Full wwPDB X-ray Structure Validation Report (i)

Oct 16, 2023 – 07:04 PM EDT

PDB ID	:	2YXR
Title	:	The plug domain of the SecY protein stablizes the closed state of the translocation channel and maintains a membrane seal
Authors	:	Li, W.; Schulman, S.
Deposited on	:	2007-04-27
Resolution	:	3.60 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:

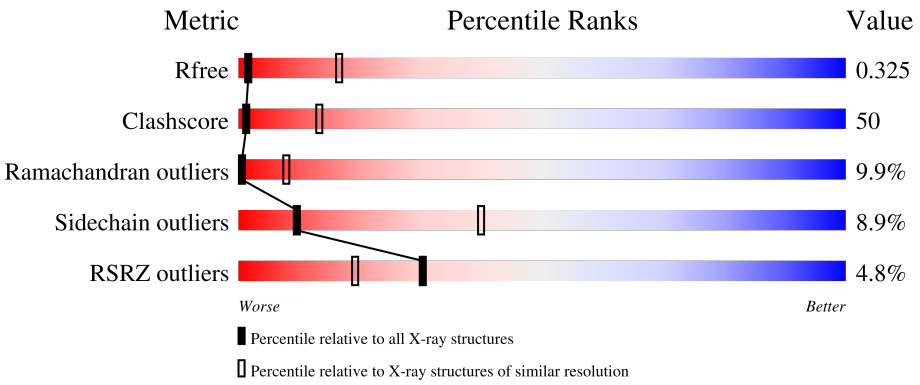
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 3.60 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	1257 (3.70-3.50)
Clashscore	141614	1353 (3.70-3.50)
Ramachandran outliers	138981	1307 (3.70-3.50)
Sidechain outliers	138945	1307 (3.70-3.50)
RSRZ outliers	127900	1161 (3.70-3.50)

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	426	6%; 30%; 57%; 12%; • •
2	B	74	32%; 46%; 8%; •; 12%
3	C	53	2%; 25%; 30%; 6%; 40%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3999 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 Preprotein translocase subunit secY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	422	3218	2150	506	543	19	0	0	0

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	GLU	deletion	UNP Q60175
A	?	-	PHE	deletion	UNP Q60175
A	?	-	TRP	deletion	UNP Q60175
A	?	-	GLN	deletion	UNP Q60175
A	?	-	THR	deletion	UNP Q60175
A	?	-	ILE	deletion	UNP Q60175
A	?	-	THR	deletion	UNP Q60175
A	?	-	ALA	deletion	UNP Q60175
A	?	-	SER	deletion	UNP Q60175
A	?	-	ARG	deletion	UNP Q60175
A	67	GLY	ILE	SEE REMARK 999	UNP Q60175

- Molecule 2 is a protein called Preprotein translocase subunit secE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	65	524	348	85	90	1	0	0	0

- Molecule 3 is a protein called Preprotein translocase secG subunit.

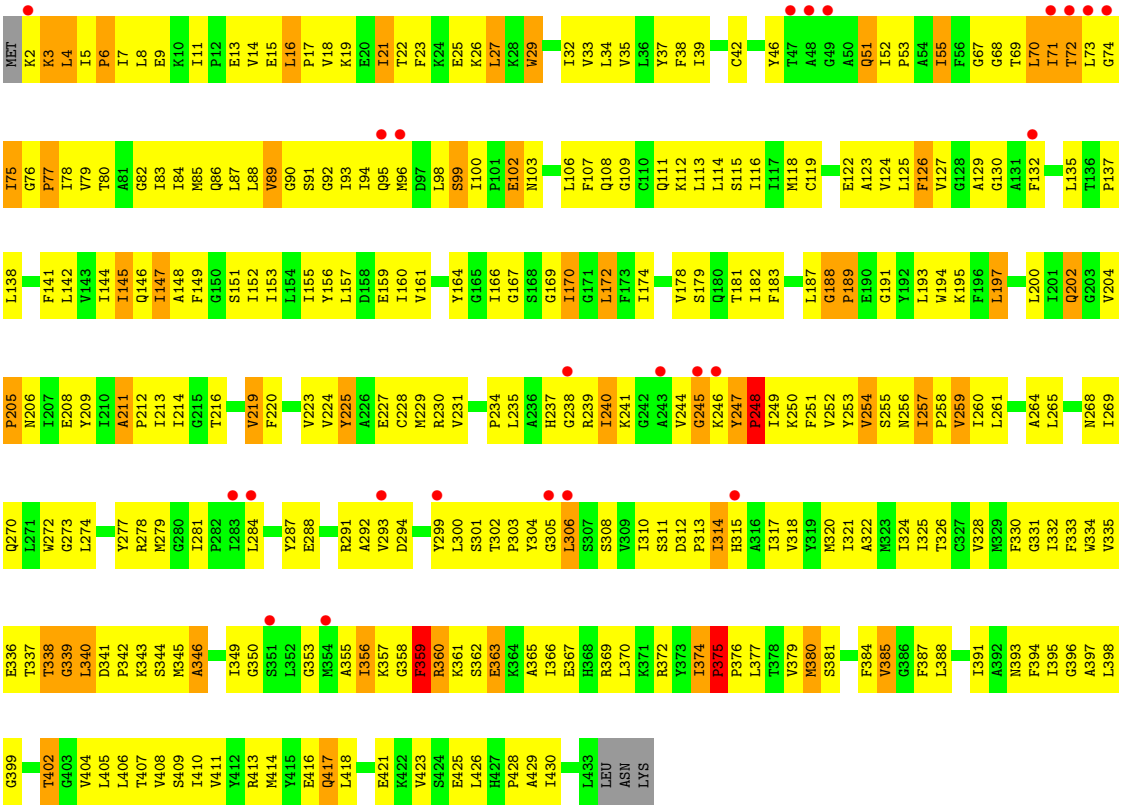
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	C	32	257	172	42	43	0	0	0

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.

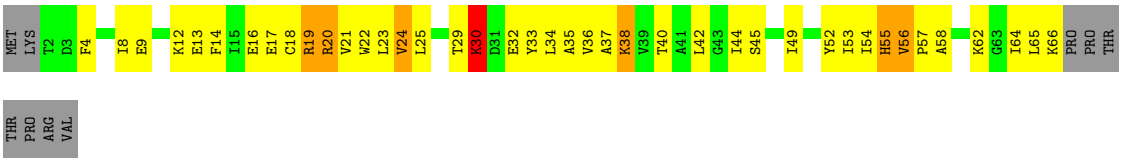
• Molecule 1: Preprotein translocase subunit secY

Chain A: 6% 30% 57% 12%



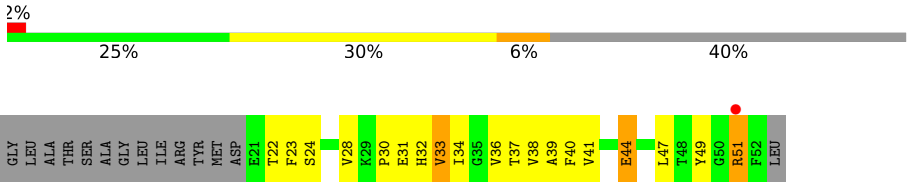
• Molecule 2: Preprotein translocase subunit secE

Chain B: 32% 46% 8% 12%



• Molecule 3: Preprotein translocase secG subunit

Chain C:


2%
25%
30%
6%
40%
MET SER LYS ARG GLU GLU THR GLY LEU ALA THR SER ALA GLY LEU ILE ARG TYR MET ASP E21 T22 F23 S24 V28 K29 P30 E31 H32 V33 I34 G35 V36 T37 V38 A39 F40 V41 E44 L47 T48 Y49 G50 R51 F52 LEU

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	92.51Å 150.36Å 79.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.07 – 3.60 44.07 – 3.60	Depositor EDS
% Data completeness (in resolution range)	65.7 (44.07-3.60) 65.7 (44.07-3.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	5.52 (at 3.57Å)	Xtriage
Refinement program	REFMAC 5.3.0026	Depositor
R, R_{free}	0.298 , 0.335 0.300 , 0.325	Depositor DCC
R_{free} test set	465 reflections (5.27%)	wwPDB-VP
Wilson B-factor (Å^2)	116.5	Xtriage
Anisotropy	0.126	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.26 , 100.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.85	EDS
Total number of atoms	3999	wwPDB-VP
Average B, all atoms (Å^2)	119.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% 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	0.59	0/3289	0.65	0/4464
2	B	0.60	0/533	0.58	0/719
3	C	0.59	0/262	0.59	0/354
All	All	0.59	0/4084	0.64	0/5537

There are no bond length outliers.

There are no bond angle outliers.

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	3218	0	3434	369	0
2	B	524	0	567	53	0
3	C	257	0	272	28	0
All	All	3999	0	4273	411	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 50.

All (411) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:ILE:HD12	1:A:114:LEU:HD21	1.38	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:TRP:O	1:A:32:ILE:HG22	1.62	0.99
1:A:11:ILE:HD11	1:A:113:LEU:HD23	1.43	0.97
1:A:16:LEU:HD23	1:A:16:LEU:H	1.30	0.96
1:A:240:ILE:HG22	1:A:241:LYS:H	1.31	0.93
1:A:340:LEU:H	1:A:340:LEU:HD12	1.38	0.88
1:A:33:VAL:HG21	1:A:161:VAL:HG22	1.56	0.88
1:A:227:GLU:HA	1:A:252:VAL:HG21	1.55	0.87
1:A:314:ILE:HD12	1:A:315:HIS:H	1.39	0.87
1:A:426:LEU:O	1:A:428:PRO:HD3	1.75	0.86
1:A:377:LEU:HA	2:B:21:VAL:HG11	1.57	0.86
1:A:13:GLU:HG3	1:A:14:VAL:H	1.42	0.85
1:A:287:TYR:HA	1:A:292:ALA:HA	1.58	0.83
1:A:46:TYR:HB3	1:A:146:GLN:HE22	1.41	0.83
1:A:200:LEU:HD23	1:A:205:PRO:HG3	1.61	0.83
1:A:216:THR:HA	1:A:397:ALA:HB1	1.60	0.83
1:A:219:VAL:O	1:A:223:VAL:HG23	1.78	0.83
1:A:333:PHE:O	1:A:337:THR:HG22	1.78	0.82
1:A:356:ILE:HG22	1:A:358:GLY:H	1.45	0.82
1:A:393:ASN:HD21	1:A:402:THR:H	1.27	0.82
1:A:73:LEU:HD13	1:A:126:PHE:CE2	2.17	0.80
1:A:380:MET:HG2	2:B:18:CYS:HB3	1.65	0.79
1:A:73:LEU:C	1:A:75:ILE:H	1.87	0.78
1:A:68:GLY:HA2	1:A:73:LEU:H	1.49	0.78
1:A:124:VAL:HG22	1:A:144:ILE:HD13	1.64	0.78
1:A:183:PHE:CE1	2:B:49:ILE:HG13	2.18	0.77
1:A:360:ARG:HB3	1:A:365:ALA:HB1	1.65	0.77
1:A:52:ILE:HG12	1:A:132:PHE:HA	1.67	0.76
1:A:234:PRO:HB2	1:A:357:LYS:CB	2.15	0.76
1:A:356:ILE:HG22	1:A:357:LYS:N	2.00	0.76
3:C:36:VAL:O	3:C:39:ALA:HB3	1.86	0.76
1:A:231:VAL:HB	1:A:249:ILE:CG1	2.16	0.76
1:A:52:ILE:HD12	1:A:52:ILE:O	1.86	0.74
1:A:358:GLY:O	1:A:359:PHE:HB2	1.86	0.74
1:A:380:MET:SD	2:B:22:TRP:HB2	2.28	0.74
1:A:374:ILE:HB	1:A:375:PRO:HD3	1.70	0.74
1:A:223:VAL:HG11	1:A:405:LEU:HA	1.69	0.73
1:A:76:GLY:O	1:A:78:ILE:N	2.21	0.73
1:A:183:PHE:HE1	2:B:49:ILE:HG13	1.52	0.73
1:A:273:GLY:CA	1:A:284:LEU:HD12	2.20	0.72
1:A:149:PHE:O	1:A:152:ILE:HB	1.89	0.72
1:A:3:LYS:HE3	1:A:3:LYS:HA	1.70	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:LEU:HD23	1:A:160:ILE:HD12	1.71	0.72
1:A:273:GLY:HA2	1:A:284:LEU:HD12	1.71	0.72
1:A:234:PRO:HB2	1:A:357:LYS:HB2	1.71	0.72
1:A:68:GLY:HA2	1:A:72:THR:HB	1.71	0.71
1:A:33:VAL:HG13	1:A:157:LEU:HD22	1.72	0.71
1:A:98:LEU:HA	1:A:103:ASN:HB2	1.71	0.71
1:A:251:PHE:CE1	1:A:381:SER:HA	2.26	0.71
1:A:328:VAL:O	1:A:332:ILE:HG13	1.92	0.69
1:A:82:GLY:HA2	1:A:111:GLN:NE2	2.07	0.69
1:A:393:ASN:HD21	1:A:402:THR:N	1.91	0.69
1:A:253:TYR:HE1	1:A:413:ARG:HH11	1.43	0.67
1:A:5:ILE:N	1:A:6:PRO:HD2	2.10	0.67
1:A:174:ILE:O	1:A:178:VAL:HG23	1.95	0.67
1:A:408:VAL:HG13	2:B:40:THR:HG21	1.77	0.67
1:A:103:ASN:HA	1:A:106:LEU:HD12	1.76	0.67
1:A:149:PHE:HB3	3:C:40:PHE:CZ	2.31	0.66
1:A:299:TYR:O	1:A:300:LEU:HD23	1.96	0.66
1:A:231:VAL:HB	1:A:249:ILE:HD11	1.77	0.66
1:A:255:SER:C	1:A:258:PRO:HD2	2.16	0.66
3:C:32:HIS:O	3:C:36:VAL:HG23	1.96	0.65
1:A:281:ILE:N	1:A:281:ILE:HD12	2.11	0.65
1:A:164:TYR:CZ	3:C:30:PRO:HG2	2.31	0.65
1:A:388:LEU:O	1:A:391:ILE:HG22	1.96	0.64
1:A:82:GLY:HA2	1:A:111:GLN:HE21	1.62	0.64
1:A:102:GLU:O	1:A:106:LEU:HG	1.98	0.64
1:A:356:ILE:HG22	1:A:357:LYS:H	1.63	0.64
1:A:127:VAL:HG11	1:A:144:ILE:HD11	1.79	0.64
1:A:52:ILE:CG1	1:A:132:PHE:HA	2.28	0.63
1:A:164:TYR:CE1	3:C:30:PRO:HG2	2.34	0.63
1:A:82:GLY:O	1:A:86:GLN:HG2	1.97	0.63
1:A:257:ILE:HG21	1:A:334:TRP:CH2	2.33	0.63
1:A:189:PRO:HD3	2:B:56:VAL:HG22	1.81	0.63
1:A:76:GLY:C	1:A:78:ILE:N	2.49	0.63
1:A:157:LEU:O	1:A:161:VAL:HG23	1.99	0.63
1:A:240:ILE:HG22	1:A:241:LYS:N	2.09	0.62
1:A:355:ALA:HB2	1:A:361:LYS:HA	1.80	0.62
1:A:138:LEU:O	1:A:142:LEU:HG	1.99	0.62
1:A:228:CYS:SG	2:B:36:VAL:HG21	2.40	0.62
1:A:250:LYS:HD2	1:A:253:TYR:HD2	1.64	0.62
1:A:16:LEU:HD23	1:A:16:LEU:N	2.10	0.62
1:A:231:VAL:O	1:A:249:ILE:HG12	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:79:VAL:O	1:A:83:ILE:HG13	1.99	0.62
1:A:231:VAL:HB	1:A:249:ILE:HG13	1.80	0.62
1:A:231:VAL:HB	1:A:249:ILE:CD1	2.30	0.62
1:A:73:LEU:C	1:A:75:ILE:N	2.53	0.61
1:A:195:LYS:HB3	1:A:209:TYR:CD2	2.34	0.61
1:A:308:SER:HB3	1:A:394:PHE:O	2.00	0.61
1:A:14:VAL:HG21	1:A:159:GLU:HB3	1.81	0.61
1:A:98:LEU:HA	1:A:103:ASN:CB	2.30	0.61
1:A:156:TYR:O	1:A:160:ILE:HG13	2.00	0.61
1:A:235:LEU:HB2	1:A:245:GLY:HA3	1.83	0.61
1:A:152:ILE:HD13	1:A:155:ILE:HD12	1.82	0.61
1:A:188:GLY:HA2	2:B:52:TYR:HE1	1.65	0.61
1:A:379:VAL:CG1	2:B:18:CYS:SG	2.89	0.60
1:A:76:GLY:C	1:A:78:ILE:H	2.04	0.60
1:A:141:PHE:CZ	1:A:145:ILE:HD11	2.36	0.60
1:A:270:GLN:NE2	1:A:301:SER:HB3	2.17	0.60
1:A:254:VAL:HG13	1:A:381:SER:OG	2.02	0.60
1:A:123:ALA:HB2	1:A:148:ALA:HB2	1.84	0.60
1:A:153:ILE:HD13	3:C:37:THR:HG23	1.84	0.59
1:A:340:LEU:HD12	1:A:340:LEU:N	2.14	0.59
1:A:250:LYS:NZ	1:A:416:GLU:OE1	2.32	0.59
1:A:287:TYR:CE2	1:A:292:ALA:HB2	2.37	0.59
1:A:374:ILE:CB	1:A:375:PRO:HD3	2.33	0.59
1:A:23:PHE:HD1	1:A:421:GLU:HB3	1.67	0.59
1:A:247:TYR:CD1	1:A:247:TYR:C	2.76	0.59
1:A:160:ILE:HD13	3:C:34:ILE:HD11	1.85	0.58
1:A:379:VAL:HG11	2:B:18:CYS:SG	2.43	0.58
1:A:281:ILE:HD12	1:A:281:ILE:H	1.67	0.58
1:A:216:THR:HA	1:A:397:ALA:CB	2.32	0.58
1:A:257:ILE:HG22	1:A:258:PRO:HD3	1.85	0.58
1:A:356:ILE:CG2	1:A:357:LYS:N	2.67	0.58
2:B:58:ALA:O	2:B:62:LYS:HG3	2.03	0.58
1:A:404:VAL:O	1:A:408:VAL:HG23	2.02	0.58
1:A:273:GLY:HA3	1:A:287:TYR:OH	2.03	0.58
1:A:342:PRO:O	1:A:343:LYS:C	2.41	0.58
1:A:13:GLU:HG3	1:A:14:VAL:N	2.15	0.58
2:B:49:ILE:O	2:B:53:ILE:HG13	2.04	0.58
1:A:86:GLN:O	1:A:333:PHE:HD2	1.87	0.58
1:A:116:ILE:HA	1:A:119:CYS:HB2	1.85	0.58
1:A:279:MET:O	1:A:279:MET:HG2	2.03	0.58
2:B:16:GLU:HA	2:B:16:GLU:OE2	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:LEU:C	1:A:90:GLY:H	2.08	0.57
1:A:379:VAL:HG11	2:B:14:PHE:HE1	1.69	0.57
1:A:250:LYS:O	1:A:254:VAL:HG12	2.04	0.57
1:A:328:VAL:HG13	1:A:379:VAL:HG22	1.87	0.57
1:A:68:GLY:CA	1:A:72:THR:HB	2.33	0.57
1:A:304:TYR:HD1	1:A:305:GLY:N	2.02	0.57
1:A:425:GLU:O	1:A:426:LEU:HD12	2.05	0.56
2:B:17:GLU:O	2:B:20:ARG:HB2	2.05	0.56
1:A:23:PHE:CD1	1:A:421:GLU:HB3	2.40	0.56
1:A:125:LEU:HD21	1:A:272:TRP:CD1	2.41	0.56
1:A:268:ASN:O	1:A:272:TRP:HB2	2.04	0.56
1:A:375:PRO:HB2	1:A:376:PRO:CD	2.36	0.56
1:A:70:LEU:O	1:A:71:ILE:HB	2.04	0.56
1:A:324:ILE:HD11	1:A:387:PHE:HB2	1.88	0.56
2:B:52:TYR:CE1	2:B:56:VAL:HG21	2.41	0.56
1:A:73:LEU:HD13	1:A:126:PHE:CD2	2.41	0.56
1:A:80:THR:HG23	1:A:268:ASN:ND2	2.20	0.56
1:A:325:ILE:HG13	1:A:326:THR:N	2.20	0.56
1:A:355:ALA:HB3	1:A:361:LYS:HG2	1.88	0.56
1:A:15:GLU:OE1	1:A:17:PRO:HG3	2.04	0.56
1:A:321:ILE:O	1:A:325:ILE:HG23	2.06	0.55
1:A:55:ILE:O	1:A:69:THR:HB	2.06	0.55
1:A:306:LEU:N	1:A:306:LEU:HD12	2.22	0.55
2:B:4:PHE:O	2:B:8:ILE:HG13	2.07	0.55
1:A:89:VAL:HG21	1:A:107:PHE:CD1	2.41	0.55
1:A:345:MET:O	1:A:349:ILE:HG12	2.07	0.55
1:A:4:LEU:C	1:A:6:PRO:HD2	2.27	0.55
1:A:116:ILE:N	1:A:116:ILE:HD12	2.22	0.55
1:A:85:MET:HB3	1:A:107:PHE:CE1	2.42	0.55
1:A:91:SER:OG	1:A:93:ILE:HG13	2.06	0.55
1:A:317:ILE:O	1:A:321:ILE:HG13	2.08	0.54
3:C:22:THR:HG22	3:C:24:SER:H	1.72	0.54
1:A:257:ILE:HG21	1:A:334:TRP:CZ2	2.42	0.54
2:B:66:LYS:HB2	2:B:66:LYS:NZ	2.22	0.54
1:A:247:TYR:C	1:A:247:TYR:HD1	2.09	0.54
1:A:29:TRP:NE1	1:A:164:TYR:HD2	2.05	0.54
1:A:98:LEU:O	1:A:99:SER:C	2.45	0.54
1:A:211:ALA:HB3	1:A:212:PRO:CD	2.38	0.54
1:A:274:LEU:O	1:A:277:TYR:HB3	2.08	0.54
1:A:340:LEU:H	1:A:340:LEU:CD1	2.14	0.54
1:A:356:ILE:CG2	1:A:357:LYS:H	2.21	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:PRO:HA	1:A:345:MET:HB2	1.90	0.53
1:A:148:ALA:O	1:A:152:ILE:HG12	2.08	0.53
1:A:182:ILE:HA	1:A:398:LEU:HD23	1.89	0.53
1:A:255:SER:HA	1:A:258:PRO:HG2	1.90	0.53
1:A:260:ILE:O	1:A:264:ALA:HB2	2.07	0.53
1:A:355:ALA:CB	1:A:361:LYS:HG2	2.39	0.53
2:B:16:GLU:O	2:B:19:ARG:HB2	2.08	0.53
1:A:188:GLY:HA2	2:B:52:TYR:CE1	2.43	0.53
1:A:85:MET:HG3	1:A:111:GLN:HG3	1.90	0.53
1:A:72:THR:O	1:A:75:ILE:HB	2.09	0.52
1:A:256:ASN:HD21	1:A:409:SER:CB	2.21	0.52
1:A:259:VAL:CG1	1:A:406:LEU:HD21	2.38	0.52
1:A:15:GLU:O	1:A:17:PRO:HD3	2.09	0.52
1:A:166:ILE:HD13	1:A:418:LEU:CD2	2.39	0.52
1:A:379:VAL:HG12	1:A:380:MET:N	2.25	0.52
1:A:379:VAL:HG12	2:B:18:CYS:SG	2.50	0.52
1:A:146:GLN:HG2	3:C:44:GLU:OE1	2.10	0.52
2:B:32:GLU:O	2:B:35:ALA:HB3	2.09	0.52
1:A:312:ASP:N	1:A:313:PRO:HD3	2.25	0.52
3:C:34:ILE:O	3:C:38:VAL:HG23	2.10	0.52
1:A:257:ILE:HG22	1:A:258:PRO:CD	2.40	0.52
1:A:80:THR:HG23	1:A:268:ASN:HD21	1.74	0.52
1:A:94:ILE:O	1:A:94:ILE:HG22	2.10	0.52
1:A:250:LYS:CD	1:A:253:TYR:HD2	2.22	0.52
2:B:62:LYS:O	2:B:66:LYS:HB2	2.09	0.52
1:A:257:ILE:N	1:A:258:PRO:CD	2.72	0.51
1:A:160:ILE:HD11	3:C:33:VAL:CG2	2.41	0.51
1:A:260:ILE:HG12	1:A:406:LEU:HD13	1.91	0.51
1:A:135:LEU:H	1:A:135:LEU:HD12	1.75	0.51
1:A:320:MET:O	1:A:324:ILE:HG12	2.11	0.51
1:A:343:LYS:O	1:A:346:ALA:HB3	2.10	0.51
1:A:160:ILE:HD13	3:C:34:ILE:CD1	2.41	0.51
1:A:425:GLU:O	1:A:425:GLU:HG3	2.10	0.51
1:A:256:ASN:HD21	1:A:409:SER:HB2	1.76	0.51
1:A:84:ILE:CD1	1:A:114:LEU:HD21	2.27	0.51
1:A:93:ILE:HG22	1:A:94:ILE:HG13	1.92	0.51
1:A:362:SER:HB2	1:A:365:ALA:CB	2.41	0.51
1:A:332:ILE:O	1:A:335:VAL:HB	2.12	0.50
1:A:80:THR:O	1:A:83:ILE:HB	2.11	0.50
1:A:71:ILE:HG22	1:A:71:ILE:O	2.10	0.50
1:A:260:ILE:HG12	1:A:406:LEU:CD1	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:362:SER:HB2	1:A:365:ALA:HB2	1.94	0.50
1:A:86:GLN:O	1:A:90:GLY:HA3	2.11	0.50
1:A:179:SER:OG	2:B:44:ILE:HG23	2.12	0.50
1:A:250:LYS:HB3	1:A:253:TYR:HB3	1.92	0.50
1:A:70:LEU:O	1:A:71:ILE:CB	2.60	0.50
1:A:76:GLY:O	1:A:77:PRO:C	2.49	0.50
1:A:102:GLU:C	1:A:106:LEU:HG	2.32	0.50
1:A:274:LEU:HD23	1:A:287:TYR:CD2	2.47	0.50
2:B:49:ILE:O	2:B:52:TYR:HB3	2.12	0.50
1:A:89:VAL:HG21	1:A:107:PHE:HD1	1.77	0.49
1:A:340:LEU:O	1:A:342:PRO:HD3	2.12	0.49
2:B:29:THR:O	2:B:32:GLU:HB3	2.12	0.49
1:A:374:ILE:CB	1:A:375:PRO:CD	2.89	0.49
1:A:252:VAL:HG23	1:A:253:TYR:H	1.77	0.49
2:B:55:HIS:O	2:B:56:VAL:C	2.50	0.49
1:A:35:VAL:HG13	2:B:54:ILE:HD11	1.95	0.49
1:A:88:LEU:C	1:A:90:GLY:N	2.65	0.49
1:A:152:ILE:HD13	1:A:155:ILE:CD1	2.41	0.49
1:A:293:VAL:O	1:A:294:ASP:HB2	2.13	0.49
1:A:374:ILE:HB	1:A:375:PRO:CD	2.40	0.49
1:A:8:LEU:HD21	3:C:23:PHE:HB2	1.94	0.49
1:A:240:ILE:C	1:A:241:LYS:HG3	2.32	0.49
1:A:206:ASN:OD1	1:A:208:GLU:HB2	2.13	0.49
1:A:362:SER:O	1:A:365:ALA:HB3	2.13	0.49
1:A:21:ILE:HD12	1:A:21:ILE:H	1.78	0.49
1:A:100:ILE:HG22	1:A:102:GLU:H	1.76	0.49
1:A:84:ILE:O	1:A:87:LEU:HB2	2.13	0.49
1:A:142:LEU:O	1:A:145:ILE:HB	2.12	0.49
1:A:187:LEU:CD2	1:A:193:LEU:HD23	2.42	0.48
1:A:204:VAL:O	1:A:204:VAL:HG23	2.12	0.48
1:A:252:VAL:HG23	1:A:253:TYR:N	2.28	0.48
1:A:287:TYR:HB3	1:A:291:ARG:O	2.13	0.48
1:A:293:VAL:O	1:A:293:VAL:HG12	2.13	0.48
1:A:225:TYR:C	1:A:225:TYR:CD2	2.85	0.48
3:C:30:PRO:O	3:C:33:VAL:HG23	2.14	0.48
1:A:216:THR:CA	1:A:397:ALA:HB1	2.39	0.48
2:B:8:ILE:O	2:B:12:LYS:HG3	2.12	0.48
1:A:21:ILE:HD13	1:A:26:LYS:HE3	1.94	0.48
1:A:337:THR:O	1:A:338:THR:C	2.51	0.48
1:A:410:ILE:HA	1:A:413:ARG:HG2	1.95	0.48
1:A:304:TYR:CD1	1:A:305:GLY:N	2.82	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:314:ILE:HD12	1:A:315:HIS:N	2.18	0.48
1:A:374:ILE:HG22	1:A:375:PRO:N	2.29	0.48
2:B:36:VAL:O	2:B:37:ALA:C	2.52	0.48
1:A:34:LEU:O	1:A:37:TYR:HB3	2.14	0.48
1:A:229:MET:CE	2:B:25:LEU:HD13	2.43	0.48
1:A:314:ILE:CD1	1:A:315:HIS:H	2.19	0.48
2:B:53:ILE:O	2:B:57:PRO:HG3	2.14	0.47
1:A:73:LEU:O	1:A:75:ILE:N	2.47	0.47
1:A:274:LEU:C	1:A:274:LEU:HD13	2.34	0.47
1:A:274:LEU:O	1:A:274:LEU:HD13	2.14	0.47
1:A:151:SER:O	1:A:155:ILE:HG13	2.13	0.47
1:A:223:VAL:CG1	1:A:405:LEU:HA	2.40	0.47
1:A:359:PHE:O	1:A:360:ARG:HB2	2.14	0.47
1:A:23:PHE:CE2	1:A:27:LEU:HD12	2.50	0.47
1:A:129:ALA:O	1:A:278:ARG:HD3	2.15	0.47
1:A:149:PHE:HB3	3:C:40:PHE:HZ	1.77	0.47
1:A:238:GLY:O	1:A:239:ARG:HB2	2.15	0.47
1:A:94:ILE:O	1:A:96:MET:HG3	2.14	0.47
1:A:149:PHE:HD2	3:C:40:PHE:HZ	1.63	0.47
1:A:73:LEU:HD13	1:A:126:PHE:CZ	2.50	0.47
1:A:407:THR:O	1:A:411:VAL:HG23	2.15	0.47
2:B:13:GLU:O	2:B:16:GLU:HB3	2.16	0.46
1:A:69:THR:O	1:A:70:LEU:HG	2.15	0.46
1:A:83:ILE:HG23	1:A:330:PHE:CZ	2.51	0.46
1:A:112:LYS:O	1:A:115:SER:HB3	2.15	0.46
1:A:253:TYR:HD1	1:A:256:ASN:HD22	1.64	0.46
1:A:52:ILE:CD1	1:A:132:PHE:HA	2.45	0.46
1:A:2:LYS:O	1:A:3:LYS:HG2	2.16	0.46
1:A:5:ILE:N	1:A:6:PRO:CD	2.78	0.46
1:A:98:LEU:O	1:A:100:ILE:N	2.49	0.46
1:A:14:VAL:CG2	1:A:159:GLU:HB3	2.45	0.46
1:A:90:GLY:C	1:A:92:GLY:H	2.18	0.46
1:A:13:GLU:CG	1:A:14:VAL:H	2.23	0.46
1:A:35:VAL:O	1:A:39:ILE:HG12	2.16	0.46
1:A:206:ASN:HB3	1:A:209:TYR:HD1	1.81	0.46
1:A:46:TYR:HB3	1:A:146:GLN:NE2	2.21	0.46
1:A:304:TYR:CD1	1:A:304:TYR:C	2.89	0.45
1:A:256:ASN:N	1:A:258:PRO:HD2	2.31	0.45
1:A:328:VAL:HG12	1:A:332:ILE:HD11	1.98	0.45
1:A:381:SER:O	1:A:384:PHE:HB3	2.16	0.45
1:A:391:ILE:O	1:A:395:ILE:HG22	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:64:ILE:O	2:B:65:LEU:HD23	2.16	0.45
1:A:338:THR:HG22	1:A:339:GLY:N	2.31	0.45
1:A:189:PRO:C	1:A:191:GLY:H	2.19	0.45
2:B:21:VAL:C	2:B:23:LEU:H	2.18	0.45
1:A:22:THR:OG1	1:A:25:GLU:HG3	2.15	0.45
1:A:85:MET:HE1	1:A:111:GLN:N	2.32	0.45
1:A:195:LYS:HD3	1:A:209:TYR:HE2	1.82	0.45
1:A:76:GLY:O	1:A:79:VAL:N	2.50	0.45
1:A:194:TRP:HA	1:A:194:TRP:CE3	2.52	0.45
1:A:342:PRO:HG3	1:A:374:ILE:HD11	1.99	0.45
1:A:46:TYR:CB	1:A:146:GLN:HE22	2.22	0.45
1:A:92:GLY:O	1:A:95:GLN:HG3	2.16	0.45
3:C:44:GLU:HA	3:C:47:LEU:HB3	1.99	0.45
1:A:98:LEU:O	1:A:100:ILE:O	2.35	0.45
2:B:35:ALA:HA	2:B:38:LYS:HE3	1.99	0.45
1:A:3:LYS:C	1:A:5:ILE:H	2.19	0.44
1:A:77:PRO:HG3	1:A:119:CYS:HA	1.98	0.44
1:A:211:ALA:HA	1:A:214:ILE:HD12	1.99	0.44
1:A:281:ILE:N	1:A:281:ILE:CD1	2.80	0.44
1:A:370:LEU:C	1:A:372:ARG:N	2.70	0.44
3:C:30:PRO:O	3:C:34:ILE:HG12	2.18	0.44
1:A:74:GLY:O	1:A:76:GLY:N	2.50	0.44
1:A:89:VAL:HG12	1:A:89:VAL:O	2.17	0.44
1:A:195:LYS:HD3	1:A:209:TYR:CE2	2.52	0.44
1:A:423:VAL:C	1:A:425:GLU:N	2.71	0.44
2:B:24:VAL:HG12	2:B:24:VAL:O	2.17	0.44
1:A:166:ILE:HG22	1:A:167:GLY:H	1.83	0.44
1:A:265:LEU:O	1:A:269:ILE:HG13	2.18	0.44
1:A:34:LEU:HD11	1:A:38:PHE:CZ	2.52	0.44
1:A:42:CYS:SG	2:B:55:HIS:HA	2.58	0.44
1:A:51:GLN:CD	1:A:51:GLN:H	2.20	0.44
2:B:29:THR:O	2:B:32:GLU:N	2.51	0.44
1:A:86:GLN:NE2	1:A:107:PHE:HZ	2.15	0.44
1:A:234:PRO:CB	1:A:357:LYS:HB2	2.46	0.44
1:A:188:GLY:CA	2:B:52:TYR:HE1	2.30	0.44
1:A:259:VAL:HG12	1:A:406:LEU:HD21	2.00	0.44
1:A:240:ILE:O	1:A:241:LYS:HG3	2.18	0.44
1:A:405:LEU:O	1:A:406:LEU:C	2.55	0.44
1:A:14:VAL:CG1	3:C:30:PRO:HG3	2.48	0.44
1:A:16:LEU:H	1:A:16:LEU:CD2	2.11	0.44
1:A:149:PHE:HB3	3:C:40:PHE:CE2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:374:ILE:O	1:A:375:PRO:C	2.56	0.44
1:A:376:PRO:HB3	2:B:17:GLU:HB3	2.00	0.44
1:A:14:VAL:HA	3:C:28:VAL:O	2.17	0.43
1:A:187:LEU:HD21	1:A:193:LEU:CD2	2.48	0.43
1:A:417:GLN:HE21	1:A:417:GLN:HB2	1.52	0.43
1:A:88:LEU:O	1:A:90:GLY:N	2.50	0.43
1:A:135:LEU:HD12	1:A:135:LEU:N	2.33	0.43
1:A:187:LEU:HD21	1:A:193:LEU:HD23	2.01	0.43
1:A:169:GLY:O	1:A:170:ILE:C	2.54	0.43
1:A:270:GLN:HE22	1:A:301:SER:HB3	1.82	0.43
3:C:49:TYR:CD1	3:C:49:TYR:N	2.85	0.43
1:A:197:LEU:O	1:A:200:LEU:HB2	2.19	0.43
1:A:384:PHE:O	1:A:385:VAL:C	2.56	0.43
1:A:359:PHE:O	1:A:360:ARG:CB	2.66	0.43
1:A:350:GLY:O	1:A:353:GLY:N	2.51	0.43
1:A:360:ARG:O	1:A:366:ILE:HG13	2.19	0.43
1:A:153:ILE:HD11	3:C:40:PHE:CD1	2.53	0.43
1:A:8:LEU:HD23	1:A:8:LEU:C	2.39	0.43
1:A:14:VAL:HG21	1:A:159:GLU:CB	2.48	0.43
1:A:71:ILE:O	1:A:71:ILE:CG2	2.67	0.43
1:A:166:ILE:HG22	1:A:167:GLY:N	2.34	0.43
1:A:194:TRP:HA	1:A:194:TRP:HE3	1.84	0.43
1:A:231:VAL:CB	1:A:249:ILE:HD11	2.48	0.43
1:A:247:TYR:CD1	1:A:248:PRO:N	2.87	0.43
1:A:314:ILE:O	1:A:318:VAL:HG23	2.18	0.43
1:A:331:GLY:O	1:A:334:TRP:HB3	2.18	0.43
3:C:31:GLU:H	3:C:31:GLU:CD	2.20	0.43
1:A:118:MET:O	1:A:122:GLU:HG2	2.19	0.43
1:A:230:ARG:HG3	1:A:230:ARG:HH11	1.82	0.43
1:A:341:ASP:O	1:A:341:ASP:OD1	2.37	0.43
1:A:370:LEU:C	1:A:372:ARG:H	2.22	0.43
1:A:5:ILE:O	1:A:9:GLU:HG3	2.19	0.42
1:A:322:ALA:O	1:A:325:ILE:HG12	2.19	0.42
2:B:66:LYS:O	2:B:66:LYS:HG2	2.17	0.42
1:A:75:ILE:HD13	1:A:122:GLU:OE2	2.19	0.42
3:C:30:PRO:HA	3:C:33:VAL:CG2	2.49	0.42
1:A:384:PHE:O	1:A:387:PHE:HB3	2.19	0.42
2:B:30:LYS:HD3	2:B:30:LYS:HA	1.91	0.42
1:A:13:GLU:O	3:C:28:VAL:HB	2.19	0.42
1:A:51:GLN:OE1	1:A:51:GLN:N	2.52	0.42
1:A:83:ILE:HG12	1:A:261:LEU:HD22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:34:LEU:O	2:B:38:LYS:HG3	2.19	0.42
1:A:26:LYS:HA	1:A:164:TYR:O	2.20	0.42
1:A:85:MET:HG3	1:A:111:GLN:CG	2.49	0.42
1:A:310:ILE:O	1:A:312:ASP:N	2.44	0.42
1:A:55:ILE:O	1:A:69:THR:CB	2.68	0.42
1:A:202:GLN:OE1	1:A:202:GLN:HA	2.20	0.42
1:A:225:TYR:C	1:A:225:TYR:HD2	2.22	0.42
1:A:317:ILE:HD12	1:A:317:ILE:H	1.85	0.42
1:A:410:ILE:O	1:A:414:MET:HB2	2.20	0.42
2:B:56:VAL:HB	2:B:57:PRO:CD	2.50	0.42
1:A:166:ILE:HD13	1:A:418:LEU:HD21	2.01	0.42
2:B:35:ALA:O	2:B:38:LYS:HB2	2.20	0.42
1:A:130:GLY:HA3	1:A:278:ARG:NH1	2.34	0.41
1:A:423:VAL:C	1:A:425:GLU:H	2.22	0.41
2:B:14:PHE:CD1	2:B:14:PHE:C	2.94	0.41
1:A:274:LEU:N	1:A:287:TYR:CZ	2.89	0.41
1:A:108:GLN:O	1:A:109:GLY:C	2.59	0.41
1:A:219:VAL:HG12	1:A:220:PHE:N	2.36	0.41
1:A:18:VAL:HG12	1:A:18:VAL:O	2.20	0.41
1:A:5:ILE:HA	1:A:8:LEU:HB3	2.02	0.41
1:A:157:LEU:HA	1:A:160:ILE:HD12	2.03	0.41
1:A:6:PRO:HG2	1:A:7:ILE:H	1.86	0.41
1:A:313:PRO:O	1:A:317:ILE:HD13	2.21	0.41
3:C:31:GLU:OE2	3:C:31:GLU:N	2.37	0.41
1:A:375:PRO:HB2	1:A:376:PRO:HD3	2.02	0.41
2:B:16:GLU:O	2:B:20:ARG:HD3	2.20	0.41
1:A:52:ILE:HD11	1:A:132:PHE:HA	2.03	0.41
1:A:147:ILE:HG22	1:A:148:ALA:N	2.35	0.41
1:A:183:PHE:CZ	2:B:45:SER:HA	2.56	0.41
1:A:224:VAL:HG12	2:B:33:TYR:CE1	2.56	0.41
2:B:14:PHE:CE1	2:B:18:CYS:SG	3.14	0.41
1:A:91:SER:C	1:A:93:ILE:H	2.23	0.41
1:A:156:TYR:HB3	3:C:33:VAL:HG11	2.03	0.41
1:A:342:PRO:HG3	1:A:374:ILE:CD1	2.51	0.41
1:A:393:ASN:ND2	1:A:402:THR:H	2.07	0.41
2:B:9:GLU:O	2:B:12:LYS:HB2	2.21	0.40
1:A:181:THR:HG21	1:A:399:GLY:HA2	2.04	0.40
1:A:302:THR:HA	1:A:303:PRO:HD3	1.83	0.40
1:A:310:ILE:C	1:A:312:ASP:H	2.23	0.40
1:A:230:ARG:HD3	1:A:248:PRO:HB2	2.04	0.40
1:A:333:PHE:N	1:A:333:PHE:CD1	2.89	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	420/426 (99%)	290 (69%)	86 (20%)	44 (10%)	0 7
2	B	63/74 (85%)	51 (81%)	7 (11%)	5 (8%)	1 11
3	C	30/53 (57%)	22 (73%)	6 (20%)	2 (7%)	1 15
All	All	513/553 (93%)	363 (71%)	99 (19%)	51 (10%)	0 8

All (51) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	53	PRO
1	A	71	ILE
1	A	72	THR
1	A	75	ILE
1	A	99	SER
1	A	246	LYS
1	A	338	THR
1	A	360	ARG
1	A	70	LEU
1	A	145	ILE
1	A	147	ILE
1	A	188	GLY
1	A	189	PRO
1	A	237	HIS
1	A	245	GLY
1	A	311	SER
1	A	339	GLY
1	A	359	PHE
1	A	385	VAL
1	A	396	GLY
2	B	55	HIS

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Mol	Chain	Res	Type
1	A	363	GLU
1	A	4	LEU
1	A	19	LYS
1	A	67	GLY
1	A	77	PRO
1	A	102	GLU
1	A	240	ILE
1	A	346	ALA
1	A	367	GLU
1	A	374	ILE
1	A	429	ALA
1	A	430	ILE
2	B	38	LYS
1	A	172	LEU
1	A	205	PRO
1	A	244	VAL
1	A	288	GLU
1	A	356	ILE
2	B	30	LYS
3	C	51	ARG
1	A	6	PRO
3	C	41	VAL
2	B	24	VAL
2	B	56	VAL
1	A	89	VAL
1	A	137	PRO
1	A	211	ALA
1	A	375	PRO
1	A	170	ILE
1	A	248	PRO

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	341/345 (99%)	310 (91%)	31 (9%)	9 39

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	57/66 (86%)	53 (93%)	4 (7%)	15	48
3	C	28/45 (62%)	25 (89%)	3 (11%)	6	32
All	All	426/456 (93%)	388 (91%)	38 (9%)	9	40

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	16	LEU
1	A	21	ILE
1	A	27	LEU
1	A	29	TRP
1	A	51	GLN
1	A	55	ILE
1	A	126	PHE
1	A	172	LEU
1	A	197	LEU
1	A	202	GLN
1	A	213	ILE
1	A	219	VAL
1	A	225	TYR
1	A	247	TYR
1	A	248	PRO
1	A	254	VAL
1	A	257	ILE
1	A	259	VAL
1	A	306	LEU
1	A	314	ILE
1	A	336	GLU
1	A	340	LEU
1	A	344	SER
1	A	359	PHE
1	A	363	GLU
1	A	369	ARG
1	A	375	PRO
1	A	380	MET
1	A	402	THR
1	A	417	GLN
2	B	19	ARG
2	B	20	ARG
2	B	30	LYS

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Mol	Chain	Res	Type
2	B	42	LEU
3	C	33	VAL
3	C	44	GLU
3	C	51	ARG

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

Mol	Chain	Res	Type
1	A	111	GLN
1	A	146	GLN
1	A	256	ASN
1	A	268	ASN
1	A	315	HIS
1	A	393	ASN
1	A	417	GLN
2	B	5	ASN
2	B	6	GLN
2	B	10	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no 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	422/426 (99%)	-0.04	24 (5%) 23 14	48, 116, 167, 198	0
2	B	65/74 (87%)	-0.51	0 100 100	53, 120, 165, 180	0
3	C	32/53 (60%)	-0.02	1 (3%) 49 33	111, 144, 176, 188	0
All	All	519/553 (93%)	-0.10	25 (4%) 30 19	48, 119, 170, 198	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	47	THR	6.6
3	C	51	ARG	4.6
1	A	48	ALA	3.9
1	A	306	LEU	3.4
1	A	2	LYS	3.4
1	A	351	SER	3.4
1	A	293	VAL	3.3
1	A	243	ALA	3.3
1	A	96	MET	3.3
1	A	95	GLN	3.2
1	A	245	GLY	2.9
1	A	246	LYS	2.8
1	A	354	MET	2.7
1	A	71	ILE	2.7
1	A	284	LEU	2.7
1	A	283	ILE	2.6
1	A	305	GLY	2.6
1	A	238	GLY	2.6
1	A	132	PHE	2.4
1	A	299	TYR	2.3
1	A	49	GLY	2.3
1	A	315	HIS	2.1
1	A	74	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	73	LEU	2.1
1	A	72	THR	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.