GLN	2.5
1	C	249	GLU	2.5
1	B	158	ASP	2.5
1	F	186	ASP	2.4
1	C	301	ALA	2.4
1	E	222	GLN	2.4
1	A	287	VAL	2.4
1	C	49	ASP	2.3
1	F	213	SER	2.3
1	A	199	GLU	2.3
1	A	282	GLN	2.3
1	E	200	CYS	2.3
1	E	184	GLU	2.3
1	B	222	GLN	2.3
1	C	211	ALA	2.2
1	B	199	GLU	2.2
1	D	49	ASP	2.2
1	C	199	GLU	2.1
1	F	222	GLN	2.1
1	A	145	ASN	2.1
1	B	226	GLN	2.1
1	B	94	LYS	2.1
1	A	226	GLN	2.1
1	D	184	GLU	2.1
1	D	168	GLN	2.1
1	D	177	GLU	2.0
1	E	89	LEU	2.0
1	B	213	SER	2.0
1	F	209	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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.