Full wwPDB X-ray Structure Validation Report (i)

Jun 19, 2024 – 02:07 PM EDT

PDB ID	:	3WBK
Title	:	crystal structure analysis of eukaryotic translation initiation factor 5B and 1A complex
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Deposited on	:	2013-05-20
Resolution	:	3.30 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 (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

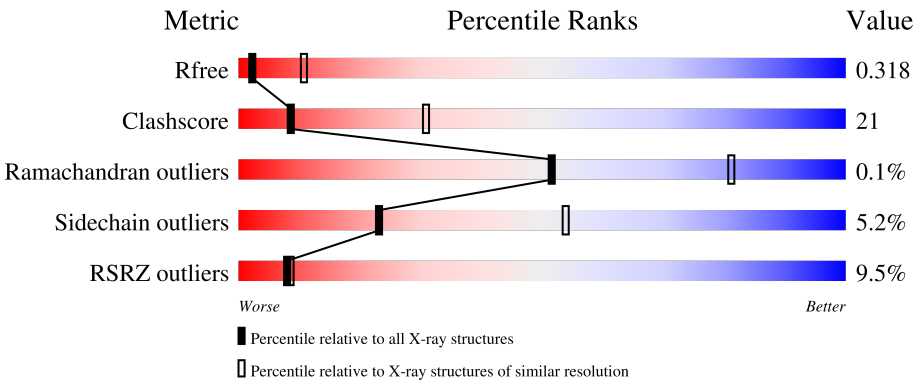
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 3.30 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	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	606	6% 65% 30%
1	B	606	13% 62% 33%
2	C	131	% 6% 92%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 9309 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 Eukaryotic translation initiation factor 5B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	590	4588	2918	778	872	20	0	0	0
1	B	590	4634	2949	785	879	21	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP P39730
A	-2	SER	-	expression tag	UNP P39730
A	-1	HIS	-	expression tag	UNP P39730
A	0	MET	-	expression tag	UNP P39730
B	-3	GLY	-	expression tag	UNP P39730
B	-2	SER	-	expression tag	UNP P39730
B	-1	HIS	-	expression tag	UNP P39730
B	0	MET	-	expression tag	UNP P39730

- Molecule 2 is a protein called Eukaryotic translation initiation factor 1A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	C	11	87	51	11	25	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

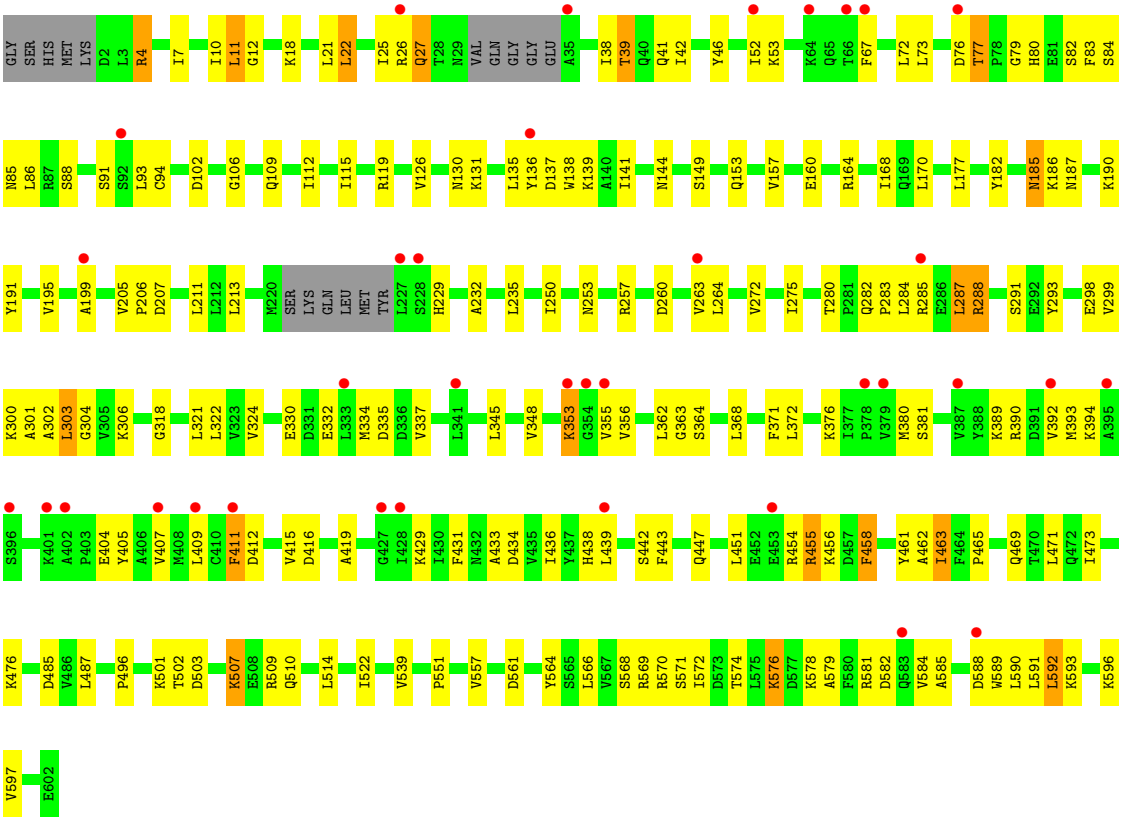
Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	GLY	-	expression tag	UNP P38912
C	-2	SER	-	expression tag	UNP P38912
C	-1	HIS	-	expression tag	UNP P38912
C	0	MET	-	expression tag	UNP P38912

3 Residue-property plots (i)

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

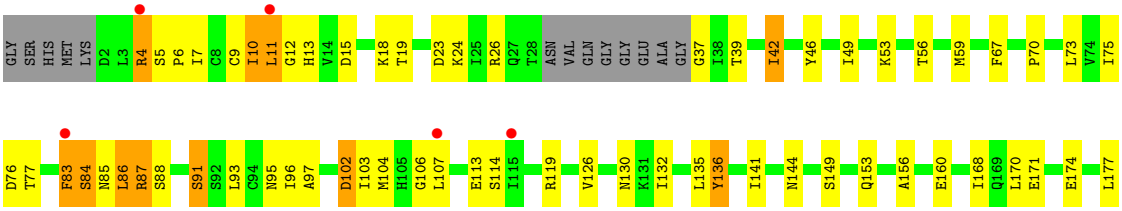
• Molecule 1: Eukaryotic translation initiation factor 5B

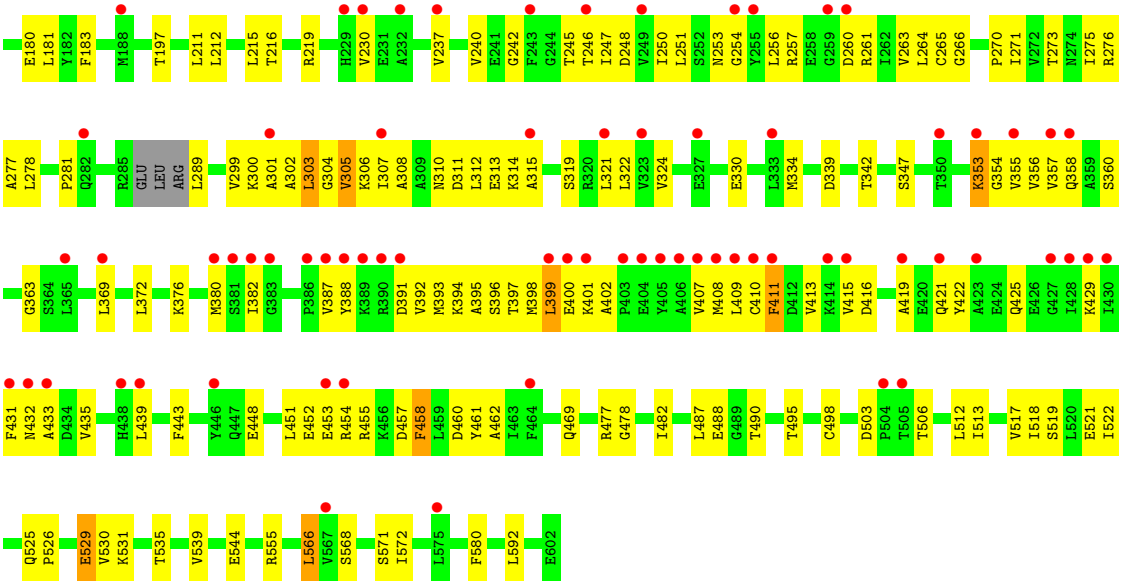
Chain A: 6% 65% 30%



• Molecule 1: Eukaryotic translation initiation factor 5B

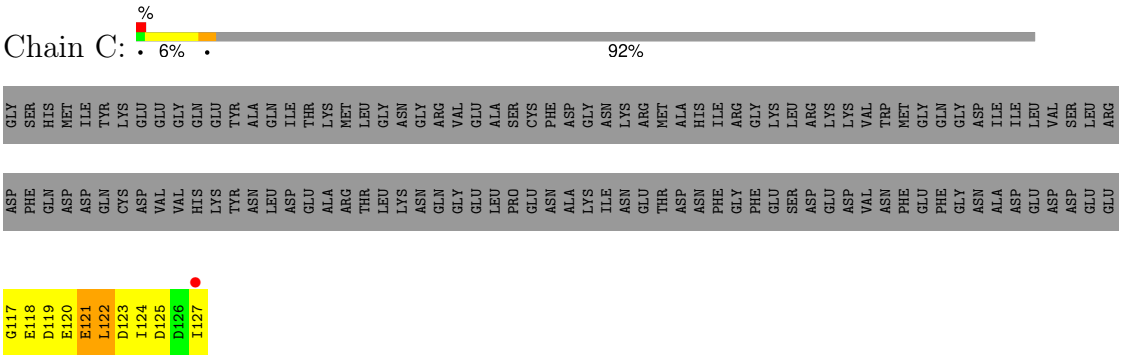
Chain B: 13% 62% 33%





- Molecule 2: Eukaryotic translation initiation factor 1A

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	101.94Å 120.94Å 132.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 3.30 44.70 – 3.30	Depositor EDS
% Data completeness (in resolution range)	98.2 (40.00-3.30) 98.4 (44.70-3.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.84 (at 3.32Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.257 , 0.317 0.254 , 0.318	Depositor DCC
R_{free} test set	1769 reflections (7.13%)	wwPDB-VP
Wilson B-factor (Å^2)	112.3	Xtriage
Anisotropy	0.667	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 97.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	9309	wwPDB-VP
Average B, all atoms (Å^2)	159.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% of the height of the origin peak. No significant pseudotranslation is detected.*

[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 (i)

5.1 Standard geometry (i)

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	0.43	0/4655	0.67	0/6295
1	B	0.41	0/4704	0.65	1/6356 (0.0%)
2	C	0.34	0/86	0.55	0/115
All	All	0.42	0/9445	0.66	1/12766 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	102	ASP	CB-CG-OD1	5.05	122.85	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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	4588	0	4694	175	0
1	B	4634	0	4766	207	0
2	C	87	0	69	21	0
All	All	9309	0	9529	389	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 21.

All (389) 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)
2:C:120:GLU:HB3	2:C:121:GLU:HB2	1.36	1.07
1:A:287:LEU:HD12	1:A:291:SER:HB3	1.31	1.07
1:B:9:CYS:HB3	1:B:83:PHE:CE1	1.90	1.06
1:B:11:LEU:HB3	1:B:83:PHE:CE2	1.95	1.02
1:A:86:LEU:HA	1:A:363:GLY:HA3	1.44	0.97
1:B:11:LEU:HB3	1:B:83:PHE:HE2	1.28	0.96
1:A:287:LEU:CD1	1:A:291:SER:HB3	1.96	0.95
1:B:299:VAL:HB	1:B:303:LEU:HD21	1.54	0.89
1:B:9:CYS:HB2	1:B:87:ARG:NH1	1.88	0.89
1:B:9:CYS:HB3	1:B:83:PHE:HE1	1.33	0.89
1:B:83:PHE:CE1	1:B:87:ARG:NH1	2.41	0.88
1:B:275:ILE:HG23	1:B:307:ILE:HG23	1.56	0.87
1:A:461:TYR:HB3	1:A:574:THR:HG21	1.57	0.86
2:C:120:GLU:HB3	2:C:121:GLU:CB	2.06	0.85
1:B:77:THR:HG21	1:B:83:PHE:CG	2.12	0.85
1:B:83:PHE:CZ	1:B:87:ARG:NH1	2.48	0.81
1:A:501:LYS:O	1:A:510:GLN:HB2	1.81	0.81
1:B:9:CYS:HB3	1:B:83:PHE:CZ	2.15	0.81
1:A:232:ALA:HB3	1:A:321:LEU:HB3	1.64	0.79
1:B:85:ASN:HA	1:B:363:GLY:HA3	1.64	0.79
1:B:356:VAL:HG13	1:B:408:MET:HA	1.63	0.79
1:A:42:ILE:HD13	1:A:306:LYS:HB2	1.65	0.78
1:B:245:THR:HB	1:B:312:LEU:HA	1.67	0.76
1:A:431:PHE:CZ	1:A:443:PHE:HA	2.20	0.76
1:A:455:ARG:HA	1:A:458:PHE:HD2	1.52	0.75
1:B:300:LYS:O	1:B:303:LEU:HD23	1.86	0.75
1:A:300:LYS:HG2	1:A:301:ALA:H	1.51	0.75
1:A:11:LEU:HD22	1:A:11:LEU:O	1.87	0.75
1:A:287:LEU:HD12	1:A:287:LEU:O	1.85	0.75
1:B:369:LEU:HD21	1:B:382:ILE:HG21	1.69	0.74
1:A:284:LEU:HB2	1:A:287:LEU:HB2	1.70	0.74
1:B:407:VAL:HG12	1:B:429:LYS:HB2	1.69	0.74
1:B:522:ILE:HG12	1:B:539:VAL:HG22	1.68	0.74
1:A:300:LYS:HG2	1:A:301:ALA:N	2.03	0.74
1:A:407:VAL:HG12	1:A:429:LYS:HB3	1.69	0.73
1:B:568:SER:H	1:B:571:SER:HB3	1.53	0.73
1:A:501:LYS:O	1:A:510:GLN:CB	2.37	0.73
1:A:11:LEU:HD13	1:A:11:LEU:N	2.05	0.72
1:A:431:PHE:HB3	1:A:439:LEU:HG	1.71	0.72
1:B:77:THR:HG21	1:B:83:PHE:CD1	2.25	0.71
1:B:9:CYS:CB	1:B:87:ARG:HH11	2.02	0.71
1:B:11:LEU:HD22	1:B:11:LEU:O	1.90	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:576:LYS:NZ	2:C:127:ILE:C	2.45	0.70
1:A:285:ARG:O	1:A:288:ARG:HG3	1.92	0.69
1:B:314:LYS:HD2	1:B:314:LYS:O	1.92	0.69
1:A:86:LEU:HA	1:A:363:GLY:CA	2.19	0.69
1:B:458:PHE:HA	1:B:461:TYR:HB2	1.73	0.69
1:A:42:ILE:HG22	1:A:304:GLY:O	1.92	0.69
1:B:126:VAL:HG21	1:B:211:LEU:HD23	1.73	0.69
1:A:263:VAL:HG12	1:A:272:VAL:HG12	1.74	0.69
1:A:284:LEU:HD22	1:A:287:LEU:HD22	1.74	0.68
1:A:514:LEU:HD21	1:A:557:VAL:HG21	1.75	0.68
1:A:187:ASN:O	1:A:191:TYR:HB2	1.93	0.68
1:A:284:LEU:HB2	1:A:287:LEU:CB	2.23	0.68
1:B:11:LEU:HD22	1:B:11:LEU:C	2.14	0.68
1:B:245:THR:CB	1:B:312:LEU:HA	2.23	0.67
1:B:83:PHE:O	1:B:87:ARG:HG3	1.95	0.67
2:C:124:ILE:HG12	2:C:124:ILE:O	1.94	0.67
1:B:237:VAL:HG22	1:B:247:ILE:HG22	1.75	0.67
1:B:357:VAL:HG11	1:B:372:LEU:HD11	1.77	0.67
1:B:339:ASP:HA	1:B:342:THR:HB	1.76	0.66
1:A:7:ILE:HG12	1:A:73:LEU:HD23	1.75	0.66
1:A:284:LEU:CB	1:A:287:LEU:HB2	2.25	0.66
1:A:451:LEU:HB3	1:A:578:LYS:HE2	1.77	0.66
1:A:514:LEU:HD23	1:A:551:PRO:HG2	1.78	0.66
1:B:353:LYS:HD2	1:B:402:ALA:HB2	1.78	0.66
1:B:9:CYS:CB	1:B:83:PHE:CE1	2.75	0.65
1:B:411:PHE:HB3	1:B:433:ALA:O	1.96	0.65
1:A:11:LEU:HD13	1:A:11:LEU:H	1.61	0.65
1:B:42:ILE:HG13	1:B:281:PRO:HD3	1.77	0.65
1:A:53:LYS:HG2	1:A:67:PHE:CZ	2.31	0.65
1:A:576:LYS:HG3	2:C:124:ILE:HD11	1.77	0.65
1:A:260:ASP:HB2	1:A:275:ILE:HD12	1.79	0.65
1:A:568:SER:H	1:A:571:SER:HB3	1.62	0.65
1:A:102:ASP:OD1	1:A:131:LYS:HD3	1.97	0.64
1:A:303:LEU:HD12	1:A:303:LEU:O	1.98	0.64
1:B:9:CYS:CB	1:B:87:ARG:NH1	2.58	0.64
1:B:119:ARG:HD2	1:B:177:LEU:HD21	1.80	0.64
1:A:284:LEU:HD22	1:A:287:LEU:CD2	2.28	0.64
1:A:38:ILE:HD12	1:A:38:ILE:O	1.98	0.64
1:A:85:ASN:O	1:A:88:SER:HB3	1.97	0.64
1:B:23:ASP:HA	1:B:26:ARG:NH1	2.12	0.64
1:B:11:LEU:O	1:B:11:LEU:HD13	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:LYS:CG	1:A:301:ALA:H	2.10	0.63
1:B:93:LEU:HD11	1:B:250:ILE:HG21	1.80	0.63
1:B:263:VAL:HG22	1:B:322:LEU:HB2	1.80	0.63
1:B:498:CYS:HB2	1:B:512:LEU:O	1.99	0.63
1:B:9:CYS:CB	1:B:83:PHE:HE1	2.09	0.62
1:A:502:THR:HG22	1:A:509:ARG:HA	1.80	0.62
1:A:318:GLY:C	1:A:362:LEU:HD12	2.19	0.62
1:B:330:GLU:O	1:B:334:MET:HG3	1.99	0.62
1:A:593:LYS:O	1:A:596:LYS:HG2	1.99	0.62
2:C:119:ASP:OD1	2:C:119:ASP:N	2.29	0.62
1:A:79:GLY:O	1:A:82:SER:HB3	2.00	0.62
1:B:11:LEU:CB	1:B:83:PHE:HE2	2.08	0.62
1:B:452:GLU:O	1:B:455:ARG:HB2	1.99	0.62
1:A:282:GLN:CD	1:A:283:PRO:HD2	2.20	0.61
1:B:39:THR:OG1	1:B:46:TYR:HB3	2.01	0.61
1:A:590:LEU:HD12	2:C:122:LEU:HD11	1.83	0.61
1:A:590:LEU:CD1	2:C:122:LEU:HD11	2.30	0.61
1:B:277:ALA:HB3	1:B:308:ALA:HB3	1.83	0.61
2:C:118:GLU:O	2:C:118:GLU:HG2	2.01	0.61
1:A:451:LEU:HG	1:A:578:LYS:HD3	1.83	0.61
1:B:522:ILE:O	1:B:525:GLN:HB2	2.01	0.60
1:B:357:VAL:CG2	1:B:382:ILE:HG22	2.31	0.60
1:A:503:ASP:HB2	1:A:507:LYS:H	1.66	0.60
1:A:588:ASP:O	1:A:592:LEU:HD12	2.02	0.60
1:A:160:GLU:O	1:A:164:ARG:HG3	2.02	0.60
1:A:149:SER:O	1:A:153:GLN:HG2	2.02	0.60
1:B:271:ILE:CD1	1:B:313:GLU:HG2	2.31	0.60
1:A:364:SER:HB2	1:A:411:PHE:CD2	2.38	0.59
1:A:280:THR:O	1:A:293:TYR:HB3	2.02	0.59
1:A:411:PHE:CE1	1:A:412:ASP:HB2	2.37	0.59
1:B:230:VAL:HA	1:B:254:GLY:HA3	1.85	0.59
1:B:360:SER:HA	1:B:413:VAL:HG13	1.84	0.59
1:B:462:ALA:HB2	1:B:580:PHE:CE1	2.37	0.59
1:A:465:PRO:HB2	1:A:591:LEU:HD23	1.83	0.59
1:A:569:ARG:HH22	2:C:127:ILE:CG1	2.16	0.59
1:A:263:VAL:HG22	1:A:322:LEU:HB2	1.84	0.58
1:A:447:GLN:O	1:A:451:LEU:HD13	2.03	0.58
1:A:569:ARG:HH22	2:C:127:ILE:HG12	1.68	0.58
1:B:113:GLU:HG3	1:B:435:VAL:HG13	1.86	0.58
1:B:247:ILE:O	1:B:306:LYS:HA	2.03	0.58
1:B:136:TYR:CE2	1:B:156:ALA:HB1	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:ILE:HG23	1:A:10:ILE:O	2.04	0.58
1:B:83:PHE:O	1:B:87:ARG:CG	2.52	0.58
1:A:126:VAL:HG21	1:A:211:LEU:HD23	1.86	0.57
1:B:421:GLN:O	1:B:425:GLN:HG3	2.04	0.57
2:C:121:GLU:OE1	2:C:122:LEU:N	2.36	0.57
1:A:576:LYS:HE2	2:C:124:ILE:CG1	2.35	0.57
1:B:11:LEU:HD13	1:B:11:LEU:H	1.69	0.57
1:B:215:LEU:HA	1:B:219:ARG:HD2	1.85	0.57
1:B:353:LYS:HB3	1:B:398:MET:SD	2.45	0.57
1:A:235:LEU:HD22	1:A:250:ILE:HD13	1.87	0.57
1:B:410:CYS:O	1:B:432:ASN:HA	2.05	0.57
1:B:136:TYR:HE2	1:B:156:ALA:HB1	1.69	0.57
1:A:503:ASP:CG	1:A:507:LYS:HB2	2.25	0.57
1:B:37:GLY:O	1:B:39:THR:HG23	2.04	0.57
1:B:409:LEU:HB3	1:B:439:LEU:HD21	1.87	0.57
1:B:457:ASP:O	1:B:460:ASP:HB2	2.05	0.56
1:B:42:ILE:HD13	1:B:306:LYS:HG3	1.88	0.56
1:B:431:PHE:CZ	1:B:443:PHE:HA	2.40	0.56
1:B:469:GLN:HB2	1:B:487:LEU:HD11	1.86	0.56
1:B:358:GLN:NE2	1:B:387:VAL:HA	2.19	0.56
1:B:4:ARG:HD2	1:B:5:SER:O	2.06	0.56
1:A:21:LEU:O	1:A:25:ILE:HG13	2.06	0.56
1:A:301:ALA:O	1:A:302:ALA:HB3	2.05	0.56
1:B:12:GLY:H	1:B:18:LYS:HD3	1.70	0.56
1:A:264:LEU:HA	1:A:337:VAL:HG11	1.87	0.56
1:B:411:PHE:HE2	1:B:439:LEU:HD22	1.72	0.55
1:A:409:LEU:HD23	1:A:431:PHE:HD2	1.72	0.55
2:C:123:ASP:HB3	2:C:125:ASP:OD1	2.07	0.55
1:A:503:ASP:HB2	1:A:507:LYS:HB2	1.87	0.55
1:B:5:SER:OG	1:B:70:PRO:HG2	2.07	0.55
1:B:264:LEU:HD13	1:B:315:ALA:HB1	1.89	0.55
1:A:569:ARG:NH2	2:C:127:ILE:O	2.40	0.54
1:B:260:ASP:HB2	1:B:275:ILE:HD12	1.89	0.54
1:A:232:ALA:CB	1:A:321:LEU:HB3	2.37	0.54
1:A:52:ILE:HD11	1:A:72:LEU:HB2	1.90	0.54
1:A:83:PHE:O	1:A:86:LEU:CB	2.55	0.54
1:B:314:LYS:O	1:B:314:LYS:CD	2.55	0.54
1:A:22:LEU:HA	1:A:25:ILE:HD12	1.89	0.54
1:A:318:GLY:O	1:A:362:LEU:HD12	2.08	0.54
1:B:240:VAL:HG12	1:B:242:GLY:H	1.73	0.54
1:B:256:LEU:HD13	1:B:307:ILE:HD11	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:264:LEU:HB3	1:B:321:LEU:HD23	1.88	0.54
1:B:454:ARG:HA	1:B:457:ASP:HB3	1.88	0.54
1:A:345:LEU:O	1:A:348:VAL:HG22	2.08	0.54
2:C:117:GLY:O	2:C:118:GLU:HB3	2.08	0.54
1:B:387:VAL:CG2	1:B:416:ASP:H	2.21	0.54
1:B:10:ILE:HG23	1:B:10:ILE:O	2.08	0.53
1:A:411:PHE:CE1	1:A:434:ASP:O	2.62	0.53
1:B:458:PHE:HA	1:B:461:TYR:CB	2.37	0.53
1:A:455:ARG:HG3	1:A:456:LYS:N	2.24	0.53
1:B:84:SER:O	1:B:87:ARG:O	2.26	0.53
1:B:314:LYS:O	1:B:314:LYS:CG	2.56	0.53
1:B:95:ASN:O	1:B:96:ILE:HG13	2.09	0.53
1:A:411:PHE:CD1	1:A:412:ASP:HB2	2.44	0.53
1:B:353:LYS:HA	1:B:398:MET:CE	2.39	0.53
2:C:121:GLU:C	2:C:122:LEU:HD23	2.30	0.52
1:A:463:ILE:HD13	1:A:566:LEU:O	2.09	0.52
1:B:149:SER:O	1:B:153:GLN:HG2	2.09	0.52
1:A:42:ILE:CD1	1:A:306:LYS:HB2	2.37	0.52
1:B:251:LEU:HG	1:B:301:ALA:HA	1.92	0.52
1:A:284:LEU:HD22	1:A:287:LEU:HB2	1.92	0.52
1:A:299:VAL:HB	1:A:303:LEU:HD21	1.91	0.52
1:B:314:LYS:HD2	1:B:314:LYS:C	2.30	0.52
1:B:11:LEU:H	1:B:11:LEU:CD1	2.23	0.52
1:A:53:LYS:HG2	1:A:67:PHE:CE2	2.45	0.52
1:A:141:ILE:HG22	1:A:144:ASN:HB2	1.92	0.51
1:B:357:VAL:HG12	1:B:409:LEU:HB2	1.92	0.51
1:B:11:LEU:CD1	1:B:11:LEU:N	2.73	0.51
1:B:394:LYS:O	1:B:397:THR:HG22	2.10	0.51
1:A:353:LYS:O	1:A:353:LYS:HG2	2.10	0.51
1:A:84:SER:CB	1:A:93:LEU:HD12	2.40	0.51
1:A:572:ILE:HG12	1:A:592:LEU:CD2	2.41	0.51
1:B:392:VAL:O	1:B:396:SER:N	2.39	0.51
1:B:49:ILE:HD11	1:B:67:PHE:HD1	1.76	0.51
1:A:185:ASN:OD1	1:A:191:TYR:HB3	2.11	0.51
1:B:248:ASP:OD1	1:B:306:LYS:HG2	2.11	0.51
1:B:251:LEU:CD2	1:B:301:ALA:HA	2.41	0.50
1:B:357:VAL:HG23	1:B:382:ILE:HG22	1.91	0.50
1:B:448:GLU:O	1:B:451:LEU:HB3	2.11	0.50
1:A:330:GLU:O	1:A:334:MET:HG3	2.12	0.50
1:B:260:ASP:HB2	1:B:275:ILE:CD1	2.41	0.50
1:B:132:ILE:O	1:B:135:LEU:HG	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:LEU:O	1:A:26:ARG:HG3	2.11	0.50
1:A:364:SER:HB2	1:A:411:PHE:CE2	2.46	0.50
1:A:458:PHE:O	1:A:462:ALA:N	2.43	0.50
1:A:503:ASP:CB	1:A:507:LYS:HB2	2.42	0.50
1:A:11:LEU:N	1:A:11:LEU:CD1	2.74	0.50
1:B:13:HIS:HD2	1:B:15:ASP:H	1.60	0.50
1:B:42:ILE:HG13	1:B:281:PRO:CD	2.42	0.50
1:B:212:LEU:O	1:B:216:THR:HG22	2.12	0.50
1:B:270:PRO:C	1:B:271:ILE:HD13	2.32	0.50
1:B:303:LEU:HD12	1:B:304:GLY:O	2.12	0.50
1:A:522:ILE:HG12	1:A:539:VAL:HB	1.92	0.49
1:B:313:GLU:O	1:B:313:GLU:HG3	2.12	0.49
1:A:4:ARG:O	1:A:4:ARG:HD3	2.12	0.49
1:B:107:LEU:HD22	1:B:171:GLU:HB3	1.92	0.49
1:B:518:ILE:HD11	1:B:544:GLU:HB2	1.94	0.49
1:A:119:ARG:HH21	1:B:376:LYS:NZ	2.10	0.49
1:B:530:VAL:HG21	1:B:539:VAL:HG21	1.95	0.49
1:A:80:HIS:O	1:A:84:SER:N	2.34	0.49
1:B:12:GLY:N	1:B:18:LYS:HD3	2.27	0.49
1:A:404:GLU:HG2	1:A:405:TYR:N	2.27	0.49
1:A:584:VAL:HG12	1:A:585:ALA:O	2.12	0.49
1:B:237:VAL:HG21	1:B:315:ALA:O	2.13	0.49
1:B:477:ARG:HG2	1:B:478:GLY:N	2.27	0.49
1:A:257:ARG:HG2	1:A:298:GLU:HB2	1.95	0.49
1:A:371:PHE:CD2	1:A:436:ILE:HD13	2.48	0.49
1:A:469:GLN:HB2	1:A:487:LEU:HD11	1.95	0.49
1:B:7:ILE:HG12	1:B:73:LEU:HD23	1.94	0.49
1:A:390:ARG:O	1:A:393:MET:HB2	2.13	0.48
1:B:393:MET:HA	1:B:396:SER:OG	2.13	0.48
1:A:462:ALA:O	1:A:463:ILE:HG13	2.13	0.48
1:A:39:THR:HB	1:A:46:TYR:HB3	1.96	0.48
1:B:358:GLN:CG	1:B:413:VAL:HG21	2.44	0.48
1:B:521:GLU:HA	1:B:525:GLN:O	2.13	0.48
1:A:205:VAL:HB	1:A:206:PRO:HD3	1.95	0.48
1:B:301:ALA:O	1:B:302:ALA:HB3	2.12	0.48
1:B:358:GLN:HG3	1:B:413:VAL:HG21	1.95	0.48
1:A:190:LYS:HE3	1:B:347:SER:HA	1.95	0.48
1:A:409:LEU:HD23	1:A:431:PHE:CD2	2.49	0.48
1:A:471:LEU:HD21	1:A:485:ASP:HB2	1.96	0.48
1:B:132:ILE:HG12	1:B:197:THR:O	2.13	0.48
1:B:513:ILE:CD1	1:B:566:LEU:HG	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:ILE:HG21	1:A:22:LEU:CD1	2.44	0.48
1:A:185:ASN:HD22	1:A:186:LYS:N	2.11	0.48
1:A:141:ILE:O	1:A:144:ASN:HB2	2.14	0.48
1:A:501:LYS:O	1:A:510:GLN:HB3	2.12	0.48
1:B:355:VAL:HG11	1:B:409:LEU:HD12	1.95	0.48
1:B:416:ASP:HB2	1:B:419:ALA:HB3	1.96	0.48
1:B:503:ASP:HB2	1:B:506:THR:OG1	2.13	0.48
1:B:265:CYS:HB3	1:B:319:SER:OG	2.14	0.47
1:A:576:LYS:HZ2	2:C:127:ILE:C	2.17	0.47
1:A:581:ARG:HD2	1:A:582:ASP:N	2.30	0.47
1:A:293:TYR:CD1	1:A:293:TYR:N	2.78	0.47
1:A:570:ARG:O	1:A:574:THR:HG23	2.14	0.47
1:B:273:THR:CB	1:B:311:ASP:HB3	2.45	0.47
1:B:251:LEU:HD23	1:B:301:ALA:HA	1.95	0.47
1:B:53:LYS:HG2	1:B:67:PHE:CZ	2.49	0.47
1:A:411:PHE:CD1	1:A:411:PHE:C	2.87	0.47
1:B:23:ASP:HA	1:B:26:ARG:HH12	1.77	0.47
1:B:181:LEU:HD13	1:B:183:PHE:CZ	2.50	0.47
1:B:253:ASN:C	1:B:301:ALA:HB2	2.34	0.47
1:B:9:CYS:HB2	1:B:87:ARG:HH11	1.59	0.47
1:B:42:ILE:HD13	1:B:306:LYS:N	2.30	0.47
1:B:416:ASP:HB2	1:B:419:ALA:CB	2.44	0.47
1:A:454:ARG:O	1:A:458:PHE:HB3	2.15	0.47
1:B:106:GLY:HA2	1:B:168:ILE:HG12	1.96	0.47
1:A:115:ILE:HG22	1:A:119:ARG:HH11	1.80	0.46
1:A:463:ILE:HD11	1:A:496:PRO:HG2	1.97	0.46
1:A:380:MET:CE	1:A:394:LYS:HB3	2.45	0.46
1:B:513:ILE:HD11	1:B:566:LEU:HG	1.96	0.46
1:A:284:LEU:HB2	1:A:287:LEU:HB3	1.97	0.46
1:B:141:ILE:HG22	1:B:144:ASN:HB2	1.98	0.46
1:B:355:VAL:HG11	1:B:409:LEU:CD1	2.46	0.46
1:B:360:SER:HA	1:B:413:VAL:CG1	2.46	0.46
1:A:431:PHE:HE1	1:A:442:SER:HG	1.63	0.46
1:B:398:MET:SD	1:B:401:LYS:HD3	2.56	0.46
1:B:11:LEU:HD13	1:B:11:LEU:N	2.31	0.46
1:B:42:ILE:HG13	1:B:281:PRO:CG	2.45	0.45
1:B:477:ARG:HG2	1:B:478:GLY:H	1.80	0.45
1:B:273:THR:HG21	1:B:311:ASP:OD2	2.16	0.45
1:B:399:LEU:HG	1:B:400:GLU:N	2.30	0.45
1:B:257:ARG:HA	1:B:278:LEU:HD11	1.98	0.45
1:A:288:ARG:HA	1:A:291:SER:OG	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:433:ALA:CB	1:A:438:HIS:HB3	2.46	0.45
1:A:584:VAL:HG11	1:A:589:TRP:CE2	2.50	0.45
1:B:42:ILE:HG22	1:B:42:ILE:O	2.16	0.45
1:A:42:ILE:HG23	1:A:42:ILE:O	2.16	0.45
1:B:572:ILE:HG12	1:B:592:LEU:HD11	1.99	0.45
1:B:9:CYS:HB3	1:B:87:ARG:HH11	1.80	0.45
1:A:288:ARG:HD2	1:A:288:ARG:O	2.17	0.45
1:B:299:VAL:HB	1:B:303:LEU:CD2	2.37	0.45
1:B:495:THR:O	1:B:517:VAL:HG23	2.17	0.45
1:A:77:THR:HB	1:A:80:HIS:H	1.81	0.45
1:A:458:PHE:CE1	1:A:579:ALA:HB2	2.52	0.45
1:A:41:GLN:OE1	1:A:303:LEU:HA	2.16	0.44
1:A:303:LEU:HD12	1:A:303:LEU:C	2.38	0.44
1:B:180:GLU:OE1	1:B:180:GLU:HA	2.17	0.44
1:B:276:ARG:HB3	1:B:308:ALA:O	2.17	0.44
1:A:10:ILE:CG2	1:A:76:ASP:HA	2.46	0.44
1:A:263:VAL:HG23	1:A:337:VAL:HG21	1.99	0.44
1:B:305:VAL:HG23	1:B:306:LYS:N	2.32	0.44
1:A:83:PHE:HA	1:A:86:LEU:CB	2.47	0.44
1:A:507:LYS:HD2	1:A:507:LYS:HA	1.39	0.44
1:B:170:LEU:O	1:B:174:GLU:HG3	2.17	0.44
1:B:266:GLY:HA2	1:B:315:ALA:HA	1.99	0.44
1:A:368:LEU:O	1:A:372:LEU:HG	2.17	0.44
1:A:356:VAL:HG13	1:A:381:SER:OG	2.17	0.44
1:B:518:ILE:CD1	1:B:544:GLU:HB2	2.48	0.44
1:B:75:ILE:HG22	1:B:76:ASP:O	2.18	0.44
1:B:85:ASN:CA	1:B:363:GLY:HA3	2.41	0.44
1:A:12:GLY:H	1:A:18:LYS:HD3	1.83	0.43
1:B:358:GLN:O	1:B:410:CYS:HA	2.18	0.43
1:A:135:LEU:HB2	1:A:138:TRP:HB2	1.99	0.43
1:A:182:TYR:CD2	1:A:195:VAL:HG22	2.52	0.43
1:A:284:LEU:CD2	1:A:287:LEU:HB2	2.48	0.43
1:A:324:VAL:HG11	1:A:330:GLU:HG2	2.00	0.43
1:B:271:ILE:HD12	1:B:313:GLU:HG2	1.99	0.43
1:B:275:ILE:HG23	1:B:307:ILE:CG2	2.40	0.43
1:B:555:ARG:HA	1:B:555:ARG:HD3	1.89	0.43
1:A:27:GLN:OE1	1:A:27:GLN:HA	2.19	0.43
1:A:501:LYS:HD2	1:A:561:ASP:OD2	2.18	0.43
1:B:88:SER:N	1:B:91:SER:O	2.51	0.43
1:B:261:ARG:O	1:B:324:VAL:HG12	2.18	0.43
1:B:12:GLY:O	1:B:18:LYS:HD3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:387:VAL:HG21	1:B:416:ASP:H	1.83	0.43
1:B:85:ASN:HA	1:B:363:GLY:CA	2.41	0.43
1:B:86:LEU:HD13	1:B:86:LEU:N	2.34	0.43
1:B:264:LEU:HB3	1:B:321:LEU:CD2	2.49	0.43
1:A:284:LEU:HB3	1:A:287:LEU:H	1.83	0.43
1:B:86:LEU:N	1:B:86:LEU:CD1	2.82	0.43
1:B:357:VAL:HG23	1:B:382:ILE:HA	2.00	0.43
1:A:596:LYS:HG3	1:A:597:VAL:N	2.33	0.43
1:B:4:ARG:O	1:B:4:ARG:HG3	2.19	0.43
1:A:300:LYS:CG	1:A:301:ALA:N	2.67	0.43
1:A:303:LEU:C	1:A:303:LEU:CD1	2.87	0.43
1:B:250:ILE:HG22	1:B:250:ILE:O	2.18	0.43
1:B:251:LEU:CG	1:B:301:ALA:HA	2.49	0.43
1:B:487:LEU:C	1:B:488:GLU:HG2	2.40	0.43
1:A:473:ILE:HG12	1:A:476:LYS:HG2	2.01	0.43
1:A:576:LYS:HE2	2:C:124:ILE:HG13	2.00	0.42
1:A:88:SER:O	1:A:91:SER:O	2.38	0.42
1:A:137:ASP:HB3	1:A:157:VAL:HG23	2.00	0.42
1:B:42:ILE:CD1	1:B:306:LYS:HG3	2.49	0.42
1:A:455:ARG:HA	1:A:458:PHE:CD2	2.42	0.42
1:B:353:LYS:HD3	1:B:401:LYS:HG2	2.00	0.42
1:B:354:GLY:C	1:B:380:MET:HB2	2.39	0.42
1:B:388:TYR:HB2	1:B:391:ASP:OD2	2.19	0.42
1:B:136:TYR:HB3	1:B:160:GLU:OE2	2.19	0.42
1:B:353:LYS:HA	1:B:398:MET:HE3	2.01	0.42
1:A:38:ILE:HD12	1:A:38:ILE:C	2.40	0.42
1:A:106:GLY:HA2	1:A:168:ILE:HG12	2.00	0.42
1:B:254:GLY:O	1:B:301:ALA:HB2	2.19	0.42
1:B:11:LEU:HD21	1:B:114:SER:HB3	2.01	0.42
1:A:229:HIS:O	1:A:253:ASN:HB3	2.20	0.42
1:A:389:LYS:HA	1:A:392:VAL:HG12	2.02	0.42
1:B:103:ILE:HG22	1:B:130:ASN:O	2.20	0.42
1:A:119:ARG:HD2	1:A:177:LEU:HD21	2.01	0.41
1:B:357:VAL:HG11	1:B:372:LEU:CD1	2.47	0.41
1:B:525:GLN:HA	1:B:526:PRO:HD3	1.90	0.41
1:A:141:ILE:HG22	1:A:141:ILE:O	2.20	0.41
1:B:56:THR:O	1:B:59:MET:HB2	2.20	0.41
1:B:490:THR:CG2	1:B:529:GLU:HG3	2.50	0.41
1:B:102:ASP:OD1	1:B:104:MET:HB2	2.21	0.41
1:B:356:VAL:HB	1:B:395:ALA:CB	2.51	0.41
1:A:416:ASP:HB3	1:A:419:ALA:H	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:5:SER:HA	1:B:6:PRO:HD3	1.88	0.41
1:A:332:GLU:O	1:A:335:ASP:HB2	2.20	0.41
1:A:130:ASN:HD21	1:A:199:ALA:H	1.67	0.41
1:A:355:VAL:HG12	1:A:405:TYR:HA	2.02	0.41
1:B:422:TYR:HA	1:B:425:GLN:OE1	2.21	0.41
1:B:531:LYS:HE2	1:B:531:LYS:HB2	1.93	0.41
1:A:52:ILE:HG21	1:A:213:LEU:HD11	2.02	0.41
1:A:576:LYS:HE2	2:C:124:ILE:HG12	2.01	0.41
1:B:6:PRO:HG2	1:B:216:THR:HG21	2.03	0.41
1:B:87:ARG:HH22	1:B:97:ALA:HB2	1.86	0.41
1:B:304:GLY:C	1:B:305:VAL:HG12	2.41	0.41
1:B:264:LEU:C	1:B:264:LEU:HD12	2.41	0.41
1:A:596:LYS:HD3	2:C:127:ILE:HD12	2.03	0.40
1:B:88:SER:HB2	1:B:91:SER:O	2.20	0.40
1:A:109:GLN:O	1:A:112:ILE:HB	2.21	0.40
1:A:139:LYS:HE3	1:A:139:LYS:HB3	1.98	0.40
1:B:9:CYS:CB	1:B:83:PHE:CZ	2.96	0.40
1:B:95:ASN:C	1:B:96:ILE:HG13	2.42	0.40
1:B:230:VAL:HG23	1:B:251:LEU:HD11	2.02	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	584/606 (96%)	561 (96%)	22 (4%)	1 (0%)	47 77
1	B	584/606 (96%)	566 (97%)	18 (3%)	0	100 100
2	C	9/131 (7%)	9 (100%)	0	0	100 100
All	All	1177/1343 (88%)	1136 (96%)	40 (3%)	1 (0%)	51 81

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	463	ILE

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	506/529 (96%)	482 (95%)	24 (5%)	26 57
1	B	518/529 (98%)	490 (95%)	28 (5%)	22 53
2	C	10/113 (9%)	8 (80%)	2 (20%)	1 5
All	All	1034/1171 (88%)	980 (95%)	54 (5%)	23 54

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ARG
1	A	11	LEU
1	A	22	LEU
1	A	27	GLN
1	A	39	THR
1	A	77	THR
1	A	94	CYS
1	A	136	TYR
1	A	170	LEU
1	A	185	ASN
1	A	207	ASP
1	A	287	LEU
1	A	288	ARG
1	A	303	LEU
1	A	353	LYS
1	A	376	LYS
1	A	411	PHE
1	A	415	VAL
1	A	455	ARG
1	A	458	PHE
1	A	507	LYS
1	A	564	TYR

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Mol	Chain	Res	Type
1	A	576	LYS
1	A	592	LEU
1	B	4	ARG
1	B	10	ILE
1	B	11	LEU
1	B	19	THR
1	B	24	LYS
1	B	42	ILE
1	B	83	PHE
1	B	84	SER
1	B	86	LEU
1	B	87	ARG
1	B	91	SER
1	B	136	TYR
1	B	246	THR
1	B	289	LEU
1	B	303	LEU
1	B	305	VAL
1	B	310	ASN
1	B	353	LYS
1	B	399	LEU
1	B	411	PHE
1	B	415	VAL
1	B	453	GLU
1	B	458	PHE
1	B	482	ILE
1	B	519	SER
1	B	529	GLU
1	B	535	THR
1	B	566	LEU
2	C	121	GLU
2	C	122	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	130	ASN
1	A	268	ASN
1	B	13	HIS
1	B	130	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	590/606 (97%)	0.42	36 (6%) 21 20	71, 144, 219, 298	0
1	B	590/606 (97%)	0.66	76 (12%) 3 3	75, 161, 262, 324	0
2	C	11/131 (8%)	0.59	1 (9%) 9 9	183, 204, 232, 238	0
All	All	1191/1343 (88%)	0.54	113 (9%) 8 8	71, 151, 246, 324	0

All (113) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	407	VAL	8.3
1	B	358	GLN	6.8
1	B	430	ILE	6.3
1	B	406	ALA	5.6
1	B	355	VAL	5.1
1	B	409	LEU	5.0
1	B	404	GLU	4.9
1	B	429	LYS	4.7
1	B	431	PHE	4.6
1	B	438	HIS	4.5
1	A	427	GLY	4.4
1	B	387	VAL	4.3
1	B	410	CYS	4.3
1	B	249	VAL	4.2
1	B	403	PRO	4.1
1	A	396	SER	4.1
1	B	446	TYR	4.1
1	A	66	THR	4.1
1	B	415	VAL	4.0
1	A	387	VAL	4.0
1	A	64	LYS	4.0
1	A	353	LYS	3.9
1	A	76	ASP	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	421	GLN	3.8
1	B	333	LEU	3.8
1	B	382	ILE	3.8
1	B	427	GLY	3.8
1	B	391	ASP	3.7
1	B	428	ILE	3.7
1	B	255	TYR	3.6
1	B	230	VAL	3.5
1	B	357	VAL	3.4
1	A	35	ALA	3.4
1	B	307	ILE	3.4
1	A	392	VAL	3.4
1	B	107	LEU	3.4
1	B	423	ALA	3.4
1	B	243	PHE	3.4
1	A	354	GLY	3.4
1	B	414	LYS	3.3
1	B	350	THR	3.3
1	B	353	LYS	3.3
1	B	188	MET	3.2
1	A	583	GLN	3.1
1	B	408	MET	3.1
1	A	401	LYS	3.1
1	B	365	LEU	3.1
1	B	386	PRO	3.0
1	A	26	ARG	3.0
1	A	409	LEU	3.0
1	B	260	ASP	3.0
1	A	67	PHE	2.9
1	A	228	SER	2.9
1	B	399	LEU	2.9
1	B	439	LEU	2.9
1	A	378	PRO	2.9
1	B	282	GLN	2.8
1	B	323	VAL	2.8
1	A	341	LEU	2.8
1	B	381	SER	2.7
1	B	4	ARG	2.7
1	B	389	LYS	2.7
1	B	432	ASN	2.6
1	A	52	ILE	2.6
1	B	232	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	400	GLU	2.6
1	B	380	MET	2.6
1	A	588	ASP	2.6
1	B	246	THR	2.6
1	A	411	PHE	2.5
1	A	407	VAL	2.5
1	A	263	VAL	2.4
1	A	333	LEU	2.4
1	B	505	THR	2.4
1	A	285	ARG	2.3
1	B	301	ALA	2.3
1	A	395	ALA	2.3
1	A	355	VAL	2.3
1	B	575	LEU	2.3
1	B	237	VAL	2.3
1	B	453	GLU	2.3
1	A	379	VAL	2.3
1	B	321	LEU	2.3
1	B	464	PHE	2.3
1	A	92	SER	2.3
1	A	199	ALA	2.3
1	A	428	ILE	2.2
1	B	454	ARG	2.2
1	B	419	ALA	2.2
1	A	453	GLU	2.2
1	A	227	LEU	2.2
1	B	405	TYR	2.2
1	B	369	LEU	2.2
1	B	390	ARG	2.2
2	C	127	ILE	2.2
1	B	433	ALA	2.1
1	B	411	PHE	2.1
1	B	567	VAL	2.1
1	B	115	ILE	2.1
1	B	254	GLY	2.1
1	B	315	ALA	2.1
1	B	401	LYS	2.1
1	B	327	GLU	2.1
1	B	504	PRO	2.1
1	B	259	GLY	2.1
1	B	388	TYR	2.1
1	B	11	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	136	TYR	2.0
1	B	383	GLY	2.0
1	B	229	HIS	2.0
1	A	402	ALA	2.0
1	A	439	LEU	2.0
1	B	83	PHE	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 monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.