



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 04:40 AM GMT

PDB ID : 4JDM  
Title : Secreted Chlamydial Protein PGP3, full-length  
Authors : Galaleldeen, A.; Taylor, A.B.; Chen, D.; Holloway, S.P.; Zhong, G.; Hart, P.J.  
Deposited on : 2013-02-25  
Resolution : 3.10 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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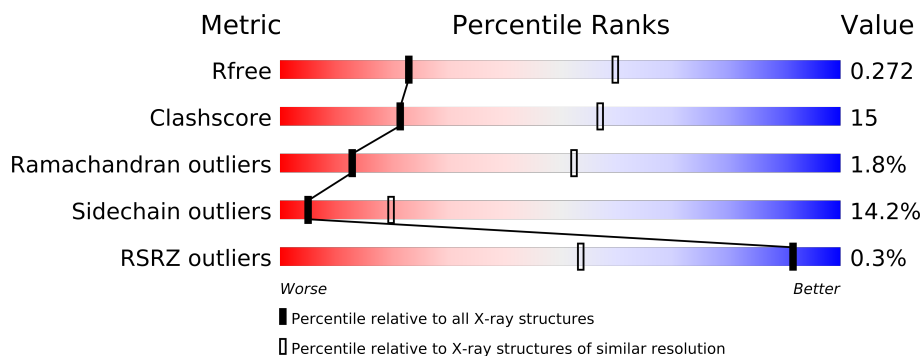
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 3.10 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))
$R_{free}$	66092	1007 (3.18-3.02)
Clashscore	79885	1078 (3.16-3.04)
Ramachandran outliers	78287	1044 (3.16-3.04)
Sidechain outliers	78261	1044 (3.16-3.04)
RSRZ outliers	66119	1008 (3.18-3.02)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	269	
1	B	269	
1	C	269	
1	D	269	
1	E	269	
1	F	269	

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 11712 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Virulence plasmid protein pGP3-D.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	Se	0	0	0
			1939	1215	317	399	4	4			
1	B	263	Total	C	N	O	S	Se	0	0	0
			1941	1217	317	398	4	5			
1	C	269	Total	C	N	O	S	Se	0	0	0
			1976	1238	323	406	4	5			
1	D	263	Total	C	N	O	S	Se	0	0	0
			1939	1215	317	399	4	4			
1	E	263	Total	C	N	O	S	Se	0	0	0
			1941	1217	317	398	4	5			
1	F	269	Total	C	N	O	S	Se	0	0	0
			1976	1238	323	406	4	5			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	EXPRESSION TAG	UNP D7DHH5
A	-3	PRO	-	EXPRESSION TAG	UNP D7DHH5
A	-2	LEU	-	EXPRESSION TAG	UNP D7DHH5
A	-1	GLY	-	EXPRESSION TAG	UNP D7DHH5
A	0	SER	-	EXPRESSION TAG	UNP D7DHH5
B	-4	GLY	-	EXPRESSION TAG	UNP D7DHH5
B	-3	PRO	-	EXPRESSION TAG	UNP D7DHH5
B	-2	LEU	-	EXPRESSION TAG	UNP D7DHH5
B	-1	GLY	-	EXPRESSION TAG	UNP D7DHH5
B	0	SER	-	EXPRESSION TAG	UNP D7DHH5
C	-4	GLY	-	EXPRESSION TAG	UNP D7DHH5
C	-3	PRO	-	EXPRESSION TAG	UNP D7DHH5
C	-2	LEU	-	EXPRESSION TAG	UNP D7DHH5
C	-1	GLY	-	EXPRESSION TAG	UNP D7DHH5
C	0	SER	-	EXPRESSION TAG	UNP D7DHH5
D	-4	GLY	-	EXPRESSION TAG	UNP D7DHH5
D	-3	PRO	-	EXPRESSION TAG	UNP D7DHH5

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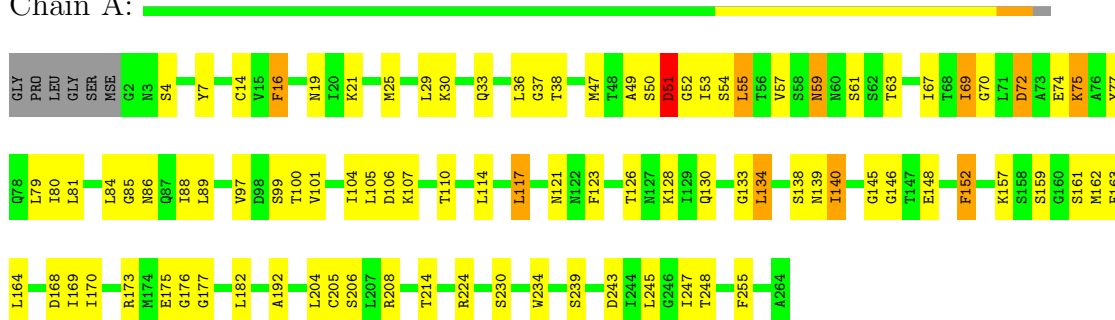
Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	LEU	-	EXPRESSION TAG	UNP D7DHH5
D	-1	GLY	-	EXPRESSION TAG	UNP D7DHH5
D	0	SER	-	EXPRESSION TAG	UNP D7DHH5
E	-4	GLY	-	EXPRESSION TAG	UNP D7DHH5
E	-3	PRO	-	EXPRESSION TAG	UNP D7DHH5
E	-2	LEU	-	EXPRESSION TAG	UNP D7DHH5
E	-1	GLY	-	EXPRESSION TAG	UNP D7DHH5
E	0	SER	-	EXPRESSION TAG	UNP D7DHH5
F	-4	GLY	-	EXPRESSION TAG	UNP D7DHH5
F	-3	PRO	-	EXPRESSION TAG	UNP D7DHH5
F	-2	LEU	-	EXPRESSION TAG	UNP D7DHH5
F	-1	GLY	-	EXPRESSION TAG	UNP D7DHH5
F	0	SER	-	EXPRESSION TAG	UNP D7DHH5

### 3 Residue-property plots

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.

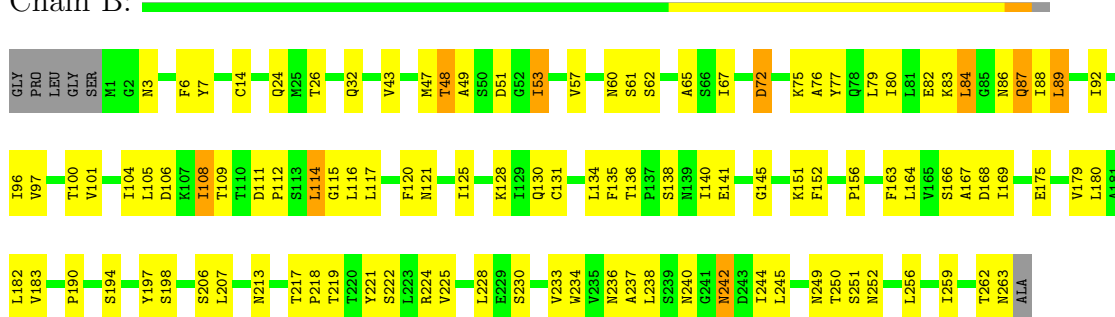
#### • Molecule 1: Virulence plasmid protein pGP3-D

Chain A:



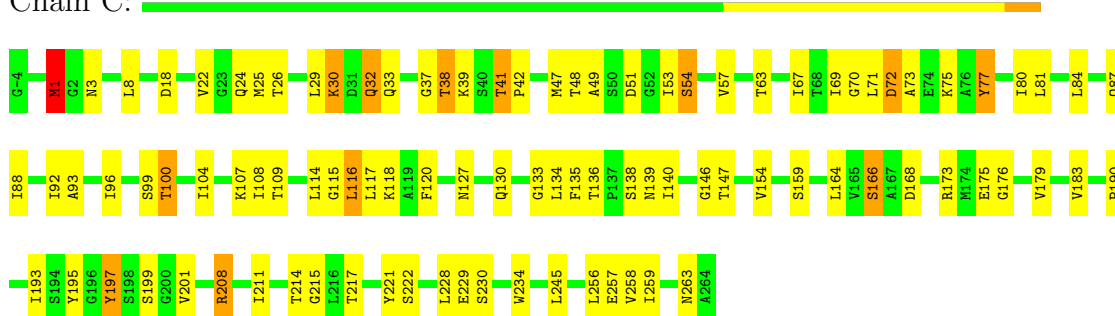
#### • Molecule 1: Virulence plasmid protein pGP3-D

Chain B:



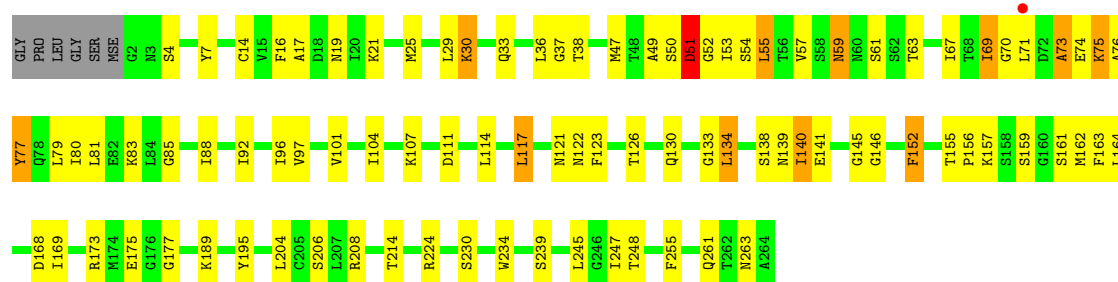
#### • Molecule 1: Virulence plasmid protein pGP3-D

Chain C:



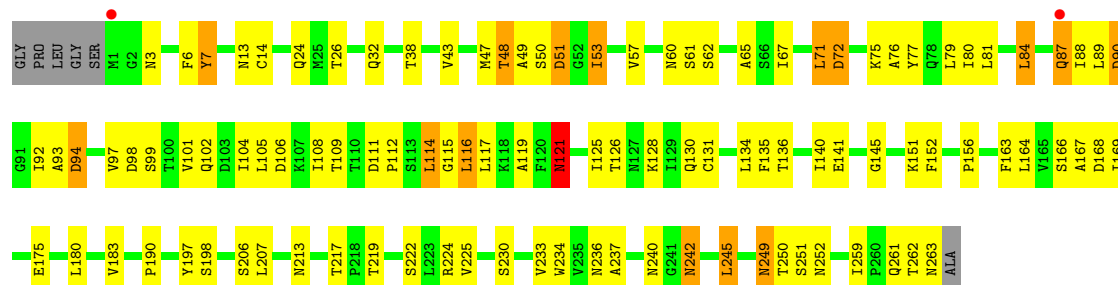
#### • Molecule 1: Virulence plasmid protein pGP3-D

Chain D:



• Molecule 1: Virulence plasmid protein pGP3-D

Chain E:



• Molecule 1: Virulence plasmid protein pGP3-D

Chain F:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	188.70Å 223.79Å 79.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.12 – 3.10 48.12 – 3.10	Depositor EDS
% Data completeness (in resolution range)	91.4 (48.12-3.10) 90.8 (48.12-3.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.66 (at 3.12Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1161)	Depositor
R, $R_{free}$	0.231 , 0.264 0.237 , 0.272	Depositor DCC
$R_{free}$ test set	2798 reflections (4.99%)	DCC
Wilson B-factor (Å <sup>2</sup> )	62.2	Xtriage
Anisotropy	0.728	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.25 , 16.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 56739 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	11712	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	71.0	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 44.12 % 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 1.5988e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

## 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	0.59	1/1960 (0.1%)	0.73	0/2660
1	B	0.58	2/1962 (0.1%)	0.73	0/2663
1	C	0.57	0/1997	0.73	0/2708
1	D	0.56	0/1960	0.72	0/2660
1	E	0.55	1/1962 (0.1%)	0.72	1/2663 (0.0%)
1	F	0.56	0/1997	0.72	0/2708
All	All	0.57	4/11838 (0.0%)	0.73	1/16062 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	121	ASN	CG-OD1	-5.96	1.10	1.24
1	B	121	ASN	CG-ND2	-5.76	1.18	1.32
1	A	205	CYS	CB-SG	-5.49	1.72	1.81
1	E	121	ASN	CG-ND2	-5.22	1.19	1.32

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	71	LEU	CB-CG-CD2	5.44	120.25	111.00

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1939	0	1957	76	0
1	B	1941	0	1964	64	0
1	C	1976	0	1998	57	0
1	D	1939	0	1957	73	0
1	E	1941	0	1964	79	0
1	F	1976	0	1998	61	0
All	All	11712	0	11838	361	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 15.

All (361) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:123:PHE:CE2	1:A:152:PHE:HB3	1.71	1.25
1:D:123:PHE:CE2	1:D:152:PHE:HB3	1.74	1.21
1:A:123:PHE:HE2	1:A:152:PHE:HB3	1.11	0.96
1:A:123:PHE:CD2	1:A:152:PHE:HB3	2.05	0.92
1:E:156:PRO:HG3	1:E:213:ASN:HD22	1.35	0.92
1:A:97:VAL:HG22	1:C:100:THR:HG22	1.52	0.91
1:D:123:PHE:CD2	1:D:152:PHE:HB3	2.05	0.90
1:D:123:PHE:HD2	1:D:152:PHE:HD1	1.21	0.89
1:E:88:ILE:CG2	1:E:92:ILE:HD13	2.03	0.89
1:B:156:PRO:HG3	1:B:213:ASN:HD22	1.38	0.89
1:A:123:PHE:HD2	1:A:152:PHE:CD1	1.91	0.88
1:C:96:ILE:HA	1:C:99:SER:HB2	1.56	0.88
1:D:123:PHE:HD2	1:D:152:PHE:CD1	1.93	0.87
1:A:168:ASP:HB3	1:A:204:LEU:HD21	1.56	0.86
1:E:53:ILE:HG13	1:F:70:GLY:HA2	1.57	0.85
1:E:38:THR:HB	1:F:59:ASN:OD1	1.76	0.84
1:D:123:PHE:HE2	1:D:152:PHE:HB3	1.42	0.83
1:E:111:ASP:HB3	1:E:114:LEU:HG	1.59	0.82
1:D:168:ASP:HB3	1:D:204:LEU:HD21	1.61	0.81
1:B:53:ILE:HG13	1:C:70:GLY:HA2	1.63	0.81
1:E:88:ILE:HG23	1:E:92:ILE:HD13	1.64	0.78
1:B:87:GLN:H	1:B:87:GLN:CD	1.89	0.76
1:E:72:ASP:HB2	1:E:76:ALA:H	1.51	0.75
1:C:136:THR:OG1	1:C:139:ASN:ND2	2.20	0.74
1:C:140:ILE:HD11	1:C:193:ILE:HD13	1.69	0.74
1:E:72:ASP:N	1:E:72:ASP:OD2	2.21	0.73
1:A:123:PHE:HE2	1:A:152:PHE:CB	1.98	0.72
1:B:111:ASP:HB3	1:B:114:LEU:HG	1.71	0.72
1:E:76:ALA:HA	1:E:79:LEU:HD12	1.71	0.72
1:D:195:TYR:OH	1:E:249:ASN:ND2	2.23	0.71

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:123:PHE:CD2	1:D:152:PHE:CD1	2.80	0.70
1:C:208:ARG:HB3	1:C:208:ARG:NH1	2.07	0.70
1:A:47:MSE:HG2	1:B:67:ILE:HD11	1.73	0.70
1:E:128:LYS:HZ2	1:E:242:ASN:HA	1.57	0.69
1:E:88:ILE:HG22	1:E:92:ILE:HD13	1.74	0.69
1:F:37:GLY:HA2	1:F:42:PRO:HB3	1.75	0.69
1:A:16:PHE:O	1:A:16:PHE:HD1	1.76	0.69
1:F:136:THR:OG1	1:F:139:ASN:ND2	2.26	0.68
1:B:53:ILE:HD11	1:C:71:LEU:HG	1.75	0.68
1:D:57:VAL:HG12	1:D:67:ILE:HG22	1.76	0.67
1:D:123:PHE:CE2	1:D:152:PHE:CB	2.67	0.67
1:C:37:GLY:HA2	1:C:42:PRO:HB3	1.76	0.67
1:B:128:LYS:HD2	1:B:237:ALA:HB2	1.75	0.67
1:A:67:ILE:HD11	1:C:47:MSE:HE3	1.77	0.66
1:C:208:ARG:HH11	1:C:208:ARG:HB3	1.59	0.66
1:E:128:LYS:HD2	1:E:237:ALA:HB2	1.76	0.65
1:F:140:ILE:HD11	1:F:193:ILE:HD13	1.79	0.65
1:D:47:MSE:HG2	1:E:67:ILE:HD11	1.78	0.65
1:E:183:VAL:HG22	1:E:190:PRO:HA	1.79	0.65
1:F:208:ARG:NH1	1:F:208:ARG:HB3	2.11	0.65
1:E:87:GLN:N	1:E:87:GLN:OE1	2.30	0.64
1:F:72:ASP:OD1	1:F:72:ASP:N	2.30	0.64
1:E:57:VAL:HG12	1:E:67:ILE:HG22	1.80	0.64
1:A:72:ASP:OD1	1:A:72:ASP:N	2.31	0.64
1:E:128:LYS:NZ	1:E:242:ASN:HA	2.13	0.64
1:D:121:ASN:HB3	1:D:123:PHE:CE1	2.33	0.64
1:B:87:GLN:HG3	1:C:77:TYR:OH	1.98	0.63
1:A:57:VAL:HG12	1:A:67:ILE:HG22	1.80	0.63
1:B:49:ALA:HB1	1:B:53:ILE:HB	1.81	0.63
1:A:16:PHE:O	1:A:16:PHE:CD1	2.52	0.63
1:B:72:ASP:OD1	1:B:72:ASP:N	2.30	0.63
1:E:72:ASP:HB3	1:E:75:LYS:HB2	1.81	0.63
1:C:72:ASP:OD1	1:C:72:ASP:N	2.30	0.62
1:F:-1:GLY:O	1:F:3:ASN:ND2	2.33	0.61
1:B:183:VAL:HG22	1:B:190:PRO:HA	1.82	0.61
1:E:49:ALA:HB1	1:E:53:ILE:HB	1.81	0.61
1:B:57:VAL:HG12	1:B:67:ILE:HG22	1.82	0.61
1:E:140:ILE:HG23	1:E:141:GLU:HG2	1.83	0.61
1:E:116:LEU:N	1:E:116:LEU:HD13	2.16	0.61
1:E:125:ILE:O	1:E:236:ASN:ND2	2.34	0.61
1:F:57:VAL:HG12	1:F:67:ILE:HG22	1.80	0.61
1:F:166:SER:HB3	1:F:208:ARG:HD2	1.82	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:32:GLN:HE21	1:C:57:VAL:HG21	1.67	0.60
1:C:57:VAL:HG12	1:C:67:ILE:HG22	1.82	0.60
1:B:140:ILE:HG23	1:B:141:GLU:HG2	1.83	0.60
1:F:208:ARG:HH11	1:F:208:ARG:HB3	1.67	0.59
1:A:4:SER:HB2	1:B:6:PHE:CD2	2.38	0.58
1:A:51:ASP:HB2	1:A:79:LEU:HD12	1.85	0.58
1:B:75:LYS:O	1:B:79:LEU:HG	2.02	0.58
1:A:123:PHE:HD2	1:A:152:PHE:HD1	1.48	0.58
1:A:123:PHE:CD2	1:A:152:PHE:CD1	2.83	0.58
1:E:93:ALA:O	1:E:97:VAL:HG23	2.04	0.57
1:F:96:ILE:HA	1:F:99:SER:HB2	1.86	0.57
1:F:32:GLN:HE21	1:F:57:VAL:HG21	1.68	0.57
1:E:109:THR:HA	1:E:117:LEU:HD11	1.87	0.57
1:D:77:TYR:HA	1:D:80:ILE:HD12	1.86	0.57
1:D:245:LEU:HD13	1:F:140:ILE:HD13	1.86	0.57
1:A:4:SER:C	1:A:16:PHE:CE1	2.78	0.57
1:B:125:ILE:O	1:B:236:ASN:ND2	2.38	0.57
1:E:89:LEU:HD12	1:E:90:ASP:N	2.20	0.56
1:F:84:LEU:O	1:F:88:ILE:HB	2.05	0.56
1:A:161:SER:OG	1:A:162:MSE:N	2.38	0.56
1:A:104:ILE:HD11	1:B:101:VAL:HG22	1.86	0.56
1:C:166:SER:HB3	1:C:208:ARG:HD2	1.87	0.56
1:C:118:LYS:HB3	1:C:257:GLU:HB3	1.88	0.56
1:F:130:GLN:HB2	1:F:234:TRP:CH2	2.40	0.56
1:D:130:GLN:HB2	1:D:234:TRP:CZ2	2.41	0.56
1:D:51:ASP:HB2	1:D:79:LEU:HD12	1.86	0.56
1:F:135:PHE:CD2	1:F:179:VAL:HB	2.42	0.55
1:D:97:VAL:HG13	1:F:100:THR:HB	1.88	0.55
1:A:16:PHE:CD1	1:A:16:PHE:C	2.79	0.55
1:D:208:ARG:HB3	1:D:208:ARG:HH11	1.70	0.55
1:C:114:LEU:N	1:C:115:GLY:HA2	2.21	0.55
1:D:81:LEU:HA	1:D:85:GLY:HA3	1.88	0.55
1:A:4:SER:C	1:A:16:PHE:HE1	2.10	0.55
1:C:25:MSE:HE1	1:C:29:LEU:HD23	1.89	0.54
1:D:161:SER:OG	1:D:162:MSE:N	2.39	0.54
1:D:122:ASN:C	1:D:123:PHE:HD1	2.09	0.54
1:C:1:MSE:HA	1:C:18:ASP:OD1	2.08	0.54
1:B:109:THR:HA	1:B:117:LEU:HD11	1.90	0.54
1:A:173:ARG:NH1	1:A:177:GLY:O	2.41	0.54
1:D:101:VAL:HG23	1:F:104:ILE:HG12	1.89	0.54
1:A:16:PHE:HD1	1:A:16:PHE:C	2.10	0.54
1:B:86:ASN:O	1:B:89:LEU:HG	2.07	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:240:ASN:OD1	1:E:242:ASN:OD1	2.25	0.53
1:F:84:LEU:HB3	1:F:88:ILE:HD13	1.89	0.53
1:D:71:LEU:HD21	1:E:71:LEU:HD21	1.90	0.53
1:D:123:PHE:HE2	1:D:152:PHE:CB	2.16	0.53
1:A:36:LEU:HD11	1:B:32:GLN:HA	1.89	0.53
1:D:75:LYS:O	1:D:79:LEU:HG	2.08	0.53
1:C:183:VAL:HG22	1:C:190:PRO:HA	1.91	0.53
1:E:156:PRO:HG3	1:E:213:ASN:ND2	2.15	0.53
1:C:130:GLN:HB2	1:C:234:TRP:CH2	2.44	0.52
1:F:111:ASP:HB3	1:F:114:LEU:HG	1.92	0.52
1:D:67:ILE:HD11	1:F:47:MSE:HE3	1.92	0.52
1:C:30:LYS:O	1:C:33:GLN:HB2	2.10	0.52
1:B:156:PRO:HG3	1:B:213:ASN:ND2	2.18	0.52
1:F:114:LEU:N	1:F:115:GLY:HA2	2.24	0.52
1:A:247:ILE:HD12	1:A:248:THR:O	2.10	0.52
1:E:116:LEU:N	1:E:116:LEU:CD1	2.73	0.52
1:E:98:ASP:O	1:E:102:GLN:HB2	2.09	0.51
1:E:131:CYS:HB2	1:E:233:VAL:O	2.10	0.51
1:B:96:ILE:O	1:B:100:THR:HG23	2.10	0.51
1:E:53:ILE:HG12	1:E:72:ASP:OD1	2.09	0.51
1:F:53:ILE:HG23	1:F:70:GLY:O	2.11	0.51
1:A:54:SER:O	1:A:69:ILE:HA	2.10	0.51
1:B:76:ALA:O	1:B:80:ILE:HG13	2.11	0.51
1:C:93:ALA:O	1:C:96:ILE:HG13	2.11	0.51
1:E:134:LEU:HD21	1:E:230:SER:HA	1.91	0.51
1:C:211:ILE:HD12	1:C:221:TYR:CE2	2.45	0.51
1:D:121:ASN:HB3	1:D:123:PHE:HE1	1.74	0.51
1:B:72:ASP:HB3	1:B:75:LYS:HB2	1.93	0.51
1:F:24:GLN:N	1:F:41:THR:OG1	2.44	0.51
1:F:183:VAL:HG22	1:F:190:PRO:HA	1.93	0.51
1:D:54:SER:O	1:D:69:ILE:HA	2.11	0.51
1:F:53:ILE:CD1	1:F:71:LEU:HD23	2.41	0.50
1:B:104:ILE:HG21	1:C:104:ILE:HD12	1.93	0.50
1:A:97:VAL:HG21	1:C:99:SER:HB3	1.93	0.50
1:B:79:LEU:HA	1:B:82:GLU:HB3	1.92	0.50
1:F:54:SER:O	1:F:69:ILE:HA	2.11	0.50
1:C:77:TYR:O	1:C:81:LEU:HG	2.12	0.50
1:D:73:ALA:O	1:D:76:ALA:HB3	2.10	0.50
1:F:118:LYS:HB3	1:F:257:GLU:HB3	1.94	0.50
1:A:106:ASP:O	1:A:110:THR:HG23	2.11	0.50
1:D:19:ASN:HB3	1:E:14:CYS:SG	2.51	0.50
1:E:72:ASP:CB	1:E:75:LYS:HB2	2.40	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:87:GLN:CA	1:E:87:GLN:OE1	2.59	0.50
1:A:123:PHE:CD2	1:A:152:PHE:CB	2.90	0.50
1:D:263:ASN:OD1	1:E:121:ASN:ND2	2.44	0.50
1:E:76:ALA:O	1:E:80:ILE:HG13	2.12	0.49
1:B:87:GLN:CD	1:B:87:GLN:N	2.61	0.49
1:A:130:GLN:HB2	1:A:234:TRP:CZ2	2.46	0.49
1:D:134:LEU:HD21	1:D:230:SER:HA	1.94	0.49
1:A:208:ARG:HH11	1:A:208:ARG:HB3	1.77	0.49
1:A:80:ILE:HD12	1:B:77:TYR:HB2	1.95	0.49
1:D:123:PHE:CD2	1:D:152:PHE:CB	2.88	0.49
1:A:133:GLY:C	1:A:146:GLY:HA2	2.33	0.49
1:E:24:GLN:HB2	1:E:43:VAL:HG12	1.94	0.49
1:E:135:PHE:CE1	1:E:145:GLY:HA3	2.47	0.49
1:D:4:SER:HB2	1:E:6:PHE:CD2	2.47	0.49
1:A:100:THR:HG22	1:B:97:VAL:HG22	1.94	0.49
1:A:4:SER:O	1:A:16:PHE:CE1	2.66	0.49
1:B:108:ILE:HG22	1:B:109:THR:HG23	1.95	0.49
1:E:130:GLN:HG3	1:E:234:TRP:CE2	2.48	0.49
1:F:25:MSE:HE1	1:F:29:LEU:HD23	1.94	0.49
1:D:111:ASP:O	1:D:117:LEU:HD21	2.13	0.48
1:E:108:ILE:HG23	1:E:116:LEU:HD23	1.94	0.48
1:F:78:GLN:O	1:F:82:GLU:HB2	2.14	0.48
1:C:54:SER:O	1:C:69:ILE:HA	2.14	0.48
1:D:123:PHE:N	1:D:123:PHE:CD1	2.81	0.48
1:D:49:ALA:HB1	1:D:53:ILE:HB	1.94	0.48
1:A:121:ASN:HB3	1:A:123:PHE:CE1	2.49	0.48
1:F:49:ALA:HB1	1:F:53:ILE:N	2.29	0.48
1:B:24:GLN:HB2	1:B:43:VAL:HG12	1.94	0.48
1:A:123:PHE:HE1	1:C:263:ASN:OD1	1.96	0.48
1:E:51:ASP:HB2	1:E:79:LEU:CD1	2.44	0.48
1:A:4:SER:O	1:A:16:PHE:HE1	1.96	0.48
1:A:54:SER:HB3	1:A:70:GLY:N	2.29	0.48
1:F:128:LYS:NZ	1:F:241:GLY:O	2.33	0.48
1:A:170:ILE:HD13	1:C:195:TYR:HD1	1.79	0.48
1:D:123:PHE:CD2	1:D:152:PHE:HD1	2.13	0.48
1:D:208:ARG:HB3	1:D:208:ARG:NH1	2.28	0.48
1:B:180:LEU:HG	1:B:225:VAL:HG22	1.96	0.48
1:F:30:LYS:O	1:F:33:GLN:HB2	2.14	0.48
1:F:228:LEU:O	1:F:230:SER:N	2.43	0.47
1:E:106:ASP:N	1:E:106:ASP:OD1	2.47	0.47
1:B:182:LEU:HD11	1:B:221:TYR:HB3	1.96	0.47
1:D:130:GLN:HB2	1:D:234:TRP:CH2	2.49	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:76:ALA:O	1:D:80:ILE:HG13	2.15	0.47
1:B:135:PHE:CE1	1:B:145:GLY:HA3	2.50	0.47
1:D:140:ILE:HG12	1:E:245:LEU:HD23	1.96	0.47
1:B:111:ASP:HA	1:B:112:PRO:HD3	1.76	0.47
1:A:49:ALA:HB1	1:A:53:ILE:HB	1.96	0.47
1:E:169:ILE:HA	1:E:251:SER:HB3	1.97	0.47
1:D:21:LYS:HE3	1:D:25:MSE:O	2.15	0.47
1:C:107:LYS:HA	1:C:107:LYS:HD3	1.75	0.47
1:D:247:ILE:HD12	1:D:248:THR:O	2.15	0.47
1:A:140:ILE:HD13	1:A:140:ILE:HA	1.74	0.46
1:D:55:LEU:HD22	1:D:69:ILE:HG22	1.96	0.46
1:A:75:LYS:O	1:A:79:LEU:HG	2.15	0.46
1:B:130:GLN:HG3	1:B:234:TRP:CE2	2.50	0.46
1:F:234:TRP:HB3	1:F:237:ALA:HB3	1.98	0.46
1:D:133:GLY:C	1:D:146:GLY:HA2	2.36	0.46
1:A:123:PHE:CD1	1:A:123:PHE:N	2.81	0.46
1:B:83:LYS:HB3	1:B:84:LEU:HG	1.98	0.46
1:C:84:LEU:O	1:C:88:ILE:HD13	2.15	0.46
1:D:104:ILE:HD11	1:E:101:VAL:HG22	1.98	0.46
1:D:134:LEU:HA	1:D:134:LEU:HD12	1.76	0.46
1:E:145:GLY:O	1:E:224:ARG:HD3	2.16	0.46
1:B:167:ALA:HB3	1:B:207:LEU:HB3	1.98	0.46
1:C:32:GLN:HE21	1:C:57:VAL:HG11	1.81	0.46
1:D:145:GLY:O	1:D:224:ARG:HD2	2.16	0.46
1:F:211:ILE:HD12	1:F:221:TYR:CE2	2.50	0.46
1:F:127:ASN:N	1:F:127:ASN:OD1	2.48	0.46
1:E:180:LEU:HG	1:E:225:VAL:HG22	1.98	0.46
1:F:38:THR:OG1	1:F:39:LYS:N	2.49	0.46
1:F:166:SER:HB3	1:F:208:ARG:CD	2.46	0.45
1:C:72:ASP:HB2	1:C:75:LYS:HB2	1.98	0.45
1:E:90:ASP:O	1:E:94:ASP:HB3	2.16	0.45
1:E:197:TYR:CE2	1:F:197:TYR:HE2	2.34	0.45
1:C:135:PHE:CD2	1:C:179:VAL:HB	2.52	0.45
1:C:24:GLN:N	1:C:41:THR:OG1	2.48	0.45
1:C:108:ILE:HG22	1:C:109:THR:HG23	1.98	0.45
1:A:51:ASP:HB2	1:A:79:LEU:CD1	2.47	0.45
1:A:134:LEU:HD21	1:A:230:SER:HA	1.97	0.45
1:E:88:ILE:HG23	1:E:92:ILE:CD1	2.42	0.45
1:D:75:LYS:H	1:D:75:LYS:HG2	1.44	0.45
1:B:105:LEU:O	1:B:109:THR:HG23	2.16	0.45
1:A:208:ARG:HB3	1:A:208:ARG:NH1	2.32	0.45
1:A:50:SER:C	1:A:52:GLY:H	2.20	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:238:LEU:HD21	1:B:244:ILE:HG23	1.99	0.45
1:E:61:SER:O	1:E:65:ALA:HB2	2.17	0.45
1:B:197:TYR:CE2	1:C:197:TYR:HE2	2.34	0.45
1:E:51:ASP:HB2	1:E:79:LEU:HD11	1.97	0.45
1:D:17:ALA:HB1	1:E:13:ASN:O	2.17	0.45
1:A:72:ASP:HB2	1:A:75:LYS:HG3	1.99	0.45
1:E:261:GLN:HB3	1:F:157:LYS:HE3	1.99	0.45
1:A:37:GLY:O	1:B:32:GLN:N	2.47	0.45
1:D:54:SER:HB3	1:D:70:GLY:N	2.32	0.45
1:A:134:LEU:HD12	1:A:134:LEU:HA	1.68	0.45
1:E:50:SER:N	1:F:68:THR:HG22	2.32	0.45
1:B:48:THR:HG22	1:B:49:ALA:H	1.81	0.44
1:A:130:GLN:HB2	1:A:234:TRP:CH2	2.52	0.44
1:D:50:SER:C	1:D:52:GLY:H	2.20	0.44
1:C:120:PHE:HD2	1:C:256:LEU:HD13	1.82	0.44
1:A:123:PHE:CD2	1:A:152:PHE:HD1	2.30	0.44
1:B:128:LYS:NZ	1:B:242:ASN:HA	2.31	0.44
1:E:104:ILE:HG21	1:F:104:ILE:HD12	1.99	0.44
1:C:77:TYR:HA	1:C:80:ILE:HD12	1.99	0.44
1:A:55:LEU:HD22	1:A:69:ILE:HG22	2.00	0.44
1:A:139:ASN:ND2	1:A:145:GLY:HA2	2.32	0.44
1:E:183:VAL:HG22	1:E:190:PRO:CA	2.47	0.44
1:F:79:LEU:O	1:F:83:LYS:HG2	2.17	0.44
1:B:61:SER:O	1:B:65:ALA:HB2	2.18	0.44
1:C:92:ILE:O	1:C:96:ILE:HG23	2.17	0.44
1:A:145:GLY:O	1:A:224:ARG:HD2	2.18	0.44
1:B:131:CYS:HB2	1:B:233:VAL:O	2.17	0.44
1:E:89:LEU:HD12	1:E:89:LEU:C	2.38	0.44
1:F:214:THR:OG1	1:F:215:GLY:N	2.50	0.44
1:A:21:LYS:HE3	1:A:25:MSE:O	2.17	0.44
1:A:97:VAL:HG13	1:C:100:THR:HA	1.99	0.44
1:C:49:ALA:HB1	1:C:53:ILE:N	2.33	0.44
1:C:114:LEU:H	1:C:115:GLY:HA2	1.83	0.43
1:B:145:GLY:O	1:B:224:ARG:HD3	2.17	0.43
1:B:240:ASN:OD1	1:B:242:ASN:OD1	2.35	0.43
1:A:19:ASN:HB3	1:B:14:CYS:SG	2.58	0.43
1:A:245:LEU:HD13	1:C:140:ILE:HD13	2.00	0.43
1:F:99:SER:O	1:F:103:ASP:HB2	2.18	0.43
1:A:84:LEU:O	1:A:86:ASN:N	2.49	0.43
1:B:240:ASN:OD1	1:B:242:ASN:CG	2.57	0.43
1:F:140:ILE:O	1:F:140:ILE:HG13	2.19	0.43
1:B:115:GLY:HA2	1:B:259:ILE:HD13	2.00	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:105:LEU:O	1:B:108:ILE:HG22	2.19	0.43
1:B:106:ASP:N	1:B:106:ASP:OD1	2.51	0.43
1:D:29:LEU:HD22	1:D:33:GLN:HB3	2.01	0.43
1:C:133:GLY:O	1:C:146:GLY:HA2	2.18	0.43
1:D:117:LEU:HD12	1:D:157:LYS:HB3	2.01	0.43
1:E:101:VAL:O	1:E:105:LEU:HG	2.17	0.43
1:B:88:ILE:H	1:B:88:ILE:HG13	1.52	0.43
1:A:128:LYS:HE2	1:A:243:ASP:CG	2.39	0.43
1:A:107:LYS:HB2	1:A:107:LYS:HE3	1.86	0.43
1:D:173:ARG:NH1	1:D:177:GLY:O	2.52	0.42
1:B:120:PHE:HD1	1:B:256:LEU:HD13	1.84	0.42
1:F:134:LEU:HD23	1:F:230:SER:HA	2.01	0.42
1:D:133:GLY:O	1:D:146:GLY:HA2	2.18	0.42
1:A:148:GLU:HA	1:A:224:ARG:HA	2.01	0.42
1:E:126:THR:HG22	1:E:250:THR:HG22	2.02	0.42
1:C:245:LEU:HD23	1:C:245:LEU:HA	1.86	0.42
1:E:48:THR:HG22	1:E:49:ALA:H	1.84	0.42
1:C:214:THR:OG1	1:C:215:GLY:N	2.52	0.42
1:C:228:LEU:O	1:C:230:SER:N	2.45	0.42
1:A:29:LEU:HD22	1:A:33:GLN:HB3	2.01	0.42
1:D:92:ILE:HA	1:D:92:ILE:HD12	1.87	0.42
1:D:36:LEU:HD11	1:E:32:GLN:HA	2.00	0.42
1:B:134:LEU:HD21	1:B:230:SER:HA	2.01	0.42
1:C:38:THR:OG1	1:C:39:LYS:N	2.52	0.42
1:D:140:ILE:HD13	1:D:140:ILE:HA	1.71	0.42
1:D:51:ASP:OD2	1:E:71:LEU:HD12	2.20	0.42
1:F:32:GLN:HE21	1:F:57:VAL:HG11	1.83	0.42
1:D:96:ILE:HD11	1:E:93:ALA:HB2	2.02	0.42
1:D:139:ASN:ND2	1:D:145:GLY:HA2	2.35	0.42
1:B:256:LEU:HD12	1:B:256:LEU:HA	1.87	0.42
1:D:155:THR:HA	1:D:156:PRO:HD2	1.81	0.42
1:A:123:PHE:HD2	1:A:152:PHE:CG	2.35	0.42
1:A:170:ILE:HD13	1:C:195:TYR:CD1	2.54	0.42
1:D:163:PHE:CD1	1:D:255:PHE:HB3	2.54	0.42
1:A:117:LEU:HD12	1:A:157:LYS:HB3	2.01	0.42
1:A:173:ARG:CZ	1:A:176:GLY:HA3	2.50	0.41
1:D:141:GLU:O	1:D:189:LYS:HB3	2.20	0.41
1:E:115:GLY:O	1:E:259:ILE:HD12	2.19	0.41
1:F:86:ASN:O	1:F:89:LEU:HG	2.20	0.41
1:B:47:MSE:HE3	1:C:67:ILE:HD12	2.02	0.41
1:F:79:LEU:HA	1:F:79:LEU:HD13	1.92	0.41
1:F:89:LEU:HA	1:F:92:ILE:HD12	2.02	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:37:GLY:O	1:E:32:GLN:N	2.52	0.41
1:D:83:LYS:HD3	1:D:83:LYS:HA	1.95	0.41
1:A:101:VAL:O	1:A:105:LEU:HG	2.20	0.41
1:E:104:ILE:O	1:E:108:ILE:HD13	2.20	0.41
1:E:47:MSE:HE3	1:F:67:ILE:HD12	2.03	0.41
1:F:133:GLY:O	1:F:146:GLY:HA2	2.21	0.41
1:D:59:ASN:N	1:D:59:ASN:OD1	2.54	0.41
1:A:123:PHE:CE1	1:C:263:ASN:OD1	2.73	0.41
1:F:32:GLN:HE21	1:F:57:VAL:CG2	2.34	0.41
1:E:167:ALA:HB3	1:E:207:LEU:HB3	2.03	0.41
1:F:53:ILE:HA	1:F:53:ILE:HD13	1.94	0.41
1:B:169:ILE:HA	1:B:251:SER:HB3	2.02	0.41
1:F:120:PHE:HD2	1:F:256:LEU:HD13	1.86	0.41
1:E:72:ASP:CG	1:E:76:ALA:HB2	2.41	0.41
1:B:128:LYS:HZ2	1:B:242:ASN:HA	1.86	0.41
1:A:104:ILE:HD11	1:B:101:VAL:HA	2.03	0.41
1:B:194:SER:HB2	1:B:207:LEU:HD12	2.03	0.41
1:D:107:LYS:HB2	1:D:107:LYS:HE3	1.82	0.41
1:C:114:LEU:HD23	1:C:114:LEU:HA	1.88	0.41
1:E:7:TYR:CE1	1:E:14:CYS:HB2	2.55	0.41
1:A:182:LEU:HB3	1:A:192:ALA:HB3	2.03	0.41
1:F:162:MSE:HG2	1:F:212:THR:HA	2.03	0.41
1:F:107:LYS:HD3	1:F:107:LYS:HA	1.83	0.41
1:A:59:ASN:OD1	1:A:59:ASN:N	2.54	0.41
1:D:71:LEU:HD23	1:D:71:LEU:HA	1.88	0.40
1:D:261:GLN:HG3	1:E:119:ALA:HB1	2.01	0.40
1:B:179:VAL:HG23	1:B:228:LEU:HB2	2.03	0.40
1:A:163:PHE:CD1	1:A:255:PHE:HB3	2.55	0.40
1:D:30:LYS:O	1:D:33:GLN:HB2	2.21	0.40
1:C:134:LEU:HD23	1:C:230:SER:HA	2.03	0.40
1:C:173:ARG:CZ	1:C:176:GLY:HA3	2.51	0.40
1:C:199:SER:OG	1:C:201:VAL:HG13	2.20	0.40
1:E:111:ASP:HA	1:E:112:PRO:HD3	1.77	0.40
1:B:114:LEU:HA	1:B:114:LEU:HD22	1.82	0.40
1:E:84:LEU:HB3	1:E:87:GLN:NE2	2.36	0.40
1:F:25:MSE:HE2	1:F:27:GLU:O	2.21	0.40
1:E:77:TYR:CZ	1:E:81:LEU:HD11	2.57	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	261/269 (97%)	244 (94%)	12 (5%)	5 (2%)	12	51	
1	B	261/269 (97%)	245 (94%)	13 (5%)	3 (1%)	21	65	
1	C	267/269 (99%)	243 (91%)	18 (7%)	6 (2%)	10	46	
1	D	261/269 (97%)	243 (93%)	13 (5%)	5 (2%)	12	51	
1	E	261/269 (97%)	241 (92%)	17 (6%)	3 (1%)	21	65	
1	F	267/269 (99%)	240 (90%)	21 (8%)	6 (2%)	10	46	
All	All	1578/1614 (98%)	1456 (92%)	94 (6%)	28 (2%)	13	52	

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	114	LEU
1	C	175	GLU
1	D	114	LEU
1	F	175	GLU
1	A	85	GLY
1	A	117	LEU
1	B	51	ASP
1	C	116	LEU
1	D	117	LEU
1	A	175	GLU
1	C	32	GLN
1	C	229	GLU
1	E	51	ASP
1	F	32	GLN
1	F	101	VAL
1	F	229	GLU
1	A	51	ASP
1	B	175	GLU
1	C	1	MSE
1	D	175	GLU
1	E	175	GLU

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Mol	Chain	Res	Type
1	F	116	LEU
1	C	73	ALA
1	D	51	ASP
1	E	99	SER
1	D	73	ALA
1	F	40	SER
1	B	218	PRO

### 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	223/222 (100%)	193 (86%)	30 (14%)	6	22
1	B	224/222 (101%)	189 (84%)	35 (16%)	4	14
1	C	227/222 (102%)	195 (86%)	32 (14%)	5	20
1	D	223/222 (100%)	197 (88%)	26 (12%)	8	29
1	E	224/222 (101%)	191 (85%)	33 (15%)	4	18
1	F	227/222 (102%)	192 (85%)	35 (15%)	4	15
All	All	1348/1332 (101%)	1157 (86%)	191 (14%)	5	20

All (191) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	TYR
1	A	14	CYS
1	A	16	PHE
1	A	30	LYS
1	A	38	THR
1	A	51	ASP
1	A	55	LEU
1	A	59	ASN
1	A	61	SER
1	A	63	THR
1	A	69	ILE
1	A	72	ASP

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Mol	Chain	Res	Type
1	A	74	GLU
1	A	75	LYS
1	A	77	TYR
1	A	81	LEU
1	A	88	ILE
1	A	89	LEU
1	A	99	SER
1	A	126	THR
1	A	134	LEU
1	A	138	SER
1	A	140	ILE
1	A	152	PHE
1	A	159	SER
1	A	164	LEU
1	A	169	ILE
1	A	206	SER
1	A	214	THR
1	A	239	SER
1	B	3	ASN
1	B	7	TYR
1	B	26	THR
1	B	48	THR
1	B	53	ILE
1	B	60	ASN
1	B	62	SER
1	B	72	ASP
1	B	84	LEU
1	B	87	GLN
1	B	89	LEU
1	B	92	ILE
1	B	108	ILE
1	B	114	LEU
1	B	116	LEU
1	B	136	THR
1	B	138	SER
1	B	151	LYS
1	B	152	PHE
1	B	163	PHE
1	B	164	LEU
1	B	166	SER
1	B	168	ASP
1	B	198	SER

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Mol	Chain	Res	Type
1	B	206	SER
1	B	217	THR
1	B	219	THR
1	B	222	SER
1	B	242	ASN
1	B	245	LEU
1	B	249	ASN
1	B	250	THR
1	B	252	ASN
1	B	262	THR
1	B	263	ASN
1	C	1	MSE
1	C	3	ASN
1	C	8	LEU
1	C	22	VAL
1	C	26	THR
1	C	30	LYS
1	C	38	THR
1	C	41	THR
1	C	48	THR
1	C	51	ASP
1	C	54	SER
1	C	63	THR
1	C	72	ASP
1	C	77	TYR
1	C	87	GLN
1	C	100	THR
1	C	116	LEU
1	C	117	LEU
1	C	127	ASN
1	C	138	SER
1	C	147	THR
1	C	154	VAL
1	C	159	SER
1	C	164	LEU
1	C	166	SER
1	C	168	ASP
1	C	197	TYR
1	C	208	ARG
1	C	217	THR
1	C	222	SER
1	C	258	VAL

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Mol	Chain	Res	Type
1	C	259	ILE
1	D	7	TYR
1	D	14	CYS
1	D	16	PHE
1	D	30	LYS
1	D	38	THR
1	D	51	ASP
1	D	55	LEU
1	D	59	ASN
1	D	61	SER
1	D	63	THR
1	D	69	ILE
1	D	74	GLU
1	D	75	LYS
1	D	77	TYR
1	D	88	ILE
1	D	126	THR
1	D	134	LEU
1	D	138	SER
1	D	140	ILE
1	D	152	PHE
1	D	159	SER
1	D	164	LEU
1	D	169	ILE
1	D	206	SER
1	D	214	THR
1	D	239	SER
1	E	3	ASN
1	E	7	TYR
1	E	26	THR
1	E	48	THR
1	E	53	ILE
1	E	60	ASN
1	E	62	SER
1	E	72	ASP
1	E	84	LEU
1	E	87	GLN
1	E	90	ASP
1	E	94	ASP
1	E	114	LEU
1	E	116	LEU
1	E	121	ASN

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Mol	Chain	Res	Type
1	E	136	THR
1	E	151	LYS
1	E	152	PHE
1	E	163	PHE
1	E	164	LEU
1	E	166	SER
1	E	168	ASP
1	E	198	SER
1	E	206	SER
1	E	217	THR
1	E	219	THR
1	E	222	SER
1	E	242	ASN
1	E	245	LEU
1	E	249	ASN
1	E	252	ASN
1	E	262	THR
1	E	263	ASN
1	F	3	ASN
1	F	22	VAL
1	F	26	THR
1	F	30	LYS
1	F	38	THR
1	F	41	THR
1	F	48	THR
1	F	51	ASP
1	F	54	SER
1	F	59	ASN
1	F	63	THR
1	F	72	ASP
1	F	77	TYR
1	F	79	LEU
1	F	87	GLN
1	F	92	ILE
1	F	98	ASP
1	F	100	THR
1	F	116	LEU
1	F	117	LEU
1	F	127	ASN
1	F	138	SER
1	F	147	THR
1	F	152	PHE

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Mol	Chain	Res	Type
1	F	154	VAL
1	F	159	SER
1	F	164	LEU
1	F	166	SER
1	F	168	ASP
1	F	197	TYR
1	F	199	SER
1	F	217	THR
1	F	222	SER
1	F	258	VAL
1	F	259	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	213	ASN
1	C	32	GLN
1	C	87	GLN
1	C	139	ASN
1	E	213	ASN
1	E	249	ASN
1	F	32	GLN
1	F	139	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

There are no ligands in this entry.



## 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	263/269 (97%)	-0.25	0	100 100	22, 53, 132, 151	0
1	B	263/269 (97%)	-0.23	0	100 100	26, 49, 141, 158	0
1	C	269/269 (100%)	-0.22	0	100 100	22, 53, 136, 146	0
1	D	263/269 (97%)	-0.18	1 (0%)	90 45	23, 55, 144, 152	0
1	E	263/269 (97%)	-0.15	2 (0%)	83 28	31, 55, 144, 159	0
1	F	269/269 (100%)	-0.17	2 (0%)	84 32	25, 52, 143, 152	0
All	All	1590/1614 (98%)	-0.20	5 (0%)	91 53	22, 53, 142, 159	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	47	MSE	4.0
1	D	71	LEU	2.6
1	E	87	GLN	2.5
1	E	1	MSE	2.5
1	F	71	LEU	2.3

### 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 ⓘ

There are no carbohydrates in this entry.

### 6.4 Ligands ⓘ

There are no ligands in this entry.

## 6.5 Other polymers ⓘ

There are no such residues in this entry.