



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:21 PM GMT

PDB ID : 3TPF  
Title : Crystal structure of anabolic ornithine carbamoyltransferase from *Campylobacter jejuni* subsp. *jejuni* NCTC 11168  
Authors : Shabalin, I.G.; Onopriyenko, O.; Grimshaw, S.; Porebski, P.J.; Grabowski, M.; Savchenko, A.; Chruszcz, M.; Anderson, W.F.; Minor, W.; Center for Structural Genomics of Infectious Diseases (CSGID)  
Deposited on : 2011-09-07  
Resolution : 2.70 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  
<http://wwpdb.org/validation/2016/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  
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

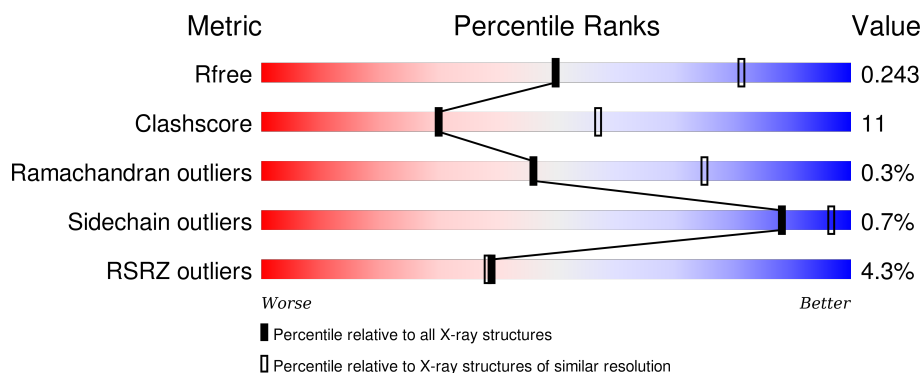
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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. 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	307	<div> <div>4%</div> <div>80%18%..</div> </div>
1	B	307	<div> <div>2%</div> <div>78%20%..</div> </div>
1	C	307	<div> <div>7%</div> <div>79%18%..</div> </div>
1	D	307	<div> <div>5%</div> <div>80%17%..</div> </div>
1	E	307	<div> <div>5%</div> <div>79%19%.</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	307	<div><div></div><div>2%</div><div>82%</div><div>14%</div><div></div></div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13981 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 Ornithine carbamoyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	303	Total	C	N	O	S	Se	0	0	0
			2334	1503	385	430	3	13			
1	B	305	Total	C	N	O	S	Se	0	0	0
			2358	1523	383	437	3	12			
1	C	301	Total	C	N	O	S	Se	0	0	0
			2284	1469	380	420	3	12			
1	D	303	Total	C	N	O	S	Se	0	0	0
			2305	1481	376	432	3	13			
1	E	300	Total	C	N	O	S	Se	0	0	0
			2281	1470	371	425	3	12			
1	F	298	Total	C	N	O	S	Se	0	0	0
			2294	1481	375	423	3	12			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	EXPRESSION TAG	UNP Q9PNU6
B	0	GLY	-	EXPRESSION TAG	UNP Q9PNU6
C	0	GLY	-	EXPRESSION TAG	UNP Q9PNU6
D	0	GLY	-	EXPRESSION TAG	UNP Q9PNU6
E	0	GLY	-	EXPRESSION TAG	UNP Q9PNU6
F	0	GLY	-	EXPRESSION TAG	UNP Q9PNU6

- Molecule 2 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			7	4	3		
2	B	1	Total	C	O	0	0
			7	4	3		
2	F	1	Total	C	O	0	0
			7	4	3		

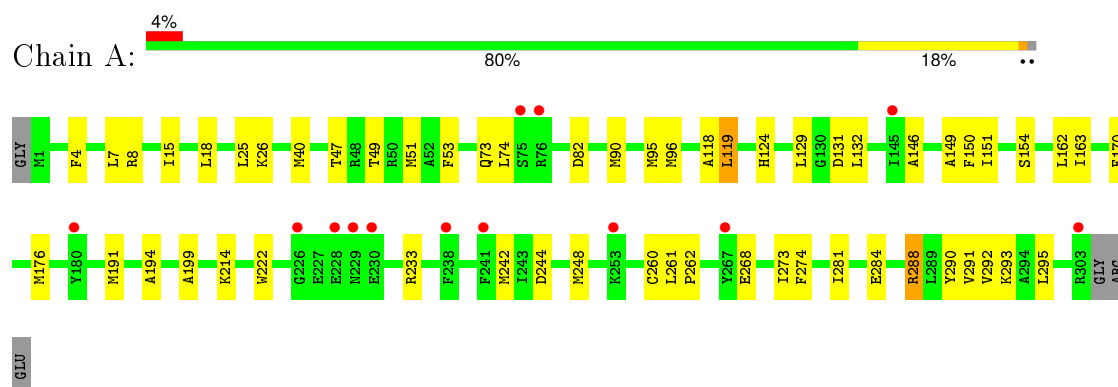
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	15	Total	O	0	0
			15	15		
3	B	32	Total	O	0	0
			32	32		
3	C	16	Total	O	0	0
			16	16		
3	D	13	Total	O	0	0
			13	13		
3	E	14	Total	O	0	0
			14	14		
3	F	14	Total	O	0	0
			14	14		

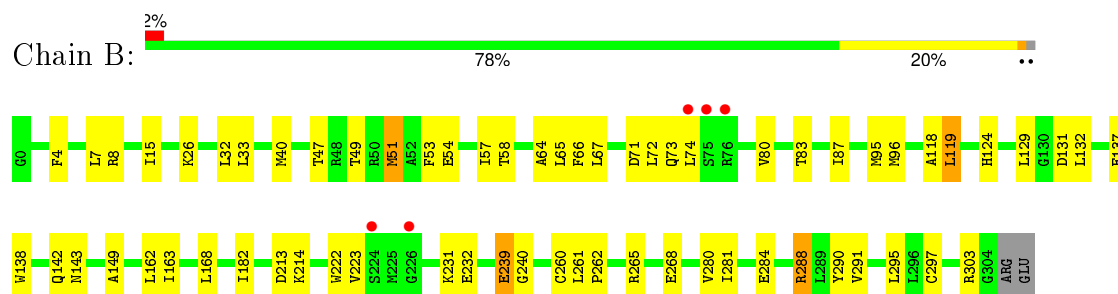
### 3 Residue-property plots [i](#)

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.

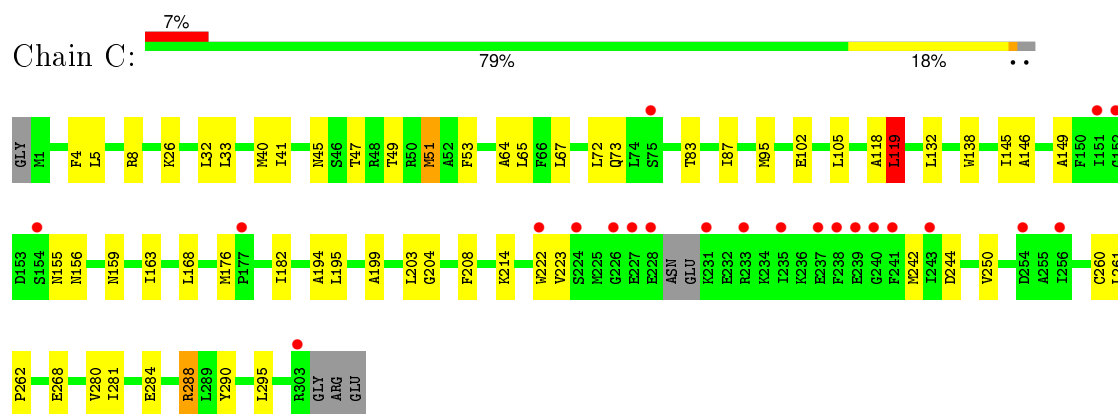
- Molecule 1: Ornithine carbamoyltransferase



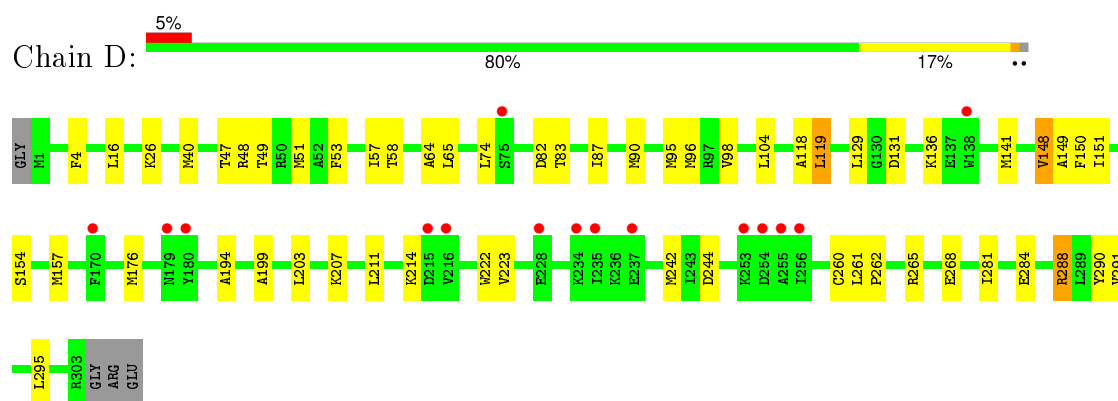
- Molecule 1: Ornithine carbamoyltransferase



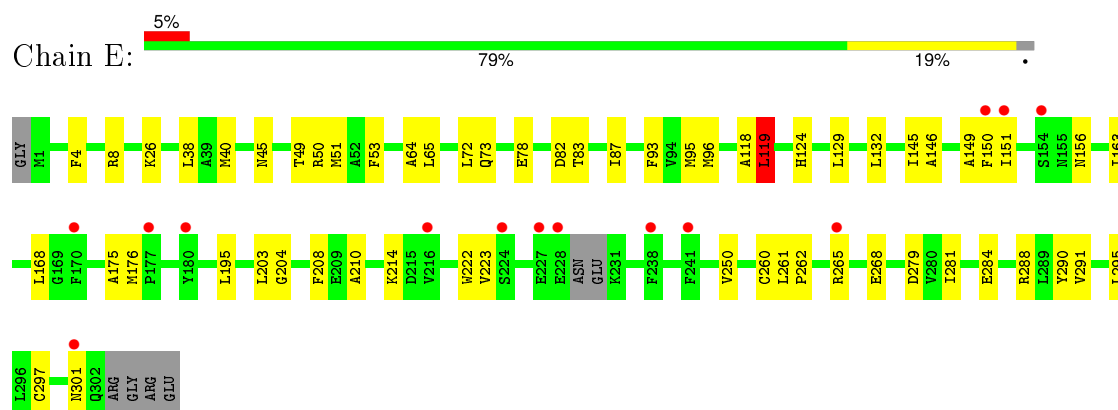
- Molecule 1: Ornithine carbamoyltransferase



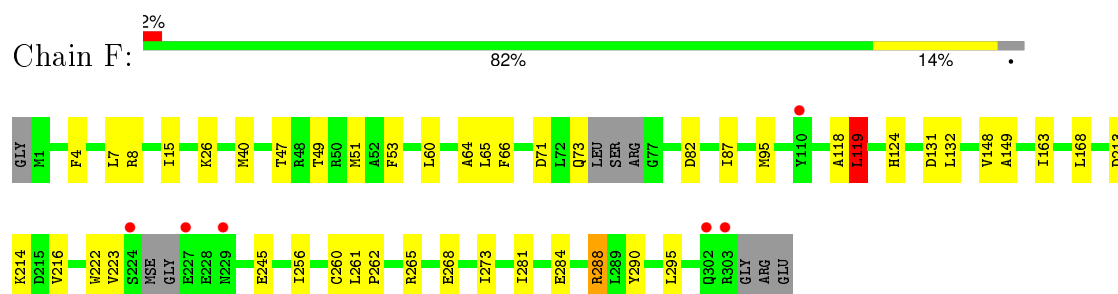
- Molecule 1: Ornithine carbamoyltransferase



- Molecule 1: Ornithine carbamoyltransferase



- Molecule 1: Ornithine carbamoyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 <sub>1</sub> 2 <sub>1</sub> 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	157.16 Å   87.40 Å   141.56 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	50.00 – 2.70 42.83 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.00-2.70) 99.8 (42.83-2.70)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.99 (at 2.69 Å)	Xtriage
Refinement program	REFMAC 5.6.0117, COOT	Depositor
R, $R_{free}$	0.202   ,   0.240 0.204   ,   0.243	Depositor DCC
$R_{free}$ test set	2758 reflections (5.36%)	DCC
Wilson B-factor (Å <sup>2</sup> )	59.8	Xtriage
Anisotropy	0.771	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 59.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	5 of 54248 reflections (0.009%)	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	13981	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.40 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.3286e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.86	0/2366	0.95	4/3182 (0.1%)
1	B	1.00	3/2389 (0.1%)	1.01	2/3206 (0.1%)
1	C	0.81	2/2312 (0.1%)	0.92	3/3108 (0.1%)
1	D	0.84	0/2337	0.93	4/3149 (0.1%)
1	E	0.77	0/2311	0.88	1/3112 (0.0%)
1	F	0.88	0/2325	0.93	0/3125
All	All	0.86	5/14040 (0.0%)	0.94	14/18882 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	54	GLU	CD-OE1	5.77	1.31	1.25
1	C	102	GLU	CD-OE1	5.42	1.31	1.25
1	C	102	GLU	CG-CD	5.18	1.59	1.51
1	B	137	GLU	CD-OE2	5.13	1.31	1.25
1	B	54	GLU	CD-OE2	5.06	1.31	1.25

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	96	MSE	CG-SE-CE	5.85	111.77	98.90
1	A	233	ARG	NE-CZ-NH2	5.80	123.20	120.30
1	C	51	MSE	CG-SE-CE	-5.76	86.23	98.90
1	A	162	LEU	CB-CG-CD2	-5.73	101.25	111.00
1	A	132	LEU	CB-CG-CD2	-5.64	101.41	111.00
1	A	90	MSE	CA-CB-CG	-5.63	103.73	113.30
1	D	65	LEU	CB-CG-CD1	5.51	120.36	111.00
1	D	157	MSE	CG-SE-CE	-5.48	86.84	98.90
1	B	232	GLU	CA-CB-CG	5.47	125.43	113.40
1	C	67	LEU	CB-CG-CD2	5.38	120.15	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	90	MSE	CA-CB-CG	-5.12	104.59	113.30
1	B	51	MSE	CG-SE-CE	5.08	110.07	98.90
1	D	148	VAL	CB-CA-C	-5.07	101.78	111.40
1	C	119	LEU	CB-CG-CD1	5.00	119.51	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	2334	0	2270	49	0
1	B	2358	0	2307	59	0
1	C	2284	0	2219	57	0
1	D	2305	0	2188	52	0
1	E	2281	0	2182	59	0
1	F	2294	0	2223	47	0
2	A	7	0	10	0	0
2	B	7	0	10	0	0
2	F	7	0	10	0	0
3	A	15	0	0	0	0
3	B	32	0	0	2	0
3	C	16	0	0	0	0
3	D	13	0	0	1	0
3	E	14	0	0	1	0
3	F	14	0	0	0	0
All	All	13981	0	13419	303	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 11.

All (303) 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:A:248:MSE:CE	1:A:248:MSE:SE	2.15	1.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:53:PHE:CD2	1:D:95:MSE:CE	2.31	1.13
1:D:53:PHE:CD2	1:D:95:MSE:HE2	1.91	1.03
1:E:53:PHE:CD2	1:E:95:MSE:CE	2.42	1.03
1:F:53:PHE:CD2	1:F:95:MSE:HE2	1.94	1.02
1:B:53:PHE:CD2	1:B:95:MSE:CE	2.50	0.94
1:A:53:PHE:CD2	1:A:95:MSE:CE	2.50	0.93
1:D:53:PHE:CG	1:D:95:MSE:CE	2.52	0.92
1:C:203:LEU:HD12	1:C:204:GLY:N	1.86	0.91
1:C:53:PHE:CD2	1:C:95:MSE:CE	2.52	0.91
1:F:65:LEU:HD12	1:F:65:LEU:N	1.85	0.90
1:D:47:THR:HG21	1:E:72:LEU:HD13	1.54	0.89
1:C:83:THR:HG22	1:C:87:ILE:HD12	1.57	0.86
1:E:53:PHE:CG	1:E:95:MSE:HE3	2.11	0.86
1:B:53:PHE:CD2	1:B:95:MSE:HE2	2.12	0.83
1:D:53:PHE:CE2	1:D:95:MSE:HE2	2.13	0.82
1:D:53:PHE:CG	1:D:95:MSE:HE2	2.14	0.81
1:F:53:PHE:CG	1:F:95:MSE:HE2	2.15	0.81
1:F:53:PHE:CD2	1:F:95:MSE:CE	2.64	0.81
1:C:73:GLN:O	1:C:73:GLN:HG2	1.80	0.81
1:B:223:VAL:HG21	1:B:231:LYS:HG3	1.62	0.80
1:E:53:PHE:CD2	1:E:95:MSE:HE3	2.15	0.80
1:F:4:PHE:CG	1:F:295:LEU:HD11	2.17	0.80
1:A:53:PHE:CG	1:A:95:MSE:CE	2.68	0.77
1:E:53:PHE:CE2	1:E:95:MSE:HE2	2.20	0.77
1:C:203:LEU:HD12	1:C:204:GLY:H	1.48	0.77
1:C:53:PHE:CG	1:C:95:MSE:CE	2.68	0.77
1:B:53:PHE:CG	1:B:95:MSE:CE	2.67	0.76
1:D:53:PHE:CG	1:D:95:MSE:HE3	2.19	0.76
1:C:65:LEU:N	1:C:65:LEU:HD12	2.00	0.76
1:C:53:PHE:CD2	1:C:95:MSE:HE2	2.22	0.75
1:A:53:PHE:CD2	1:A:95:MSE:HE2	2.21	0.73
1:C:242:MSE:HE3	1:C:244:ASP:HB3	1.69	0.73
1:B:80:VAL:HG23	3:B:331:HOH:O	1.88	0.72
1:F:64:ALA:C	1:F:65:LEU:HD12	2.11	0.71
1:D:83:THR:HG22	1:D:87:ILE:HD12	1.72	0.71
1:B:53:PHE:CE2	1:B:95:MSE:HE2	2.26	0.71
1:D:40:MSE:HE2	1:D:64:ALA:HB1	1.73	0.70
1:A:4:PHE:CG	1:A:295:LEU:HD11	2.26	0.70
1:A:51:MSE:HE1	1:C:87:ILE:HG12	1.75	0.68
1:A:4:PHE:CD1	1:A:295:LEU:HD11	2.29	0.68
1:C:132:LEU:HD21	1:C:168:LEU:HD11	1.73	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:239:GLU:HG3	1:B:240:GLY:H	1.57	0.68
1:C:40:MSE:HA	1:C:95:MSE:HG3	1.76	0.68
1:F:4:PHE:CD1	1:F:295:LEU:HD11	2.29	0.68
1:D:53:PHE:CD2	1:D:95:MSE:HE1	2.24	0.67
1:E:53:PHE:CE2	1:E:95:MSE:CE	2.77	0.67
1:B:40:MSE:HA	1:B:95:MSE:HG3	1.76	0.67
1:A:129:LEU:HB3	1:A:291:VAL:HG11	1.75	0.67
1:B:53:PHE:CG	1:B:95:MSE:HE3	2.29	0.67
1:E:4:PHE:CG	1:E:295:LEU:HD11	2.29	0.67
1:C:4:PHE:CD1	1:C:295:LEU:HD11	2.31	0.66
1:C:83:THR:CG2	1:C:87:ILE:HD12	2.26	0.66
1:F:53:PHE:CG	1:F:95:MSE:CE	2.78	0.65
1:A:53:PHE:CG	1:A:95:MSE:HE3	2.31	0.65
1:D:4:PHE:CD1	1:D:295:LEU:HD11	2.32	0.65
1:E:53:PHE:CD2	1:E:95:MSE:HE2	2.27	0.65
1:E:53:PHE:CG	1:E:95:MSE:CE	2.76	0.65
1:A:40:MSE:HA	1:A:95:MSE:HG3	1.78	0.65
1:B:239:GLU:HG3	1:B:240:GLY:N	2.12	0.65
1:C:41:ILE:HG23	1:C:72:LEU:HD23	1.79	0.65
1:E:83:THR:HG22	1:E:87:ILE:HD12	1.78	0.64
1:E:279:ASP:HB2	3:E:315:HOH:O	1.96	0.64
1:E:45:ASN:HD22	1:F:71:ASP:HA	1.62	0.64
1:E:4:PHE:CD1	1:E:295:LEU:HD11	2.32	0.64
1:A:47:THR:HG21	1:C:72:LEU:HD13	1.80	0.64
1:E:149:ALA:HB2	1:E:214:LYS:HG3	1.80	0.64
1:C:53:PHE:CD2	1:C:95:MSE:HE1	2.34	0.63
1:F:65:LEU:CD1	1:F:65:LEU:N	2.59	0.63
1:B:53:PHE:CG	1:B:95:MSE:HE2	2.33	0.63
1:B:132:LEU:HD21	1:B:168:LEU:HD11	1.80	0.63
1:A:53:PHE:CD2	1:A:95:MSE:HE1	2.33	0.63
1:C:53:PHE:CG	1:C:95:MSE:HE3	2.34	0.63
1:D:150:PHE:C	1:D:151:ILE:HD12	2.20	0.62
1:E:203:LEU:HD12	1:E:204:GLY:N	2.15	0.62
1:C:64:ALA:C	1:C:65:LEU:HD12	2.19	0.61
1:B:87:ILE:HG12	1:C:51:MSE:HE1	1.82	0.61
1:D:4:PHE:CG	1:D:295:LEU:HD11	2.35	0.61
1:C:40:MSE:HG2	1:C:95:MSE:HG2	1.82	0.60
1:B:72:LEU:HD13	1:C:47:THR:HG21	1.82	0.60
1:C:149:ALA:HB2	1:C:214:LYS:HG3	1.83	0.60
1:B:32:LEU:HD12	1:B:297:CYS:SG	2.42	0.60
1:E:132:LEU:HD21	1:E:168:LEU:HD11	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:118:ALA:O	1:F:119:LEU:HB2	2.02	0.60
1:D:53:PHE:CB	1:D:95:MSE:HE3	2.32	0.60
1:C:4:PHE:CG	1:C:295:LEU:HD11	2.37	0.60
1:B:40:MSE:HG2	1:B:95:MSE:HG2	1.84	0.59
1:B:66:PHE:C	1:B:67:LEU:HD23	2.23	0.59
1:B:83:THR:HG22	1:B:87:ILE:HD12	1.85	0.58
1:B:57:ILE:HG23	1:B:58:THR:N	2.18	0.58
1:C:53:PHE:CE2	1:C:95:MSE:HE2	2.38	0.58
1:D:40:MSE:HA	1:D:95:MSE:HG3	1.85	0.58
1:B:4:PHE:CD1	1:B:295:LEU:HD11	2.38	0.58
1:C:53:PHE:CG	1:C:95:MSE:HE2	2.38	0.58
1:A:51:MSE:CE	1:C:87:ILE:HG12	2.33	0.58
1:D:83:THR:CG2	1:D:87:ILE:HD12	2.33	0.58
1:F:47:THR:HG22	1:F:51:MSE:HE2	1.85	0.58
1:A:40:MSE:HG2	1:A:95:MSE:HG2	1.87	0.57
1:A:53:PHE:CE2	1:A:95:MSE:HE2	2.38	0.57
1:C:118:ALA:O	1:C:119:LEU:HB2	2.05	0.57
1:D:51:MSE:CE	1:E:87:ILE:HG12	2.35	0.57
1:F:149:ALA:HB2	1:F:214:LYS:HG3	1.86	0.57
1:E:65:LEU:N	1:E:65:LEU:HD12	2.20	0.57
1:E:50:ARG:NH1	1:F:66:PHE:O	2.37	0.57
1:A:53:PHE:CG	1:A:95:MSE:HE2	2.39	0.57
1:D:154:SER:HA	1:D:176:MSE:HE1	1.85	0.57
1:E:118:ALA:O	1:E:119:LEU:HB2	2.05	0.57
1:E:40:MSE:HA	1:E:95:MSE:HG3	1.87	0.56
1:B:87:ILE:HG12	1:C:51:MSE:CE	2.36	0.56
1:B:118:ALA:O	1:B:119:LEU:HB2	2.05	0.56
1:F:40:MSE:HA	1:F:95:MSE:HG3	1.88	0.56
1:E:51:MSE:CE	1:F:87:ILE:HG12	2.35	0.56
1:D:82:ASP:OD2	1:F:265:ARG:NH2	2.38	0.56
1:F:8:ARG:HG3	1:F:163:ILE:HD13	1.87	0.56
1:A:129:LEU:HB3	1:A:291:VAL:CG1	2.36	0.55
1:A:7:LEU:HD22	1:A:15:ILE:HG12	1.88	0.55
1:C:182:ILE:HD12	1:C:203:LEU:HD22	1.89	0.55
1:F:49:THR:HG22	1:F:95:MSE:HE1	1.88	0.55
1:F:53:PHE:CE2	1:F:95:MSE:HE2	2.40	0.55
1:B:260:CYS:O	1:B:261:LEU:HB2	2.07	0.54
1:A:118:ALA:O	1:A:119:LEU:HB2	2.08	0.54
1:E:45:ASN:ND2	1:F:71:ASP:OD1	2.40	0.54
1:E:119:LEU:HD12	1:E:124:HIS:CD2	2.42	0.54
1:B:4:PHE:CG	1:B:295:LEU:HD11	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:119:LEU:HD11	1:E:156:ASN:HD21	1.73	0.54
1:C:32:LEU:O	1:C:33:LEU:HD23	2.08	0.54
1:C:8:ARG:HG3	1:C:163:ILE:HD13	1.90	0.54
1:C:65:LEU:N	1:C:65:LEU:CD1	2.71	0.54
1:C:41:ILE:CG2	1:C:72:LEU:HD23	2.38	0.54
1:F:222:TRP:CZ2	1:F:268:GLU:HA	2.43	0.53
1:E:261:LEU:N	1:E:261:LEU:HD23	2.22	0.53
1:B:138:TRP:CE2	1:B:280:VAL:HG23	2.43	0.53
1:B:65:LEU:HD12	1:B:65:LEU:N	2.23	0.53
1:B:149:ALA:HB2	1:B:214:LYS:HG3	1.90	0.53
1:C:222:TRP:CZ2	1:C:268:GLU:HA	2.43	0.53
1:A:74:LEU:HD11	1:A:96:MSE:CE	2.38	0.53
1:D:149:ALA:HB2	1:D:214:LYS:HG3	1.89	0.53
1:D:98:VAL:HG21	1:D:104:LEU:HG	1.91	0.53
1:D:40:MSE:HG2	1:D:95:MSE:HG2	1.91	0.53
1:A:154:SER:HA	1:A:176:MSE:HE1	1.89	0.52
1:E:203:LEU:HD12	1:E:204:GLY:H	1.73	0.52
1:F:7:LEU:HD22	1:F:15:ILE:HG12	1.91	0.52
1:E:119:LEU:HD12	1:E:124:HIS:CG	2.45	0.52
1:B:71:ASP:HA	1:C:45:ASN:HD22	1.75	0.51
1:E:222:TRP:CZ2	1:E:268:GLU:HA	2.46	0.51
1:C:260:CYS:O	1:C:261:LEU:HB2	2.11	0.51
1:B:8:ARG:HG3	1:B:163:ILE:HD13	1.92	0.51
1:A:129:LEU:CB	1:A:291:VAL:HG11	2.40	0.50
1:A:73:GLN:HG3	1:A:73:GLN:O	2.11	0.50
1:D:222:TRP:CZ2	1:D:268:GLU:HA	2.45	0.50
1:A:150:PHE:C	1:A:151:ILE:HD12	2.32	0.50
1:F:73:GLN:HG3	1:F:73:GLN:O	2.11	0.50
1:E:260:CYS:O	1:E:261:LEU:HB2	2.12	0.50
1:A:149:ALA:HB2	1:A:214:LYS:HG3	1.94	0.50
1:D:118:ALA:O	1:D:119:LEU:HB2	2.12	0.50
1:B:83:THR:CG2	1:B:87:ILE:HD12	2.42	0.50
1:A:18:LEU:HD13	1:A:129:LEU:HD11	1.93	0.50
1:D:223:VAL:HG12	1:D:223:VAL:O	2.12	0.50
1:C:208:PHE:HD2	1:C:250:VAL:HG21	1.77	0.50
1:C:138:TRP:CE2	1:C:280:VAL:HG23	2.48	0.49
1:B:47:THR:HG22	1:B:51:MSE:CE	2.41	0.49
1:E:73:GLN:O	1:E:73:GLN:HG3	2.12	0.49
1:F:119:LEU:HD12	1:F:124:HIS:CD2	2.47	0.49
1:D:26:LYS:HD2	1:D:290:TYR:CE2	2.47	0.49
1:B:67:LEU:HD23	1:B:67:LEU:N	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:40:MSE:HG2	1:E:95:MSE:HG2	1.95	0.49
1:F:40:MSE:HG2	1:F:95:MSE:HG2	1.94	0.49
1:A:191:MSE:HA	1:A:191:MSE:HE3	1.94	0.49
1:C:145:ILE:CG2	1:C:146:ALA:N	2.74	0.49
1:E:150:PHE:C	1:E:151:ILE:HD12	2.34	0.48
1:A:194:ALA:CB	1:A:199:ALA:HB3	2.43	0.48
1:F:260:CYS:O	1:F:261:LEU:HB2	2.13	0.48
1:D:207:LYS:O	1:D:211:LEU:HD12	2.13	0.48
1:A:260:CYS:O	1:A:261:LEU:HB2	2.14	0.48
1:F:119:LEU:HD12	1:F:124:HIS:CG	2.49	0.47
1:B:26:LYS:HD2	1:B:290:TYR:CE2	2.49	0.47
1:D:260:CYS:O	1:D:261:LEU:HB2	2.14	0.47
1:E:8:ARG:HG3	1:E:163:ILE:HD13	1.95	0.47
1:C:118:ALA:O	1:C:119:LEU:CB	2.62	0.47
1:B:303:ARG:HB2	3:B:327:HOH:O	2.14	0.47
1:B:53:PHE:CD2	1:B:95:MSE:HE1	2.45	0.47
1:D:129:LEU:HB3	1:D:291:VAL:HG11	1.95	0.47
1:C:41:ILE:HG23	1:C:72:LEU:CD2	2.43	0.47
1:B:213:ASP:O	1:B:213:ASP:CG	2.53	0.47
1:C:223:VAL:O	1:C:223:VAL:HG12	2.13	0.47
1:D:51:MSE:HB2	1:D:51:MSE:HE2	1.82	0.47
1:A:119:LEU:HD12	1:A:124:HIS:CD2	2.49	0.47
1:A:222:TRP:CZ2	1:A:268:GLU:HA	2.49	0.47
1:B:222:TRP:CZ2	1:B:268:GLU:HA	2.49	0.47
1:B:261:LEU:HB3	1:B:262:PRO:HA	1.97	0.47
1:E:129:LEU:HB3	1:E:291:VAL:HG11	1.97	0.46
1:B:7:LEU:HD22	1:B:15:ILE:HG12	1.97	0.46
1:A:26:LYS:HD2	1:A:290:TYR:CE2	2.50	0.46
1:B:53:PHE:CB	1:B:95:MSE:HE3	2.45	0.46
1:B:118:ALA:O	1:B:119:LEU:CB	2.63	0.46
1:F:132:LEU:HD21	1:F:168:LEU:HD11	1.97	0.46
1:F:118:ALA:O	1:F:119:LEU:CB	2.64	0.46
1:B:162:LEU:HD22	1:B:182:ILE:HD13	1.98	0.46
1:C:95:MSE:HG3	1:C:95:MSE:O	2.16	0.46
1:D:49:THR:OG1	1:D:288:ARG:NH1	2.49	0.46
1:D:136:LYS:CA	1:D:141:MSE:HE3	2.47	0.45
1:F:4:PHE:CD1	1:F:295:LEU:CD1	2.98	0.45
1:A:273:ILE:CG2	1:A:274:PHE:N	2.78	0.45
1:B:74:LEU:HD11	1:B:96:MSE:CE	2.46	0.45
1:C:53:PHE:CB	1:C:95:MSE:HE3	2.47	0.45
1:D:242:MSE:HE3	1:D:244:ASP:HB3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:48:ARG:NH1	1:E:78:GLU:OE1	2.48	0.45
1:F:40:MSE:HE2	1:F:64:ALA:HB1	1.99	0.45
1:E:51:MSE:HE1	1:F:87:ILE:HG12	1.98	0.45
1:F:148:VAL:HG22	1:F:216:VAL:HB	1.99	0.45
1:D:87:ILE:HG12	1:F:51:MSE:CE	2.47	0.45
1:B:64:ALA:C	1:B:65:LEU:HD12	2.37	0.45
1:D:57:ILE:HG23	1:D:58:THR:N	2.31	0.45
1:D:48:ARG:HB3	3:D:318:HOH:O	2.16	0.45
1:D:281:ILE:O	1:D:284:GLU:HB3	2.17	0.45
1:B:129:LEU:HB3	1:B:291:VAL:CG1	2.47	0.45
1:E:26:LYS:HD2	1:E:290:TYR:CE2	2.51	0.45
1:B:57:ILE:CG2	1:B:58:THR:N	2.79	0.45
1:A:281:ILE:O	1:A:284:GLU:HB3	2.17	0.45
1:E:49:THR:CG2	1:E:95:MSE:HE1	2.47	0.44
1:B:142:GLN:O	1:B:143:ASN:CB	2.66	0.44
1:D:53:PHE:HB2	1:D:95:MSE:HE3	1.99	0.44
1:B:129:LEU:HB3	1:B:291:VAL:HG11	1.98	0.44
1:E:38:LEU:CD1	1:E:93:PHE:HB2	2.48	0.44
1:D:129:LEU:HB3	1:D:291:VAL:CG1	2.48	0.44
1:C:26:LYS:HD2	1:C:290:TYR:CE2	2.52	0.44
1:F:213:ASP:O	1:F:213:ASP:CG	2.56	0.44
1:F:131:ASP:OD1	1:F:284:GLU:OE2	2.35	0.44
1:A:25:LEU:HD13	1:A:293:LYS:HB3	2.01	0.43
1:D:154:SER:CA	1:D:176:MSE:HE1	2.48	0.43
1:E:129:LEU:HB3	1:E:291:VAL:CG1	2.47	0.43
1:A:8:ARG:HG3	1:A:163:ILE:HD13	1.99	0.43
1:C:155:ASN:O	1:C:159:ASN:ND2	2.52	0.43
1:B:131:ASP:OD1	1:B:284:GLU:OE2	2.36	0.43
1:E:118:ALA:O	1:E:119:LEU:CB	2.66	0.43
1:A:273:ILE:HD12	1:A:273:ILE:HA	1.89	0.43
1:B:49:THR:OG1	1:B:288:ARG:NH1	2.50	0.43
1:A:53:PHE:CB	1:A:95:MSE:HE3	2.49	0.43
1:B:66:PHE:O	1:B:67:LEU:HD23	2.18	0.43
1:C:5:LEU:HD13	1:C:105:LEU:HD23	2.01	0.43
1:A:146:ALA:HB3	1:A:170:PHE:CE2	2.54	0.43
1:E:53:PHE:CD2	1:E:95:MSE:HE1	2.45	0.43
1:C:156:ASN:HA	1:C:159:ASN:HD22	1.83	0.42
1:C:194:ALA:CB	1:C:199:ALA:HB3	2.49	0.42
1:B:47:THR:HG22	1:B:51:MSE:HE2	2.02	0.42
1:B:239:GLU:CG	1:B:240:GLY:N	2.81	0.42
1:B:32:LEU:O	1:B:33:LEU:HD23	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:64:ALA:C	1:E:65:LEU:HD12	2.40	0.42
1:D:131:ASP:OD1	1:D:284:GLU:OE2	2.38	0.42
1:E:261:LEU:HB3	1:E:262:PRO:HA	2.01	0.42
1:C:261:LEU:HB3	1:C:262:PRO:HA	2.01	0.42
1:F:26:LYS:HD2	1:F:290:TYR:CE2	2.55	0.42
1:E:53:PHE:CZ	1:E:95:MSE:HE2	2.53	0.42
1:A:261:LEU:HB3	1:A:262:PRO:HA	2.02	0.42
1:F:261:LEU:HB3	1:F:262:PRO:HA	2.02	0.42
1:D:261:LEU:HB3	1:D:262:PRO:HA	2.02	0.42
1:A:118:ALA:O	1:A:119:LEU:CB	2.67	0.42
1:A:131:ASP:OD1	1:A:284:GLU:OE2	2.38	0.42
1:E:208:PHE:HD2	1:E:250:VAL:HG21	1.84	0.42
1:C:176:MSE:HE2	1:C:176:MSE:HB3	1.93	0.42
1:D:148:VAL:CG1	1:D:149:ALA:N	2.80	0.42
1:E:195:LEU:HD23	1:E:195:LEU:HA	1.80	0.42
1:A:242:MSE:HE3	1:A:244:ASP:HB3	2.01	0.42
1:D:136:LYS:CB	1:D:141:MSE:HE3	2.49	0.41
1:E:281:ILE:O	1:E:284:GLU:HB3	2.19	0.41
1:E:176:MSE:HB3	1:E:176:MSE:HE2	1.90	0.41
1:A:49:THR:OG1	1:A:288:ARG:NH1	2.50	0.41
1:A:4:PHE:CD1	1:A:295:LEU:CD1	3.02	0.41
1:D:151:ILE:HD12	1:D:151:ILE:N	2.35	0.41
1:B:119:LEU:HD12	1:B:124:HIS:CD2	2.55	0.41
1:E:151:ILE:HD12	1:E:151:ILE:N	2.35	0.41
1:E:129:LEU:CB	1:E:291:VAL:HG11	2.50	0.41
1:A:82:ASP:OD2	1:B:265:ARG:NH2	2.52	0.41
1:D:53:PHE:CD1	1:D:95:MSE:HE2	2.52	0.41
1:A:51:MSE:HE2	1:A:51:MSE:HB2	1.81	0.41
1:C:195:LEU:HD23	1:C:195:LEU:HA	1.81	0.41
1:B:73:GLN:HG3	1:B:73:GLN:O	2.20	0.41
1:E:223:VAL:HG12	1:E:223:VAL:O	2.21	0.41
1:E:145:ILE:CG2	1:E:146:ALA:N	2.83	0.41
1:B:49:THR:HG22	1:B:95:MSE:HE1	2.02	0.41
1:D:194:ALA:CB	1:D:199:ALA:HB3	2.50	0.41
1:E:175:ALA:HB2	1:E:210:ALA:HB2	2.02	0.41
1:D:118:ALA:O	1:D:119:LEU:CB	2.68	0.41
1:F:245:GLU:HA	1:F:273:ILE:HD13	2.03	0.41
1:E:265:ARG:NH2	1:F:82:ASP:OD1	2.53	0.41
1:C:281:ILE:O	1:C:284:GLU:HB3	2.20	0.41
1:A:53:PHE:CD1	1:A:292:VAL:HG11	2.56	0.41
1:C:49:THR:OG1	1:C:288:ARG:NH1	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:256:ILE:HD13	1:F:256:ILE:HG21	1.78	0.41
1:F:223:VAL:O	1:F:223:VAL:HG12	2.20	0.41
1:D:74:LEU:HD11	1:D:96:MSE:CE	2.51	0.41
1:F:53:PHE:HB2	1:F:95:MSE:HE3	2.03	0.41
1:C:145:ILE:HG22	1:C:146:ALA:N	2.34	0.41
1:C:51:MSE:HE2	1:C:51:MSE:HB2	1.66	0.40
1:F:281:ILE:O	1:F:284:GLU:HB3	2.21	0.40
1:F:49:THR:OG1	1:F:288:ARG:NH1	2.54	0.40
1:A:146:ALA:HB3	1:A:170:PHE:CD2	2.57	0.40
1:F:60:LEU:HA	1:F:60:LEU:HD23	1.95	0.40
1:B:281:ILE:O	1:B:284:GLU:HB3	2.20	0.40
1:D:16:LEU:HA	1:D:16:LEU:HD23	1.93	0.40
1:E:297:CYS:O	1:E:301:ASN:ND2	2.50	0.40
1:E:51:MSE:HE2	1:E:51:MSE:HB2	1.78	0.40
1:D:265:ARG:NH2	1:E:82:ASP:OD2	2.54	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	301/307 (98%)	286 (95%)	14 (5%)	1 (0%)	46	75
1	B	303/307 (99%)	285 (94%)	17 (6%)	1 (0%)	46	75
1	C	297/307 (97%)	284 (96%)	12 (4%)	1 (0%)	46	75
1	D	301/307 (98%)	286 (95%)	14 (5%)	1 (0%)	46	75
1	E	296/307 (96%)	286 (97%)	9 (3%)	1 (0%)	46	75
1	F	292/307 (95%)	279 (96%)	12 (4%)	1 (0%)	46	75
All	All	1790/1842 (97%)	1706 (95%)	78 (4%)	6 (0%)	46	75

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	LEU
1	B	119	LEU
1	C	119	LEU
1	D	119	LEU
1	E	119	LEU
1	F	119	LEU

### 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	236/253 (93%)	235 (100%)	1 (0%)	93	98
1	B	238/253 (94%)	236 (99%)	2 (1%)	86	96
1	C	228/253 (90%)	227 (100%)	1 (0%)	93	98
1	D	228/253 (90%)	226 (99%)	2 (1%)	84	95
1	E	226/253 (89%)	224 (99%)	2 (1%)	84	95
1	F	230/253 (91%)	228 (99%)	2 (1%)	84	95
All	All	1386/1518 (91%)	1376 (99%)	10 (1%)	88	96

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	288	ARG
1	B	239	GLU
1	B	288	ARG
1	C	288	ARG
1	D	203	LEU
1	D	288	ARG
1	E	119	LEU
1	E	288	ARG
1	F	119	LEU
1	F	288	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	GLN
1	A	99	ASN
1	C	45	ASN
1	C	159	ASN
1	D	45	ASN
1	D	99	ASN
1	E	45	ASN
1	E	73	GLN
1	E	159	ASN
1	F	73	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

3 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	PEG	A	307	-	6,6,6	0.67	0	5,5,5	0.22	0
2	PEG	B	307	-	6,6,6	0.67	0	5,5,5	0.40	0
2	PEG	F	307	-	6,6,6	0.73	0	5,5,5	0.50	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PEG	A	307	-	-	0/4/4/4	0/0/0/0
2	PEG	B	307	-	-	0/4/4/4	0/0/0/0
2	PEG	F	307	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	290/307 (94%)	-0.09	13 (4%) 37 36	35, 68, 128, 178	0
1	B	292/307 (95%)	-0.22	5 (1%) 73 74	30, 56, 110, 157	0
1	C	288/307 (93%)	0.20	22 (7%) 17 15	41, 83, 171, 212	0
1	D	290/307 (94%)	0.12	15 (5%) 31 30	34, 84, 149, 170	0
1	E	287/307 (93%)	0.12	14 (4%) 33 32	43, 102, 172, 187	0
1	F	286/307 (93%)	-0.24	6 (2%) 67 68	36, 62, 133, 161	0
All	All	1733/1842 (94%)	-0.02	75 (4%) 39 38	30, 73, 151, 212	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	241	PHE	9.2
1	C	177	PRO	7.4
1	B	75	SER	6.7
1	E	241	PHE	6.7
1	D	235	ILE	6.4
1	B	76	ARG	5.9
1	C	151	ILE	5.1
1	E	238	PHE	4.9
1	C	227	GLU	4.8
1	D	75	SER	4.6
1	A	76	ARG	4.5
1	B	74	LEU	4.5
1	D	170	PHE	4.4
1	C	303	ARG	4.1
1	C	237	GLU	3.8
1	D	237	GLU	3.7
1	D	180	TYR	3.7
1	C	228	GLU	3.7
1	A	303	ARG	3.5

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Mol	Chain	Res	Type	RSRZ
1	C	235	ILE	3.4
1	C	222	TRP	3.4
1	B	226	GLY	3.3
1	E	177	PRO	3.3
1	A	230	GLU	3.3
1	E	170	PHE	3.2
1	F	224	SER	3.2
1	C	256	ILE	3.2
1	C	154	SER	3.2
1	A	180	TYR	3.1
1	F	229	ASN	3.1
1	D	255	ALA	3.1
1	E	180	TYR	3.1
1	D	215	ASP	3.1
1	E	224	SER	3.1
1	C	238	PHE	3.0
1	D	253	LYS	3.0
1	F	227	GLU	2.9
1	E	216	VAL	2.8
1	A	238	PHE	2.8
1	A	229	ASN	2.8
1	C	152	GLY	2.8
1	E	228	GLU	2.8
1	E	150	PHE	2.8
1	A	226	GLY	2.8
1	A	267	TYR	2.7
1	D	254	ASP	2.7
1	A	75	SER	2.6
1	E	227	GLU	2.6
1	E	151	ILE	2.6
1	C	243	ILE	2.6
1	D	216	VAL	2.6
1	C	254	ASP	2.5
1	C	231	LYS	2.5
1	A	241	PHE	2.3
1	A	253	LYS	2.3
1	A	228	GLU	2.3
1	C	239	GLU	2.3
1	E	154	SER	2.3
1	D	256	ILE	2.3
1	F	302	GLN	2.3
1	B	224	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	226	GLY	2.2
1	C	240	GLY	2.2
1	D	234	LYS	2.2
1	E	301	ASN	2.2
1	D	138	TRP	2.1
1	D	179	ASN	2.1
1	D	228	GLU	2.1
1	F	303	ARG	2.0
1	C	233	ARG	2.0
1	C	75	SER	2.0
1	C	224	SER	2.0
1	A	145	ILE	2.0
1	F	110	TYR	2.0
1	E	265	ARG	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	PEG	B	307	7/7	0.90	0.14	-	69,75,86,95	0
2	PEG	A	307	7/7	0.88	0.18	-	72,84,96,98	0
2	PEG	F	307	7/7	0.89	0.30	-	73,79,83,83	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.