Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 02:34 am BST

PDB ID	:	2G3V
Title	:	Crystal structure of CagS (HP0534, Cag13) from Helicobacter pylori
Authors	:	Cendron, L.; Tasca, E.; Angelini, A.; Seydel, A.; Battistutta, R.; Montecucco, C.; Zanotti, G.
Deposited on	:	2006-02-21
Resolution	:	2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

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

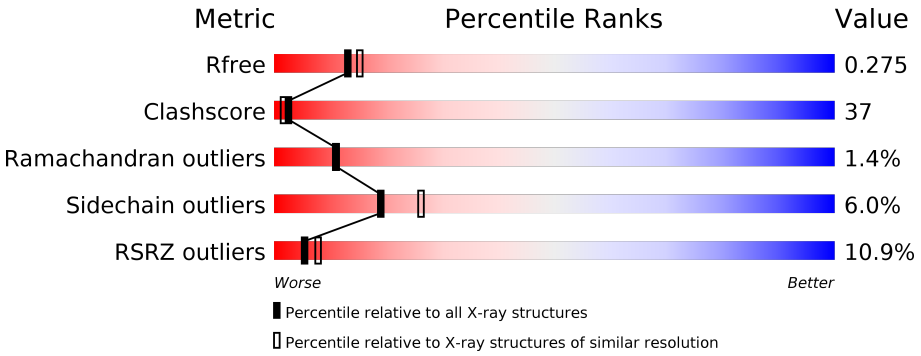
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	208	6% 37% 38% • 23%
1	B	208	7% 37% 37% • 23%
1	C	208	8% 36% 38% • 23%
1	D	208	10% 35% 39% • 23%
2	E	7	43% 57%
2	F	7	43% 57%

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Mol	Chain	Length	Quality of chain
2	G	7	43% 57%
2	H	7	29% 71%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5790 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 CAG pathogenicity island protein 13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	161	Total	C	N	O	Se	0	0	0
			1349	877	218	245	9			
1	B	161	Total	C	N	O	Se	0	0	0
			1349	877	218	245	9			
1	C	161	Total	C	N	O	Se	0	0	0
			1349	877	218	245	9			
1	D	161	Total	C	N	O	Se	0	0	0
			1349	877	218	245	9			

- Molecule 2 is a protein called (UNK)(UNK)(UNK)(UNK)(UNK)(MSE)(UNK).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	7	Total	C	N	O	Se	0	0	0
			39	23	7	8	1			
2	F	7	Total	C	N	O	Se	0	0	0
			39	23	7	8	1			
2	G	7	Total	C	N	O	Se	0	0	0
			39	23	7	8	1			
2	H	7	Total	C	N	O	Se	0	0	0
			39	23	7	8	1			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	54	Total	O	0	0
			54	54		
3	B	48	Total	O	0	0
			48	48		
3	C	57	Total	O	0	0
			57	57		
3	D	62	Total	O	0	0
			62	62		

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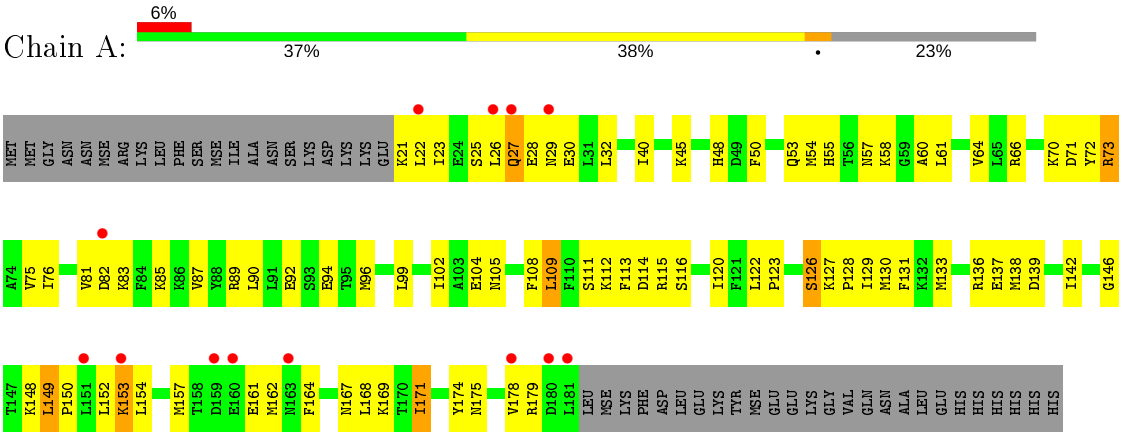
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	E	3	Total O 3 3	0	0
3	F	5	Total O 5 5	0	0
3	G	5	Total O 5 5	0	0
3	H	4	Total O 4 4	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

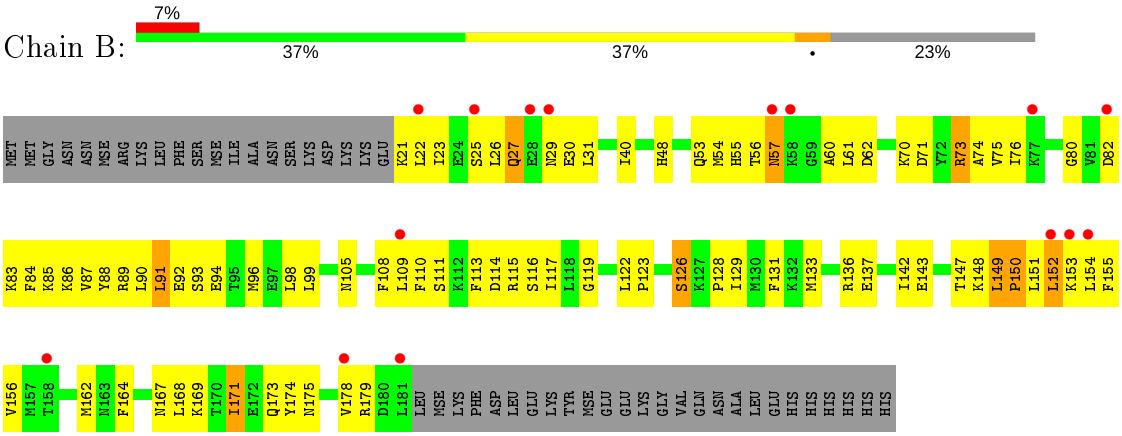
- Molecule 1: CAG pathogenicity island protein 13

Chain A:



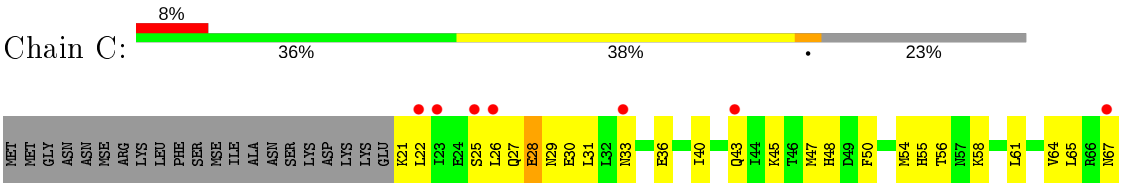
- Molecule 1: CAG pathogenicity island protein 13

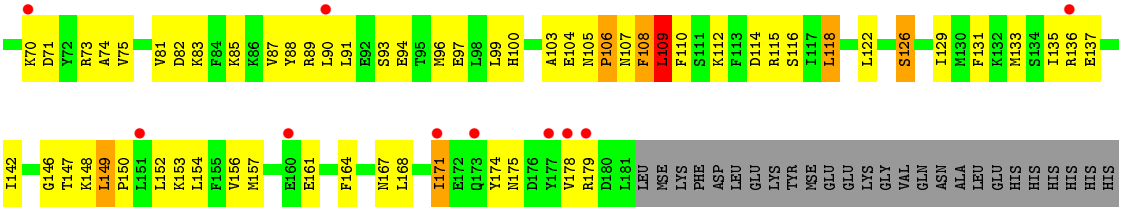
Chain B:



- Molecule 1: CAG pathogenicity island protein 13

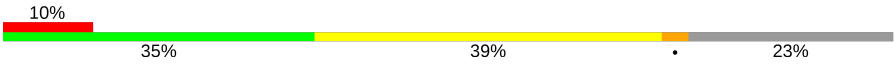
Chain C:

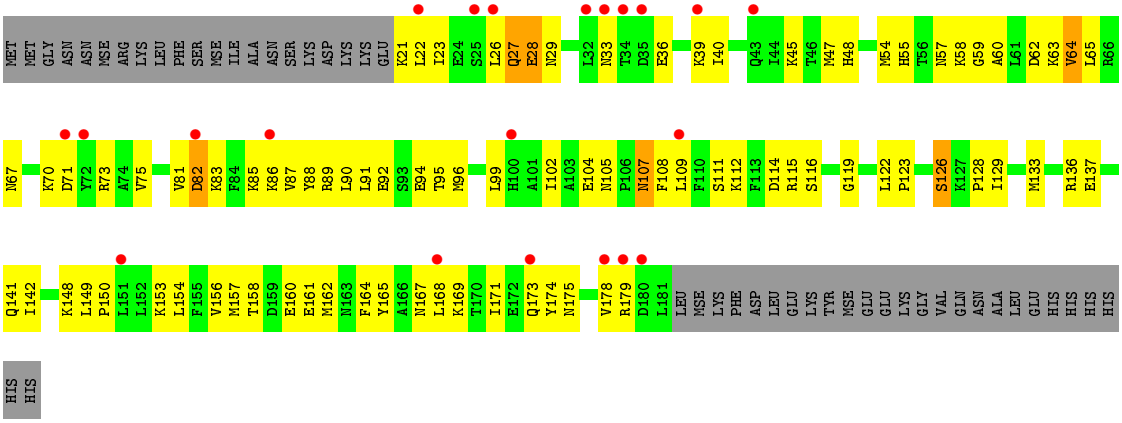




• Molecule 1: CAG pathogenicity island protein 13

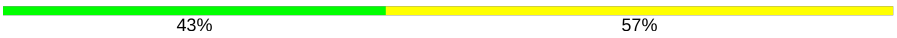
Chain D:

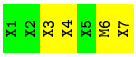




• Molecule 2: (UNK)(UNK)(UNK)(UNK)(UNK)(MSE)(UNK)

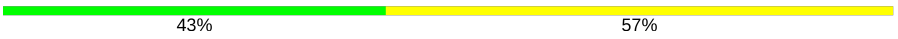
Chain E:

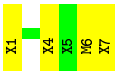




• Molecule 2: (UNK)(UNK)(UNK)(UNK)(UNK)(MSE)(UNK)

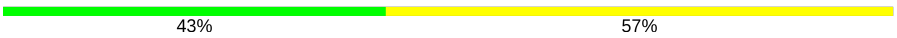
Chain F:





• Molecule 2: (UNK)(UNK)(UNK)(UNK)(UNK)(MSE)(UNK)

Chain G:



• Molecule 2: (UNK)(UNK)(UNK)(UNK)(UNK)(MSE)(UNK)

Chain H:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α, β, γ	124.36Å 124.36Å 55.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.41 – 2.30 39.41 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (39.41-2.30) 99.5 (39.41-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.01	Depositor
$< I/\sigma(I) >$ [1]	1.89 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.219 , 0.279 0.220 , 0.275	Depositor DCC
R_{free} test set	3691 reflections (6.39%)	wwPDB-VP
Wilson B-factor (Å^2)	9.3	Xtriage
Anisotropy	0.589	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 49.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.28$, $< L^2 > = 0.11$	Xtriage
Estimated twinning fraction	0.236 for h,-k,-l	Xtriage
F_o,F_c correlation	0.77	EDS
Total number of atoms	5790	wwPDB-VP
Average B, all atoms (Å^2)	10.0	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 31.65 % 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 1.0684e-03.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|>5$	RMSZ	#$\|Z\|>5$
1	A	0.44	0/1365	0.63	0/1812
1	B	0.43	0/1365	0.64	0/1812
1	C	0.44	0/1365	0.65	2/1812 (0.1%)
1	D	0.43	0/1365	0.62	0/1812
2	E	0.43	0/7	1.04	0/7
2	F	0.62	0/7	1.14	0/7
2	G	0.32	0/7	1.04	0/7
2	H	0.54	0/7	1.15	0/7
All	All	0.43	0/5488	0.64	2/7276 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	118	LEU	CB-CA-C	-6.66	97.55	110.20
1	C	108	PHE	N-CA-C	-5.27	96.78	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	1349	0	1384	94	0
1	B	1349	0	1384	101	0
1	C	1349	0	1384	103	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1349	0	1384	107	0
2	E	39	0	17	5	0
2	F	39	0	17	3	0
2	G	39	0	17	6	0
2	H	39	0	17	5	0
3	A	54	0	0	17	0
3	B	48	0	0	16	0
3	C	57	0	0	15	0
3	D	62	0	0	22	0
3	E	3	0	0	4	0
3	F	5	0	0	1	0
3	G	5	0	0	6	0
3	H	4	0	0	1	0
All	All	5790	0	5604	412	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 37.

All (412) 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:C:153:LYS:HE3	1:C:156:VAL:HG21	1.35	1.09
1:A:139:ASP:HA	3:A:218:HOH:O	1.64	0.97
1:D:21:LYS:HG3	1:D:22:LEU:HG	1.46	0.95
1:C:71:ASP:O	1:C:75:VAL:HG23	1.66	0.94
1:A:137:GLU:HG3	1:C:137:GLU:HG3	1.50	0.92
1:A:57:ASN:HB3	1:A:60:ALA:HB3	1.53	0.91
1:A:133:MSE:HE2	1:A:137:GLU:HB3	1.55	0.85
2:E:7:UNK:HA	3:E:281:HOH:O	1.75	0.85
1:A:21:LYS:HG3	1:A:22:LEU:HG	1.60	0.83
1:B:22:LEU:HA	3:B:298:HOH:O	1.77	0.83
1:D:112:LYS:HE3	1:D:115:ARG:NH1	1.94	0.82
1:C:112:LYS:HE3	1:C:115:ARG:NH2	1.95	0.82
1:D:29:ASN:HD21	1:D:111:SER:HB2	1.44	0.81
1:B:149:LEU:N	1:B:150:PRO:HD2	1.96	0.80
1:D:54:MSE:HE1	1:D:64:VAL:HG11	1.63	0.80
1:B:23:ILE:HG22	1:B:27:GLN:HE22	1.44	0.80
1:D:153:LYS:HE3	1:D:156:VAL:HG21	1.65	0.79
1:B:26:LEU:HD21	1:B:178:VAL:CG1	2.11	0.78
1:B:26:LEU:HD21	1:B:178:VAL:HG13	1.67	0.77
1:B:86:LYS:HB3	3:B:358:HOH:O	1.84	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:PHE:O	1:A:109:LEU:HB2	1.85	0.77
1:B:137:GLU:HG3	1:D:137:GLU:HG3	1.67	0.77
1:C:152:LEU:HD23	1:C:154:LEU:HG	1.64	0.77
1:B:31:LEU:HD23	1:B:171:ILE:HG23	1.66	0.76
1:C:73:ARG:HG3	1:C:74:ALA:N	1.97	0.76
1:D:133:MSE:HE2	1:D:137:GLU:HB3	1.66	0.76
1:C:26:LEU:HD11	1:C:178:VAL:HG13	1.68	0.76
1:D:63:LYS:HG2	1:D:67:ASN:HD22	1.49	0.76
2:G:7:UNK:HA	3:G:251:HOH:O	1.86	0.76
1:A:71:ASP:OD1	1:A:73:ARG:HB3	1.85	0.75
1:D:112:LYS:HE3	1:D:115:ARG:CZ	2.17	0.75
1:A:48:HIS:NE2	1:A:96:MSE:HG2	2.02	0.74
1:B:31:LEU:CD2	1:B:171:ILE:HG23	2.17	0.74
1:B:153:LYS:HB3	1:B:156:VAL:HG21	1.70	0.74
1:D:29:ASN:ND2	1:D:111:SER:HB2	2.03	0.73
1:D:83:LYS:HB2	3:D:215:HOH:O	1.86	0.73
1:B:151:LEU:HB2	3:B:233:HOH:O	1.88	0.73
1:D:96:MSE:HG3	3:D:236:HOH:O	1.89	0.73
1:C:87:VAL:CG1	1:C:142:ILE:HG23	2.18	0.73
1:D:175:ASN:O	1:D:179:ARG:HB2	1.89	0.72
1:A:23:ILE:HA	1:A:26:LEU:HB2	1.71	0.72
1:B:147:THR:HG21	1:D:148:LYS:HE2	1.72	0.71
1:C:45:LYS:HD2	3:C:227:HOH:O	1.89	0.71
1:A:175:ASN:O	1:A:179:ARG:HB2	1.91	0.71
1:C:31:LEU:HG	1:C:171:ILE:HG23	1.73	0.71
1:C:21:LYS:HG3	1:C:22:LEU:HG	1.73	0.71
1:C:175:ASN:O	1:C:179:ARG:HB2	1.91	0.70
1:B:55:HIS:HA	1:B:85:LYS:HG3	1.72	0.70
1:D:48:HIS:NE2	1:D:96:MSE:HG2	2.07	0.70
1:D:63:LYS:HG2	1:D:67:ASN:ND2	2.06	0.70
1:B:133:MSE:HE2	1:B:137:GLU:HB3	1.72	0.70
1:D:141:GLN:NE2	3:D:208:HOH:O	2.24	0.70
1:B:175:ASN:O	1:B:179:ARG:HB2	1.90	0.69
1:B:30:GLU:O	1:B:111:SER:HB2	1.92	0.69
1:C:58:LYS:HZ1	1:C:81:VAL:HG21	1.56	0.69
1:C:26:LEU:O	1:C:26:LEU:HD23	1.92	0.69
1:D:104:GLU:OE2	1:D:162:MSE:HE1	1.92	0.69
1:D:54:MSE:CE	1:D:64:VAL:HG11	2.22	0.69
1:A:55:HIS:HA	1:A:85:LYS:HG3	1.75	0.69
1:A:54:MSE:HE1	1:A:64:VAL:HB	1.75	0.68
1:D:67:ASN:HA	3:D:241:HOH:O	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:133:MSE:HE2	1:C:137:GLU:HB3	1.76	0.68
1:A:105:ASN:ND2	1:A:162:MSE:HE3	2.09	0.67
1:A:57:ASN:HB3	1:A:60:ALA:CB	2.23	0.67
1:A:105:ASN:HD21	1:A:162:MSE:HG2	1.58	0.67
1:B:23:ILE:HG22	1:B:27:GLN:NE2	2.10	0.67
1:C:167:ASN:OD1	3:C:235:HOH:O	2.12	0.66
1:B:40:ILE:HD11	1:B:115:ARG:HG2	1.75	0.66
1:B:110:PHE:N	1:B:114:ASP:OD2	2.28	0.66
1:B:56:THR:HG22	1:B:56:THR:O	1.95	0.66
1:C:108:PHE:O	1:C:109:LEU:HB2	1.96	0.66
1:A:45:LYS:HE3	3:A:215:HOH:O	1.96	0.66
1:A:87:VAL:CG1	1:A:142:ILE:HG23	2.25	0.65
1:A:71:ASP:O	1:A:75:VAL:HG23	1.95	0.65
1:C:147:THR:HB	3:C:241:HOH:O	1.94	0.65
1:A:146:GLY:O	1:A:150:PRO:HG2	1.96	0.65
1:B:119:GLY:HA3	3:B:308:HOH:O	1.96	0.64
1:A:120:ILE:HD11	3:E:429:HOH:O	1.95	0.64
1:D:86:LYS:HB3	3:D:211:HOH:O	1.96	0.64
1:B:152:LEU:HG	1:B:154:LEU:HG	1.79	0.64
1:C:87:VAL:HG13	1:C:142:ILE:HG23	1.80	0.64
1:B:21:LYS:HG3	1:B:22:LEU:N	2.13	0.64
1:D:26:LEU:HD23	1:D:26:LEU:O	1.98	0.63
1:A:55:HIS:NE2	1:A:92:GLU:OE1	2.30	0.62
1:A:112:LYS:HE2	1:A:115:ARG:NH2	2.14	0.62
1:D:160:GLU:HB2	3:D:216:HOH:O	2.00	0.61
1:B:126:SER:OG	1:B:129:ILE:HG12	2.00	0.61
1:D:75:VAL:HG12	3:D:262:HOH:O	2.01	0.61
1:B:57:ASN:HB3	1:B:60:ALA:CB	2.31	0.61
1:C:179:ARG:HD3	3:C:212:HOH:O	2.01	0.61
1:C:136:ARG:NE	1:C:136:ARG:HA	2.16	0.61
1:A:48:HIS:CD2	1:A:96:MSE:HG2	2.36	0.60
1:B:86:LYS:HG2	3:B:234:HOH:O	2.00	0.60
1:A:87:VAL:HG13	1:A:142:ILE:HG23	1.83	0.60
1:C:61:LEU:O	1:C:65:LEU:HG	2.02	0.60
1:B:55:HIS:NE2	1:B:92:GLU:OE1	2.31	0.60
1:C:157:MSE:HG3	3:C:242:HOH:O	2.00	0.60
1:C:25:SER:HA	1:C:28:GLU:OE2	2.02	0.60
1:C:48:HIS:NE2	1:C:96:MSE:HG2	2.15	0.60
1:B:117:ILE:HG23	1:B:155:PHE:CZ	2.37	0.59
1:D:87:VAL:CG1	1:D:142:ILE:HG23	2.32	0.59
1:B:117:ILE:HG23	1:B:155:PHE:HZ	1.66	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:VAL:CG1	1:B:142:ILE:HG23	2.32	0.59
1:C:126:SER:OG	1:C:129:ILE:HG12	2.02	0.59
1:D:105:ASN:ND2	1:D:162:MSE:HE3	2.16	0.59
1:A:53:GLN:NE2	3:A:224:HOH:O	2.34	0.59
1:C:115:ARG:HD2	3:C:259:HOH:O	2.02	0.59
1:A:136:ARG:HA	1:A:136:ARG:HE	1.68	0.59
1:A:40:ILE:CD1	1:A:115:ARG:HG2	2.32	0.59
1:D:90:LEU:HD21	1:D:150:PRO:HD3	1.83	0.59
1:A:40:ILE:HD11	1:A:115:ARG:HG2	1.84	0.59
1:A:54:MSE:HE3	1:A:61:LEU:HD12	1.84	0.59
1:B:53:GLN:HG2	3:B:265:HOH:O	2.02	0.59
1:B:57:ASN:HB3	1:B:60:ALA:HB3	1.84	0.59
1:B:48:HIS:NE2	1:B:96:MSE:HG2	2.18	0.59
1:D:57:ASN:OD1	1:D:59:GLY:N	2.36	0.58
1:A:148:LYS:O	1:A:152:LEU:HB2	2.03	0.58
1:A:85:LYS:O	1:A:89:ARG:HG3	2.04	0.58
1:B:136:ARG:HA	1:B:136:ARG:NE	2.19	0.58
1:A:108:PHE:O	1:A:109:LEU:CB	2.51	0.58
1:A:136:ARG:NE	1:A:136:ARG:HA	2.18	0.58
1:B:40:ILE:CD1	1:B:115:ARG:HG2	2.32	0.58
1:D:136:ARG:HA	1:D:136:ARG:NE	2.19	0.58
1:D:87:VAL:HG13	1:D:142:ILE:HG23	1.85	0.58
1:D:85:LYS:HE2	3:D:243:HOH:O	2.04	0.58
1:C:70:LYS:HG3	1:C:131:PHE:CZ	2.39	0.58
1:A:105:ASN:HD21	1:A:162:MSE:HE3	1.67	0.57
1:C:30:GLU:HG2	1:C:112:LYS:HD2	1.84	0.57
1:D:58:LYS:HE3	3:D:247:HOH:O	2.04	0.57
1:B:87:VAL:HG13	1:B:142:ILE:HG23	1.86	0.57
1:B:85:LYS:O	1:B:89:ARG:HG3	2.04	0.57
1:C:94:GLU:HA	3:C:254:HOH:O	2.04	0.57
1:C:136:ARG:HA	1:C:136:ARG:HE	1.69	0.57
1:A:54:MSE:HE3	1:A:61:LEU:CD1	2.33	0.57
1:A:73:ARG:HG2	3:A:229:HOH:O	2.04	0.57
1:B:136:ARG:HE	1:B:136:ARG:HA	1.68	0.57
1:B:26:LEU:HD21	1:B:178:VAL:HG11	1.88	0.56
1:D:85:LYS:O	1:D:89:ARG:HG3	2.05	0.56
1:D:89:ARG:NH2	3:D:212:HOH:O	2.27	0.56
1:B:26:LEU:HB3	3:B:243:HOH:O	2.06	0.56
1:C:58:LYS:NZ	1:C:81:VAL:HG21	2.21	0.56
1:C:25:SER:O	1:C:29:ASN:HB3	2.06	0.56
1:B:136:ARG:NH2	3:B:235:HOH:O	2.33	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:6:MSE:O	2:E:7:UNK:C	2.54	0.56
1:D:83:LYS:HG2	1:D:142:ILE:HG21	1.88	0.56
2:H:6:MSE:O	2:H:7:UNK:C	2.54	0.56
1:B:21:LYS:N	3:B:257:HOH:O	2.38	0.55
1:C:152:LEU:CD2	1:C:154:LEU:HG	2.33	0.55
1:A:115:ARG:NH2	3:A:239:HOH:O	2.39	0.55
1:C:152:LEU:HG	3:G:439:HOH:O	2.06	0.55
1:B:21:LYS:HD2	3:B:232:HOH:O	2.07	0.55
1:C:30:GLU:CG	1:C:112:LYS:HD2	2.37	0.55
1:D:136:ARG:HA	1:D:136:ARG:HE	1.70	0.55
1:D:105:ASN:HD21	1:D:162:MSE:HG2	1.70	0.55
1:D:47:MSE:HG3	3:D:231:HOH:O	2.06	0.55
1:D:83:LYS:HE3	3:D:220:HOH:O	2.05	0.55
2:F:6:MSE:O	2:F:7:UNK:C	2.55	0.55
1:B:143:GLU:HG3	1:D:148:LYS:NZ	2.22	0.55
1:D:54:MSE:HE1	1:D:64:VAL:CG1	2.35	0.55
1:D:90:LEU:O	1:D:94:GLU:HG2	2.07	0.55
1:A:142:ILE:HB	3:A:218:HOH:O	2.07	0.55
1:C:28:GLU:O	1:C:28:GLU:HG2	2.07	0.55
1:D:48:HIS:CD2	1:D:96:MSE:HG2	2.42	0.55
1:A:25:SER:HG	1:A:113:PHE:HD1	1.55	0.54
1:C:149:LEU:N	1:C:150:PRO:HD2	2.21	0.54
1:A:66:ARG:NH1	1:A:72:TYR:CE2	2.76	0.54
1:A:73:ARG:HG2	1:A:73:ARG:HH21	1.71	0.54
1:A:112:LYS:HE2	1:A:115:ARG:HH22	1.70	0.54
2:G:6:MSE:O	2:G:7:UNK:C	2.56	0.54
1:C:90:LEU:O	1:C:94:GLU:HG2	2.08	0.54
1:C:112:LYS:HE3	1:C:115:ARG:HH22	1.72	0.54
1:C:54:MSE:HE1	1:C:64:VAL:HB	1.90	0.54
1:C:112:LYS:HE3	1:C:115:ARG:CZ	2.38	0.54
1:C:87:VAL:HG11	1:C:142:ILE:HG23	1.88	0.54
1:C:54:MSE:CE	1:C:61:LEU:HD12	2.38	0.53
1:C:70:LYS:NZ	1:C:131:PHE:CE2	2.76	0.53
1:A:45:LYS:CE	3:A:215:HOH:O	2.55	0.53
1:A:73:ARG:HG2	1:A:73:ARG:NH2	2.24	0.53
1:B:70:LYS:NZ	1:B:131:PHE:CE2	2.77	0.53
1:A:102:ILE:CD1	1:A:109:LEU:HD11	2.39	0.52
1:C:54:MSE:SE	1:C:64:VAL:HG21	2.58	0.52
1:C:85:LYS:O	1:C:89:ARG:HG3	2.09	0.52
1:D:27:GLN:HA	1:D:27:GLN:OE1	2.09	0.52
1:D:64:VAL:HG12	1:D:65:LEU:N	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:171:ILE:HD11	3:B:428:HOH:O	2.08	0.52
1:C:108:PHE:O	1:C:109:LEU:CB	2.58	0.52
1:D:40:ILE:HD11	1:D:115:ARG:HG2	1.92	0.52
1:A:32:LEU:HG	1:A:111:SER:C	2.31	0.51
1:A:90:LEU:O	1:A:94:GLU:HG2	2.11	0.51
1:B:29:ASN:ND2	1:B:111:SER:OG	2.43	0.51
1:B:143:GLU:CG	1:D:148:LYS:NZ	2.73	0.51
1:D:39:LYS:HA	3:D:253:HOH:O	2.09	0.51
1:A:30:GLU:O	1:A:111:SER:HB2	2.11	0.51
1:B:56:THR:O	1:B:57:ASN:HB2	2.09	0.51
1:C:107:ASN:HB2	3:C:211:HOH:O	2.09	0.51
1:D:119:GLY:HA3	3:D:249:HOH:O	2.09	0.51
1:B:148:LYS:C	1:B:150:PRO:HD2	2.30	0.51
1:B:154:LEU:O	1:B:155:PHE:HB2	2.11	0.51
1:C:54:MSE:HE3	1:C:61:LEU:HD12	1.93	0.51
1:B:21:LYS:CG	1:B:22:LEU:N	2.74	0.51
1:B:80:GLY:HA2	3:B:406:HOH:O	2.10	0.51
1:B:90:LEU:O	1:B:94:GLU:HG2	2.11	0.51
1:B:54:MSE:SE	1:B:61:LEU:HA	2.61	0.51
1:D:87:VAL:HG13	3:D:211:HOH:O	2.11	0.51
1:A:126:SER:OG	1:A:128:PRO:HD2	2.11	0.50
1:C:27:GLN:HG2	3:C:233:HOH:O	2.10	0.50
1:A:152:LEU:HG	1:A:154:LEU:HG	1.93	0.50
1:A:126:SER:OG	1:A:129:ILE:HG12	2.12	0.50
1:B:113:PHE:HA	3:B:382:HOH:O	2.11	0.50
1:C:54:MSE:HE3	1:C:61:LEU:CD1	2.42	0.50
1:B:25:SER:HG	1:B:113:PHE:HD1	1.59	0.50
1:B:123:PRO:HA	1:B:129:ILE:HG21	1.94	0.50
1:D:60:ALA:O	1:D:64:VAL:HG23	2.12	0.50
1:A:123:PRO:HA	1:A:129:ILE:HG21	1.94	0.49
1:B:147:THR:HG21	1:D:148:LYS:CE	2.42	0.49
1:B:71:ASP:O	1:B:75:VAL:HG23	2.12	0.49
1:C:70:LYS:HG3	1:C:131:PHE:CE1	2.47	0.49
1:C:146:GLY:O	1:C:150:PRO:HG3	2.13	0.49
2:H:4:UNK:C	2:H:6:MSE:H	2.24	0.49
1:A:114:ASP:OD1	1:A:174:TYR:OH	2.28	0.49
1:A:22:LEU:HD23	3:A:223:HOH:O	2.12	0.49
1:C:40:ILE:HD11	1:C:115:ARG:HG2	1.94	0.49
1:C:70:LYS:NZ	1:C:131:PHE:CD2	2.81	0.49
1:C:55:HIS:HA	1:C:85:LYS:HG3	1.95	0.49
1:B:48:HIS:CD2	1:B:96:MSE:HG2	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:26:LEU:HD11	1:C:178:VAL:CG1	2.40	0.49
1:C:97:GLU:HB2	3:C:254:HOH:O	2.12	0.49
1:B:143:GLU:CG	1:D:148:LYS:HZ2	2.27	0.48
1:D:164:PHE:CZ	1:D:168:LEU:HD21	2.47	0.48
1:D:63:LYS:NZ	1:D:67:ASN:ND2	2.61	0.48
1:A:152:LEU:O	1:A:153:LYS:C	2.51	0.48
1:B:23:ILE:CG2	1:B:27:GLN:HE22	2.20	0.48
1:D:126:SER:OG	1:D:129:ILE:HG12	2.14	0.48
1:D:26:LEU:HD21	1:D:178:VAL:HG13	1.94	0.48
1:C:30:GLU:HG2	1:C:112:LYS:CD	2.44	0.48
1:A:87:VAL:HG11	1:A:142:ILE:HG23	1.95	0.48
1:A:105:ASN:ND2	1:A:162:MSE:HG2	2.24	0.48
1:D:29:ASN:HD21	1:D:111:SER:CB	2.21	0.48
1:A:112:LYS:HG2	1:A:115:ARG:NH2	2.29	0.48
1:D:83:LYS:HD3	3:D:251:HOH:O	2.14	0.48
1:C:40:ILE:CD1	1:C:115:ARG:HG2	2.43	0.48
1:D:157:MSE:HE3	1:D:161:GLU:HB3	1.96	0.48
1:B:109:LEU:HD23	1:B:114:ASP:HB3	1.95	0.48
1:B:174:TYR:O	1:B:178:VAL:HG12	2.13	0.48
1:C:114:ASP:OD1	1:C:174:TYR:OH	2.30	0.48
1:C:152:LEU:CG	3:G:439:HOH:O	2.60	0.48
1:D:26:LEU:HD11	1:D:178:VAL:HG22	1.94	0.48
1:D:21:LYS:HE3	1:D:22:LEU:CD1	2.44	0.48
1:D:102:ILE:CD1	1:D:109:LEU:HD11	2.43	0.47
1:D:70:LYS:HG2	1:D:70:LYS:O	2.13	0.47
1:C:83:LYS:HG2	1:C:142:ILE:HG21	1.96	0.47
1:B:57:ASN:O	1:B:60:ALA:HB3	2.14	0.47
1:C:48:HIS:CD2	1:C:96:MSE:HG2	2.50	0.47
1:D:104:GLU:HG3	1:D:162:MSE:HE3	1.97	0.47
1:B:98:LEU:CD1	1:B:108:PHE:HE2	2.28	0.47
1:A:58:LYS:HE2	1:A:85:LYS:NZ	2.30	0.47
1:B:87:VAL:N	3:B:234:HOH:O	2.47	0.47
1:C:100:HIS:O	1:C:103:ALA:HB3	2.14	0.47
1:A:70:LYS:HA	1:A:131:PHE:CE1	2.50	0.47
2:F:1:UNK:N	3:F:209:HOH:O	2.32	0.47
2:F:4:UNK:C	2:F:6:MSE:H	2.27	0.47
1:A:105:ASN:HD21	1:A:162:MSE:CG	2.25	0.47
1:A:73:ARG:O	1:A:76:ILE:HG12	2.15	0.47
1:A:81:VAL:HG22	3:A:210:HOH:O	2.14	0.47
1:B:71:ASP:OD1	1:B:74:ALA:CB	2.62	0.47
1:D:174:TYR:O	1:D:178:VAL:HG12	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:126:SER:OG	1:D:128:PRO:HD2	2.15	0.46
1:B:70:LYS:HE2	1:B:131:PHE:O	2.15	0.46
2:E:6:MSE:HG3	3:E:286:HOH:O	2.15	0.46
2:G:4:UNK:C	2:G:6:MSE:H	2.29	0.46
2:H:1:UNK:N	3:H:211:HOH:O	2.36	0.46
1:B:29:ASN:HD22	1:B:111:SER:CB	2.29	0.46
3:A:222:HOH:O	1:C:148:LYS:HG3	2.15	0.46
1:C:28:GLU:O	1:C:28:GLU:CG	2.63	0.46
1:B:126:SER:OG	1:B:128:PRO:HD2	2.16	0.46
1:B:30:GLU:O	1:B:111:SER:CB	2.62	0.46
1:D:54:MSE:SE	1:D:64:VAL:HG21	2.65	0.46
1:D:55:HIS:NE2	1:D:92:GLU:OE1	2.48	0.46
1:D:26:LEU:HD11	1:D:178:VAL:HG13	1.98	0.46
2:E:3:UNK:HA	3:E:429:HOH:O	2.15	0.46
1:B:109:LEU:HA	1:B:114:ASP:CB	2.46	0.46
1:B:83:LYS:HG2	1:B:142:ILE:HG21	1.98	0.46
1:B:105:ASN:HD21	1:B:162:MSE:HG2	1.81	0.46
1:C:152:LEU:HD11	3:G:439:HOH:O	2.16	0.46
1:C:31:LEU:HD21	1:C:110:PHE:CZ	2.52	0.45
1:C:21:LYS:HG3	1:C:22:LEU:N	2.31	0.45
1:C:33:ASN:ND2	1:C:36:GLU:CD	2.70	0.45
1:A:149:LEU:HG	3:A:217:HOH:O	2.15	0.45
1:B:71:ASP:OD1	1:B:74:ALA:HB2	2.16	0.45
1:B:93:SER:O	1:B:96:MSE:HB2	2.16	0.45
1:A:164:PHE:CZ	1:A:168:LEU:HD21	2.52	0.45
1:B:98:LEU:HB2	1:B:155:PHE:CD2	2.51	0.45
1:B:73:ARG:HA	1:B:76:ILE:HG12	1.98	0.45
1:D:21:LYS:O	1:D:22:LEU:HD23	2.17	0.45
1:A:174:TYR:O	1:A:178:VAL:HG12	2.16	0.45
2:E:4:UNK:C	2:E:6:MSE:H	2.28	0.45
1:A:23:ILE:O	1:A:27:GLN:HB2	2.16	0.45
1:C:71:ASP:CG	1:C:74:ALA:HB3	2.36	0.45
1:C:104:GLU:O	1:C:104:GLU:HG3	2.17	0.45
1:C:85:LYS:NZ	3:C:228:HOH:O	2.49	0.45
1:B:57:ASN:OD1	1:B:60:ALA:N	2.50	0.45
1:A:76:ILE:CD1	3:A:227:HOH:O	2.64	0.45
1:B:173:GLN:HG2	3:B:275:HOH:O	2.17	0.45
1:D:173:GLN:HA	3:D:258:HOH:O	2.16	0.45
1:A:137:GLU:HG3	1:C:137:GLU:CG	2.34	0.44
1:C:157:MSE:HE3	1:C:161:GLU:HB3	1.99	0.44
1:C:28:GLU:O	3:C:230:HOH:O	2.21	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:55:HIS:ND1	1:D:55:HIS:O	2.51	0.44
1:A:167:ASN:O	1:A:169:LYS:HG2	2.18	0.44
1:C:164:PHE:CZ	1:C:168:LEU:HD21	2.52	0.44
1:D:105:ASN:HD21	1:D:162:MSE:HE3	1.83	0.44
1:A:104:GLU:HG3	1:A:162:MSE:HE3	2.00	0.44
1:A:167:ASN:O	1:A:168:LEU:C	2.55	0.44
1:D:105:ASN:ND2	1:D:162:MSE:HG2	2.32	0.44
1:D:33:ASN:ND2	1:D:36:GLU:CD	2.71	0.44
1:C:154:LEU:HD21	2:G:2:UNK:C	2.48	0.44
1:B:105:ASN:HD21	1:B:162:MSE:HE3	1.83	0.44
2:G:2:UNK:CB	3:G:439:HOH:O	2.66	0.44
1:C:67:ASN:HB3	3:C:246:HOH:O	2.17	0.44
1:D:71:ASP:O	1:D:71:ASP:OD1	2.35	0.44
1:D:73:ARG:HG3	1:D:73:ARG:NH2	2.33	0.44
1:B:111:SER:C	1:B:113:PHE:N	2.71	0.43
1:B:21:LYS:HG3	1:B:22:LEU:H	1.83	0.43
1:B:136:ARG:HE	1:B:136:ARG:CA	2.30	0.43
1:A:21:LYS:HG3	1:A:22:LEU:N	2.33	0.43
1:C:136:ARG:HE	1:C:136:ARG:CA	2.30	0.43
1:C:167:ASN:O	1:C:168:LEU:C	2.56	0.43
1:A:50:PHE:O	1:A:54:MSE:HG2	2.18	0.43
1:B:84:PHE:O	1:B:87:VAL:HG22	2.18	0.43
1:C:87:VAL:HG23	1:C:88:TYR:N	2.33	0.43
1:D:26:LEU:HD21	1:D:178:VAL:CG1	2.48	0.43
1:A:133:MSE:HE3	3:A:213:HOH:O	2.18	0.43
1:B:164:PHE:CZ	1:B:168:LEU:HD21	2.53	0.43
1:D:168:LEU:N	3:D:226:HOH:O	2.51	0.43
1:A:29:ASN:HD22	1:A:111:SER:CB	2.32	0.43
1:B:167:ASN:O	1:B:168:LEU:C	2.56	0.43
1:C:174:TYR:O	1:C:178:VAL:HG12	2.18	0.43
1:D:85:LYS:HB2	1:D:85:LYS:HE3	1.86	0.43
1:C:43:GLN:O	1:C:47:MSE:HG3	2.19	0.43
1:D:158:THR:O	1:D:162:MSE:HG3	2.19	0.43
1:D:167:ASN:O	1:D:168:LEU:C	2.57	0.43
1:D:167:ASN:O	1:D:169:LYS:HG2	2.17	0.43
1:D:33:ASN:OD1	1:D:36:GLU:HG3	2.19	0.43
1:A:105:ASN:HD21	1:A:162:MSE:CE	2.31	0.43
1:B:55:HIS:CG	1:B:89:ARG:HG2	2.54	0.43
2:G:2:UNK:CA	3:G:439:HOH:O	2.67	0.43
1:C:48:HIS:HE1	3:C:239:HOH:O	2.01	0.42
1:C:71:ASP:OD1	1:C:74:ALA:HB3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:85:LYS:HB2	1:C:85:LYS:HE3	1.80	0.42
1:D:88:TYR:HA	1:D:91:LEU:HB2	2.01	0.42
1:B:21:LYS:CG	1:B:22:LEU:H	2.32	0.42
1:C:105:ASN:N	1:C:106:PRO:HD3	2.34	0.42
1:A:157:MSE:HE3	1:A:161:GLU:HB3	2.00	0.42
1:B:113:PHE:CE2	1:B:178:VAL:HG21	2.54	0.42
1:D:114:ASP:C	1:D:116:SER:H	2.22	0.42
1:D:136:ARG:CA	1:D:136:ARG:HE	2.33	0.42
1:D:114:ASP:C	1:D:116:SER:N	2.73	0.42
1:D:40:ILE:HD13	3:D:268:HOH:O	2.19	0.42
1:D:48:HIS:CD2	1:D:95:THR:OG1	2.73	0.42
1:A:21:LYS:NZ	3:A:245:HOH:O	2.46	0.42
1:C:93:SER:O	1:C:96:MSE:HB2	2.19	0.42
1:D:108:PHE:O	1:D:109:LEU:HB2	2.20	0.42
1:D:45:LYS:NZ	3:D:254:HOH:O	2.52	0.42
1:A:130:MSE:HE3	3:A:219:HOH:O	2.19	0.42
1:C:114:ASP:C	1:C:116:SER:N	2.73	0.42
1:D:82:ASP:HB3	3:D:228:HOH:O	2.19	0.42
1:A:138:MSE:O	1:A:142:ILE:HG13	2.19	0.42
1:B:56:THR:CG2	1:B:56:THR:O	2.66	0.42
1:C:75:VAL:HG22	1:C:135:ILE:CD1	2.49	0.42
1:A:171:ILE:HD11	3:A:216:HOH:O	2.20	0.41
1:B:167:ASN:O	1:B:169:LYS:HG2	2.19	0.41
1:D:40:ILE:CD1	1:D:115:ARG:HG2	2.50	0.41
1:D:154:LEU:HD23	2:H:2:UNK:CB	2.50	0.41
1:A:114:ASP:C	1:A:116:SER:H	2.22	0.41
1:A:136:ARG:CA	1:A:136:ARG:HE	2.31	0.41
1:B:73:ARG:HG3	1:B:73:ARG:HH21	1.85	0.41
1:D:48:HIS:HD2	1:D:95:THR:OG1	2.04	0.41
1:A:114:ASP:C	1:A:116:SER:N	2.73	0.41
1:B:88:TYR:HA	1:B:91:LEU:HB2	2.02	0.41
1:C:136:ARG:NE	1:C:136:ARG:CA	2.83	0.41
1:A:54:MSE:SE	1:A:64:VAL:HG21	2.71	0.41
1:C:71:ASP:OD1	1:C:71:ASP:C	2.59	0.41
1:C:114:ASP:C	1:C:116:SER:H	2.24	0.41
1:A:29:ASN:ND2	1:A:111:SER:CB	2.83	0.41
1:A:127:LYS:HB3	1:A:128:PRO:HD3	2.03	0.41
1:C:73:ARG:CG	1:C:74:ALA:N	2.78	0.41
1:D:108:PHE:O	1:D:109:LEU:CB	2.69	0.41
1:A:83:LYS:HG2	1:A:142:ILE:HG21	2.03	0.41
1:B:114:ASP:C	1:B:116:SER:N	2.74	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:21:LYS:HE3	1:B:22:LEU:HD12	2.03	0.41
1:D:62:ASP:O	1:D:63:LYS:C	2.58	0.41
1:D:58:LYS:NZ	1:D:81:VAL:HG21	2.35	0.41
1:D:107:ASN:OD1	1:D:165:TYR:O	2.39	0.41
1:B:57:ASN:HB3	1:B:60:ALA:HB2	2.01	0.41
1:C:36:GLU:O	1:C:40:ILE:HG12	2.20	0.41
1:D:123:PRO:HA	1:D:129:ILE:HG21	2.03	0.41
1:D:154:LEU:HD21	2:H:2:UNK:C	2.51	0.41
1:D:28:GLU:O	1:D:28:GLU:CD	2.58	0.40
1:A:76:ILE:HD13	3:A:227:HOH:O	2.20	0.40
1:C:179:ARG:NH1	3:C:212:HOH:O	2.54	0.40
1:D:23:ILE:O	1:D:27:GLN:N	2.54	0.40
1:A:23:ILE:O	1:A:23:ILE:HG22	2.22	0.40
1:B:137:GLU:HG3	1:D:137:GLU:CG	2.45	0.40
1:B:31:LEU:HG	3:B:239:HOH:O	2.21	0.40
1:C:148:LYS:O	1:C:152:LEU:HB2	2.22	0.40
1:C:71:ASP:OD1	1:C:74:ALA:CB	2.69	0.40
1:C:88:TYR:HA	1:C:91:LEU:HB2	2.04	0.40
1:A:104:GLU:HG3	1:A:162:MSE:CE	2.52	0.40
1:A:29:ASN:HD22	1:A:111:SER:HB2	1.86	0.40
1:A:54:MSE:CE	1:A:61:LEU:HD12	2.51	0.40
1:C:50:PHE:O	1:C:54:MSE:HG2	2.22	0.40
1:D:104:GLU:HG2	3:D:222:HOH:O	2.20	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	159/208 (76%)	145 (91%)	11 (7%)	3 (2%)	8 7
1	B	159/208 (76%)	140 (88%)	16 (10%)	3 (2%)	8 7

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	C	159/208 (76%)	149 (94%)	9 (6%)	1 (1%)	25 31
1	D	159/208 (76%)	141 (89%)	16 (10%)	2 (1%)	12 12
2	E	1/7 (14%)	0	1 (100%)	0	100 100
2	F	1/7 (14%)	0	1 (100%)	0	100 100
2	G	1/7 (14%)	0	1 (100%)	0	100 100
2	H	1/7 (14%)	0	1 (100%)	0	100 100
All	All	640/860 (74%)	575 (90%)	56 (9%)	9 (1%)	11 11

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	109	LEU
1	A	109	LEU
1	D	64	VAL
1	B	57	ASN
1	B	150	PRO
1	B	152	LEU
1	D	27	GLN
1	A	153	LYS
1	A	73	ARG

5.3.2 Protein sidechains (i)

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	152/182 (84%)	144 (95%)	8 (5%)	22 31
1	B	152/182 (84%)	142 (93%)	10 (7%)	16 22
1	C	152/182 (84%)	141 (93%)	11 (7%)	14 18
1	D	152/182 (84%)	144 (95%)	8 (5%)	22 31
2	E	-	1 (100%)	0	100 100
2	F	-	1 (100%)	0	100 100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	G	-	1 (100%)	0	100 100
2	H	-	1 (100%)	0	100 100
All	All	612/728 (84%)	575 (94%)	37 (6%)	19 26

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	28	GLU
1	A	82	ASP
1	A	99	LEU
1	A	122	LEU
1	A	126	SER
1	A	149	LEU
1	A	171	ILE
1	B	27	GLN
1	B	62	ASP
1	B	73	ARG
1	B	82	ASP
1	B	91	LEU
1	B	99	LEU
1	B	122	LEU
1	B	126	SER
1	B	149	LEU
1	B	171	ILE
1	C	28	GLU
1	C	56	THR
1	C	82	ASP
1	C	99	LEU
1	C	106	PRO
1	C	109	LEU
1	C	118	LEU
1	C	122	LEU
1	C	126	SER
1	C	149	LEU
1	C	171	ILE
1	D	28	GLU
1	D	82	ASP
1	D	99	LEU
1	D	107	ASN
1	D	122	LEU

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Mol	Chain	Res	Type
1	D	126	SER
1	D	149	LEU
1	D	171	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	ASN
1	A	105	ASN
1	A	107	ASN
1	B	27	GLN
1	B	29	ASN
1	B	48	HIS
1	B	105	ASN
1	B	107	ASN
1	D	29	ASN
1	D	48	HIS
1	D	67	ASN
1	D	105	ASN
1	D	107	ASN

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

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^{th} 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	152/208 (73%)	0.77	13 (8%) 10 14	1, 7, 21, 36	0
1	B	152/208 (73%)	0.86	15 (9%) 7 10	1, 8, 23, 35	0
1	C	152/208 (73%)	0.89	17 (11%) 5 7	1, 8, 21, 37	0
1	D	152/208 (73%)	1.14	21 (13%) 2 4	1, 9, 22, 36	0
2	E	0/7	-	-	-	-
2	F	0/7	-	-	-	-
2	G	0/7	-	-	-	-
2	H	0/7	-	-	-	-
All	All	608/860 (70%)	0.92	66 (10%) 5 8	1, 8, 23, 37	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	35	ASP	7.7
1	C	23	ILE	6.1
1	A	26	LEU	5.2
1	D	25	SER	5.0
1	C	151	LEU	4.6
1	A	22	LEU	4.3
1	C	179	ARG	4.2
1	D	168	LEU	3.9
1	D	178	VAL	3.9
1	C	25	SER	3.9
1	D	34	THR	3.5
1	D	33	ASN	3.4
1	B	77	LYS	3.3
1	D	26	LEU	3.3
1	D	100	HIS	3.2
1	D	179	ARG	3.2
1	C	33	ASN	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	160	GLU	3.1
1	B	82	ASP	3.0
1	D	109	LEU	3.0
1	C	26	LEU	3.0
1	D	22	LEU	2.9
1	B	178	VAL	2.8
1	D	39	LYS	2.8
1	B	152	LEU	2.8
1	C	171	ILE	2.8
1	D	180	ASP	2.7
1	B	158	THR	2.7
1	C	173	GLN	2.7
1	A	160	GLU	2.7
1	C	178	VAL	2.6
1	D	173	GLN	2.6
1	B	181	LEU	2.6
1	D	82	ASP	2.5
1	B	25	SER	2.5
1	D	71	ASP	2.5
1	D	72	TYR	2.4
1	B	154	LEU	2.4
1	A	29	ASN	2.4
1	A	82	ASP	2.4
1	D	43	GLN	2.4
1	A	181	LEU	2.3
1	D	32	LEU	2.3
1	A	151	LEU	2.3
1	A	178	VAL	2.3
1	B	57	ASN	2.3
1	C	70	LYS	2.2
1	B	28	GLU	2.2
1	B	58	LYS	2.2
1	A	159	ASP	2.2
1	C	43	GLN	2.2
1	D	151	LEU	2.2
1	A	153	LYS	2.2
1	C	22	LEU	2.2
1	A	180	ASP	2.2
1	B	153	LYS	2.2
1	C	67	ASN	2.1
1	A	163	ASN	2.1
1	B	29	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	22	LEU	2.1
1	C	136	ARG	2.0
1	C	177	TYR	2.0
1	A	27	GLN	2.0
1	D	86	LYS	2.0
1	B	109	LEU	2.0
1	C	90	LEU	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

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.