



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 10:01 PM GMT

PDB ID : 5C6N  
Title : protein A  
Authors : Lu, M.  
Deposited on : 2015-06-23  
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : unknown  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026982  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20026982

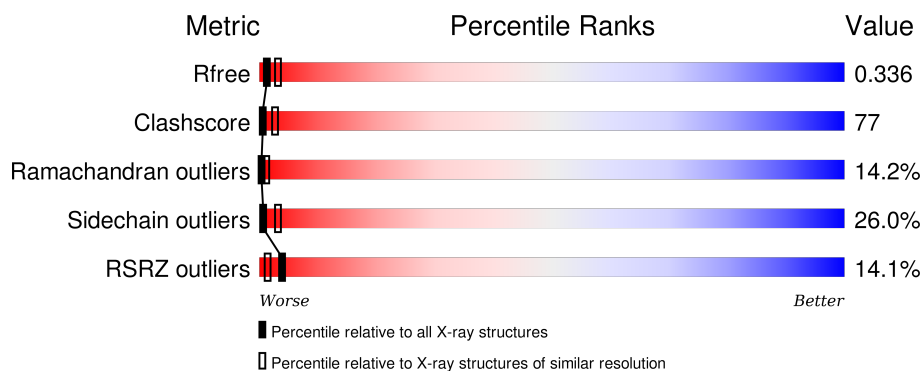
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	463	

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3413 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 BH2163 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	446	Total	C	N	O	S	0	0	0
			3413	2263	549	573	28			

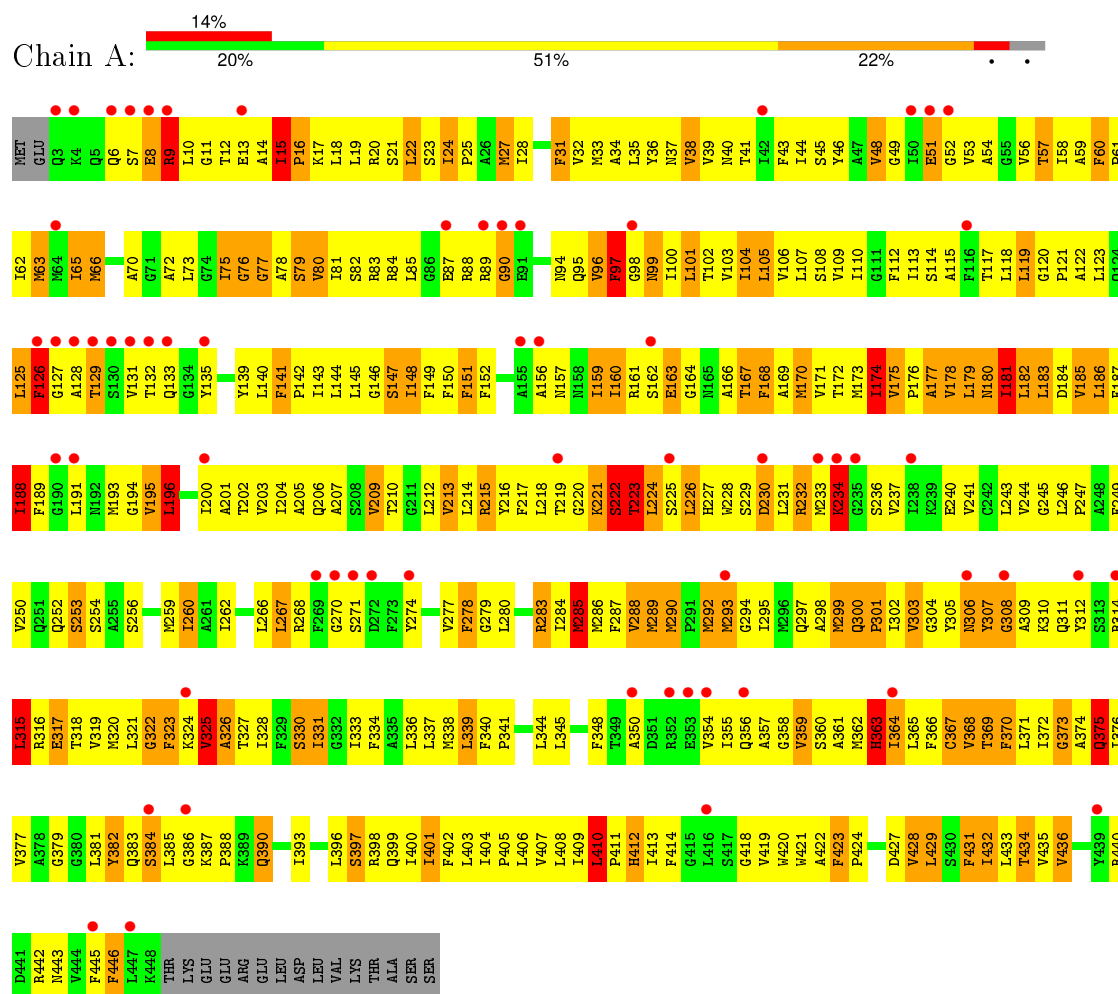
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	40	ASN	ASP	conflict	UNP Q9KAX3

### 3 Residue-property plots

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

- Molecule 1: BH2163 protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	89.68Å 93.42Å 101.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00 20.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	90.0 (20.00-3.00) 87.6 (20.00-3.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.70 (at 2.98Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.280 , 0.300 0.327 , 0.336	Depositor DCC
$R_{free}$ test set	766 reflections (4.97%)	DCC
Wilson B-factor (Å <sup>2</sup> )	72.6	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.09 , -9.2	EDS
Estimated twinning fraction	0.145 for k,h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.30$ , $\langle L^2 \rangle = 0.14$	Xtriage
Outliers	7 of 15465 reflections (0.045%)	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	3413	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	202.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.22% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 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	0.60	0/3481	0.93	2/4714 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	308	GLY	N-CA-C	-6.09	97.87	113.10
1	A	315	LEU	CB-CG-CD1	6.05	121.28	111.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	125	LEU	Peptide
1	A	15	ILE	Peptide
1	A	195	VAL	Peptide
1	A	306	ASN	Peptide
1	A	307	TYR	Peptide

### 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	3413	0	3610	544	0
All	All	3413	0	3610	544	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 77.

All (544) 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:173:MET:O	1:A:176:PRO:HD2	1.33	1.27
1:A:224:LEU:O	1:A:226:LEU:HG	1.44	1.14
1:A:404:ILE:HG13	1:A:405:PRO:CD	1.75	1.14
1:A:289:MET:HA	1:A:289:MET:CE	1.77	1.13
1:A:404:ILE:CG1	1:A:405:PRO:HD3	1.78	1.12
1:A:289:MET:HA	1:A:289:MET:HE2	1.31	1.11
1:A:78:ALA:HA	1:A:162:SER:HB2	1.20	1.11
1:A:364:ILE:HG23	1:A:367:CYS:SG	1.94	1.07
1:A:78:ALA:HA	1:A:162:SER:CB	1.84	1.07
1:A:59:ALA:HB1	1:A:140:LEU:HD13	1.30	1.07
1:A:247:PRO:HG3	1:A:388:PRO:HB2	1.34	1.06
1:A:407:VAL:O	1:A:411:PRO:HD2	1.54	1.05
1:A:15:ILE:HD13	1:A:15:ILE:O	1.56	1.04
1:A:222:SER:O	1:A:226:LEU:HD11	1.58	1.02
1:A:149:PHE:HD2	1:A:207:ALA:HA	1.21	1.01
1:A:420:TRP:O	1:A:424:PRO:HD2	1.59	1.00
1:A:173:MET:O	1:A:175:VAL:N	1.95	0.99
1:A:364:ILE:HG22	1:A:420:TRP:HZ2	1.27	0.98
1:A:156:ALA:O	1:A:160:ILE:HG23	1.64	0.96
1:A:298:ALA:O	1:A:301:PRO:HD2	1.66	0.96
1:A:300:GLN:HB3	1:A:301:PRO:HD3	1.48	0.95
1:A:187:PHE:HA	1:A:191:LEU:HD12	1.50	0.93
1:A:361:ALA:O	1:A:365:LEU:HB2	1.69	0.93
1:A:149:PHE:CD2	1:A:207:ALA:HA	2.03	0.93
1:A:24:ILE:HG13	1:A:25:PRO:HD3	1.52	0.92
1:A:59:ALA:HB1	1:A:140:LEU:CD1	1.99	0.92
1:A:183:LEU:HD13	1:A:187:PHE:HE1	1.32	0.91
1:A:19:LEU:HD12	1:A:321:LEU:HD23	1.54	0.89
1:A:115:ALA:HB2	1:A:144:LEU:HB3	1.55	0.89
1:A:87:GLU:HB2	1:A:308:GLY:HA2	1.56	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:GLY:HA3	1:A:241:VAL:HG13	1.53	0.87
1:A:106:VAL:O	1:A:110:ILE:HD12	1.74	0.87
1:A:7:SER:HB2	1:A:223:THR:O	1.74	0.86
1:A:148:ILE:HG13	1:A:149:PHE:N	1.91	0.85
1:A:13:GLU:OE2	1:A:306:ASN:ND2	2.09	0.85
1:A:35:LEU:O	1:A:38:VAL:HB	1.75	0.85
1:A:411:PRO:HD3	1:A:418:GLY:HA3	1.58	0.84
1:A:51:GLU:CD	1:A:52:GLY:H	1.79	0.84
1:A:60:PHE:HB3	1:A:61:PRO:HD3	1.59	0.84
1:A:402:PHE:CD1	1:A:429:LEU:HD21	2.13	0.83
1:A:240:GLU:HA	1:A:243:LEU:HD12	1.59	0.83
1:A:299:MET:CE	1:A:322:GLY:HA2	2.08	0.83
1:A:364:ILE:CG2	1:A:367:CYS:SG	2.67	0.83
1:A:78:ALA:CA	1:A:162:SER:HB2	2.07	0.82
1:A:36:TYR:HD1	1:A:181:ILE:CD1	1.91	0.82
1:A:308:GLY:O	1:A:310:LYS:N	2.13	0.81
1:A:314:ARG:O	1:A:317:GLU:HB3	1.78	0.81
1:A:406:LEU:HD22	1:A:422:ALA:HA	1.62	0.81
1:A:315:LEU:HD12	1:A:316:ARG:N	1.95	0.81
1:A:20:ARG:O	1:A:24:ILE:HG23	1.80	0.81
1:A:361:ALA:O	1:A:365:LEU:CB	2.28	0.80
1:A:128:ALA:HB1	1:A:133:GLN:OE1	1.79	0.80
1:A:94:ASN:O	1:A:231:LEU:HD21	1.79	0.80
1:A:299:MET:HE1	1:A:322:GLY:HA2	1.61	0.80
1:A:396:LEU:O	1:A:401:ILE:HG12	1.82	0.79
1:A:323:PHE:CD1	1:A:377:VAL:HG11	2.18	0.78
1:A:108:SER:OG	1:A:148:ILE:HA	1.83	0.78
1:A:224:LEU:O	1:A:226:LEU:CG	2.30	0.78
1:A:310:LYS:O	1:A:312:TYR:HD1	1.68	0.77
1:A:195:VAL:N	1:A:196:LEU:HD12	1.99	0.77
1:A:180:ASN:OD1	1:A:202:THR:HG23	1.84	0.77
1:A:364:ILE:HG22	1:A:420:TRP:CZ2	2.15	0.77
1:A:173:MET:O	1:A:176:PRO:CD	2.26	0.77
1:A:16:PRO:HD2	1:A:17:LYS:H	1.49	0.77
1:A:163:GLU:HG3	1:A:164:GLY:H	1.50	0.76
1:A:31:PHE:HB2	1:A:290:MET:SD	2.26	0.76
1:A:17:LYS:HA	1:A:20:ARG:HB2	1.66	0.76
1:A:325:VAL:O	1:A:328:ILE:HG22	1.86	0.75
1:A:407:VAL:O	1:A:411:PRO:CD	2.32	0.74
1:A:325:VAL:HG13	1:A:326:ALA:H	1.52	0.74
1:A:15:ILE:O	1:A:15:ILE:CD1	2.35	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:ILE:HG22	1:A:182:LEU:HD23	1.68	0.74
1:A:184:ASP:OD1	1:A:201:ALA:HB1	1.88	0.73
1:A:390:GLN:HA	1:A:393:ILE:HD12	1.68	0.73
1:A:299:MET:HE1	1:A:322:GLY:CA	2.20	0.72
1:A:397:SER:OG	1:A:398:ARG:N	2.20	0.72
1:A:183:LEU:HD13	1:A:187:PHE:CE1	2.22	0.71
1:A:350:ALA:HB3	1:A:354:VAL:HG21	1.71	0.71
1:A:139:TYR:O	1:A:142:PRO:HG2	1.91	0.71
1:A:334:PHE:HE1	1:A:367:CYS:HA	1.55	0.70
1:A:184:ASP:OD1	1:A:201:ALA:CB	2.38	0.70
1:A:382:TYR:CE1	1:A:390:GLN:HB2	2.26	0.70
1:A:420:TRP:HZ3	1:A:421:TRP:CE3	2.09	0.70
1:A:125:LEU:O	1:A:126:PHE:HB2	1.92	0.70
1:A:224:LEU:C	1:A:226:LEU:HG	2.12	0.70
1:A:247:PRO:HG2	1:A:383:GLN:NE2	2.07	0.70
1:A:365:LEU:HD23	1:A:366:PHE:CE1	2.26	0.70
1:A:110:ILE:HA	1:A:113:ILE:HD12	1.74	0.70
1:A:279:GLY:O	1:A:283:ARG:HD2	1.92	0.69
1:A:365:LEU:HD23	1:A:366:PHE:CZ	2.26	0.69
1:A:218:LEU:HA	1:A:221:LYS:HD3	1.75	0.69
1:A:410:LEU:HB3	1:A:418:GLY:HA2	1.74	0.69
1:A:88:ARG:O	1:A:90:GLY:N	2.22	0.69
1:A:410:LEU:O	1:A:414:PHE:N	2.26	0.68
1:A:129:THR:HB	1:A:132:THR:HG22	1.76	0.68
1:A:78:ALA:CA	1:A:162:SER:CB	2.69	0.68
1:A:397:SER:HB2	1:A:402:PHE:CD1	2.29	0.68
1:A:364:ILE:CG2	1:A:420:TRP:HZ2	2.03	0.67
1:A:179:LEU:CD1	1:A:205:ALA:HA	2.23	0.67
1:A:423:PHE:O	1:A:424:PRO:C	2.32	0.67
1:A:187:PHE:HB3	1:A:193:MET:HB2	1.76	0.67
1:A:320:MET:HG3	1:A:445:PHE:CE1	2.30	0.66
1:A:170:MET:O	1:A:174:ILE:HG23	1.94	0.66
1:A:249:PHE:O	1:A:253:SER:OG	2.08	0.66
1:A:222:SER:O	1:A:226:LEU:CD1	2.39	0.66
1:A:334:PHE:HB2	1:A:369:THR:HB	1.76	0.66
1:A:149:PHE:H	1:A:149:PHE:HD1	1.38	0.66
1:A:420:TRP:O	1:A:424:PRO:CD	2.40	0.66
1:A:300:GLN:HB3	1:A:301:PRO:CD	2.26	0.65
1:A:227:HIS:CD2	1:A:231:LEU:HB2	2.31	0.65
1:A:19:LEU:HD22	1:A:298:ALA:HB3	1.79	0.65
1:A:144:LEU:O	1:A:147:SER:OG	2.13	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:400:ILE:O	1:A:402:PHE:N	2.29	0.65
1:A:181:ILE:H	1:A:181:ILE:HD13	1.60	0.65
1:A:148:ILE:HG13	1:A:149:PHE:H	1.62	0.65
1:A:100:ILE:HG22	1:A:104:ILE:HD11	1.78	0.65
1:A:323:PHE:HD1	1:A:377:VAL:HG11	1.59	0.65
1:A:112:PHE:HD1	1:A:145:LEU:O	1.80	0.65
1:A:299:MET:CE	1:A:322:GLY:CA	2.73	0.65
1:A:66:MET:HB2	1:A:151:PHE:CE2	2.32	0.65
1:A:11:GLY:HA2	1:A:305:TYR:CE1	2.32	0.65
1:A:141:PHE:O	1:A:145:LEU:HG	1.97	0.64
1:A:19:LEU:HD23	1:A:298:ALA:HB1	1.80	0.64
1:A:87:GLU:H	1:A:308:GLY:CA	2.11	0.64
1:A:288:VAL:O	1:A:288:VAL:HG13	1.96	0.64
1:A:320:MET:HB2	1:A:445:PHE:HE1	1.63	0.64
1:A:364:ILE:C	1:A:420:TRP:HE1	2.01	0.64
1:A:167:THR:O	1:A:170:MET:HB3	1.97	0.64
1:A:53:VAL:O	1:A:57:THR:OG1	2.15	0.64
1:A:289:MET:CA	1:A:289:MET:CE	2.66	0.63
1:A:82:SER:HB2	1:A:304:GLY:HA3	1.79	0.63
1:A:297:GLN:HA	1:A:300:GLN:HB2	1.81	0.63
1:A:304:GLY:O	1:A:307:TYR:O	2.15	0.63
1:A:409:ILE:CG2	1:A:409:ILE:O	2.47	0.63
1:A:36:TYR:HA	1:A:181:ILE:HD11	1.81	0.63
1:A:31:PHE:HD2	1:A:32:VAL:HG23	1.64	0.63
1:A:410:LEU:HB3	1:A:418:GLY:CA	2.29	0.63
1:A:159:ILE:HD11	1:A:217:PHE:CZ	2.33	0.62
1:A:108:SER:O	1:A:148:ILE:HG22	1.99	0.62
1:A:58:ILE:O	1:A:61:PRO:HD2	1.98	0.62
1:A:331:ILE:C	1:A:331:ILE:HD12	2.20	0.62
1:A:46:TYR:N	1:A:48:VAL:O	2.32	0.62
1:A:299:MET:HE3	1:A:322:GLY:HA2	1.81	0.62
1:A:195:VAL:H	1:A:196:LEU:HD12	1.64	0.62
1:A:87:GLU:H	1:A:308:GLY:HA3	1.63	0.62
1:A:84:ARG:HA	1:A:87:GLU:O	1.99	0.62
1:A:106:VAL:O	1:A:110:ILE:CD1	2.46	0.62
1:A:16:PRO:CD	1:A:17:LYS:H	2.09	0.62
1:A:97:PHE:HE1	1:A:101:LEU:HD11	1.65	0.62
1:A:167:THR:O	1:A:170:MET:CB	2.47	0.62
1:A:65:ILE:HG22	1:A:66:MET:N	2.14	0.62
1:A:10:LEU:HD23	1:A:12:THR:HB	1.81	0.62
1:A:36:TYR:CD1	1:A:181:ILE:CD1	2.78	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:299:MET:HA	1:A:302:ILE:HB	1.82	0.61
1:A:322:GLY:O	1:A:325:VAL:CG1	2.48	0.61
1:A:370:PHE:N	1:A:370:PHE:CD2	2.68	0.61
1:A:222:SER:O	1:A:224:LEU:N	2.34	0.61
1:A:363:HIS:O	1:A:365:LEU:N	2.32	0.61
1:A:373:GLY:O	1:A:377:VAL:HG23	1.99	0.61
1:A:223:THR:HG22	1:A:224:LEU:HD12	1.82	0.61
1:A:407:VAL:HG12	1:A:411:PRO:HG2	1.83	0.61
1:A:65:ILE:CG2	1:A:66:MET:N	2.63	0.61
1:A:152:PHE:HE2	1:A:213:VAL:HB	1.66	0.61
1:A:62:ILE:N	1:A:62:ILE:HD12	2.15	0.61
1:A:385:LEU:O	1:A:387:LYS:N	2.32	0.61
1:A:24:ILE:HG13	1:A:25:PRO:CD	2.29	0.61
1:A:303:VAL:HG12	1:A:303:VAL:O	2.01	0.61
1:A:46:TYR:HB3	1:A:189:PHE:CZ	2.35	0.61
1:A:364:ILE:O	1:A:367:CYS:SG	2.56	0.60
1:A:408:LEU:N	1:A:408:LEU:HD23	2.15	0.60
1:A:247:PRO:HG3	1:A:388:PRO:CB	2.22	0.60
1:A:140:LEU:HA	1:A:143:ILE:HD12	1.82	0.60
1:A:173:MET:C	1:A:175:VAL:H	2.03	0.60
1:A:18:LEU:O	1:A:22:LEU:HB2	2.02	0.60
1:A:19:LEU:CD2	1:A:298:ALA:CB	2.80	0.60
1:A:227:HIS:HB3	1:A:231:LEU:HB2	1.84	0.60
1:A:41:THR:O	1:A:44:ILE:N	2.34	0.60
1:A:25:PRO:HB3	1:A:166:ALA:O	2.02	0.60
1:A:374:ALA:O	1:A:376:ILE:N	2.35	0.60
1:A:97:PHE:O	1:A:101:LEU:HD23	2.01	0.59
1:A:70:ALA:HA	1:A:73:LEU:HD12	1.83	0.59
1:A:370:PHE:HD2	1:A:370:PHE:H	1.48	0.59
1:A:66:MET:HB2	1:A:151:PHE:HE2	1.66	0.59
1:A:46:TYR:HB3	1:A:189:PHE:HZ	1.68	0.59
1:A:254:SER:HB3	1:A:399:GLN:HG2	1.85	0.59
1:A:254:SER:CB	1:A:399:GLN:HG2	2.33	0.59
1:A:94:ASN:HD21	1:A:227:HIS:HB2	1.66	0.59
1:A:156:ALA:HB1	1:A:213:VAL:HG11	1.85	0.59
1:A:88:ARG:C	1:A:90:GLY:H	2.07	0.59
1:A:323:PHE:CE1	1:A:377:VAL:CG1	2.86	0.58
1:A:303:VAL:HG12	1:A:385:LEU:HD21	1.85	0.58
1:A:177:ALA:O	1:A:178:VAL:C	2.42	0.58
1:A:420:TRP:HZ3	1:A:421:TRP:CD2	2.21	0.58
1:A:169:ALA:O	1:A:171:VAL:N	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:410:LEU:CB	1:A:418:GLY:HA2	2.33	0.58
1:A:322:GLY:O	1:A:325:VAL:HG13	2.03	0.58
1:A:79:SER:HB2	1:A:383:GLN:O	2.03	0.58
1:A:402:PHE:CG	1:A:429:LEU:HD21	2.38	0.58
1:A:103:VAL:O	1:A:106:VAL:HB	2.04	0.58
1:A:62:ILE:HD12	1:A:62:ILE:H	1.69	0.57
1:A:25:PRO:HA	1:A:170:MET:HE1	1.86	0.57
1:A:24:ILE:HD12	1:A:24:ILE:C	2.25	0.57
1:A:72:ALA:CB	1:A:249:PHE:HD2	2.17	0.57
1:A:149:PHE:N	1:A:149:PHE:CD1	2.69	0.57
1:A:404:ILE:C	1:A:406:LEU:H	2.08	0.57
1:A:310:LYS:O	1:A:312:TYR:CD1	2.55	0.57
1:A:409:ILE:O	1:A:409:ILE:HG22	2.04	0.57
1:A:38:VAL:HG12	1:A:39:VAL:N	2.19	0.57
1:A:149:PHE:CD2	1:A:207:ALA:CA	2.83	0.57
1:A:96:VAL:HA	1:A:99:ASN:OD1	2.04	0.57
1:A:51:GLU:CD	1:A:52:GLY:N	2.55	0.57
1:A:390:GLN:O	1:A:393:ILE:HB	2.05	0.57
1:A:315:LEU:CD1	1:A:316:ARG:N	2.67	0.57
1:A:169:ALA:O	1:A:170:MET:C	2.42	0.56
1:A:10:LEU:HD23	1:A:12:THR:CB	2.35	0.56
1:A:80:VAL:HG12	1:A:81:ILE:N	2.19	0.56
1:A:410:LEU:HD21	1:A:421:TRP:HB2	1.86	0.56
1:A:188:ILE:HG23	1:A:189:PHE:N	2.20	0.56
1:A:43:PHE:O	1:A:46:TYR:HB2	2.06	0.56
1:A:115:ALA:HB2	1:A:144:LEU:CB	2.32	0.56
1:A:330:SER:HB3	1:A:370:PHE:HA	1.86	0.56
1:A:397:SER:HB2	1:A:402:PHE:CE1	2.41	0.56
1:A:43:PHE:HB3	1:A:188:ILE:HD12	1.86	0.56
1:A:23:SER:OG	1:A:24:ILE:N	2.37	0.56
1:A:19:LEU:HD11	1:A:299:MET:HE1	1.87	0.55
1:A:114:SER:HA	1:A:117:THR:HB	1.88	0.55
1:A:96:VAL:HG12	1:A:97:PHE:N	2.21	0.55
1:A:179:LEU:O	1:A:183:LEU:N	2.39	0.55
1:A:420:TRP:CZ3	1:A:421:TRP:CD2	2.94	0.55
1:A:383:GLN:C	1:A:385:LEU:H	2.09	0.55
1:A:370:PHE:HD2	1:A:370:PHE:N	2.03	0.55
1:A:423:PHE:HB2	1:A:424:PRO:CD	2.36	0.55
1:A:176:PRO:O	1:A:179:LEU:HB3	2.06	0.55
1:A:24:ILE:HA	1:A:27:MET:HB2	1.88	0.55
1:A:302:ILE:C	1:A:304:GLY:H	2.10	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:ASN:HD21	1:A:202:THR:HA	1.72	0.55
1:A:96:VAL:O	1:A:97:PHE:C	2.44	0.55
1:A:443:ASN:O	1:A:446:PHE:CD2	2.59	0.55
1:A:100:ILE:HG22	1:A:104:ILE:CD1	2.37	0.55
1:A:330:SER:HB3	1:A:370:PHE:O	2.05	0.54
1:A:187:PHE:O	1:A:191:LEU:HB2	2.07	0.54
1:A:233:MET:HG3	1:A:234:LYS:HG3	1.88	0.54
1:A:102:THR:HG21	1:A:233:MET:O	2.08	0.54
1:A:289:MET:HA	1:A:289:MET:HE3	1.82	0.54
1:A:13:GLU:OE2	1:A:314:ARG:HB2	2.07	0.54
1:A:19:LEU:CD2	1:A:298:ALA:HB1	2.37	0.54
1:A:367:CYS:SG	1:A:367:CYS:O	2.65	0.54
1:A:184:ASP:O	1:A:187:PHE:N	2.41	0.54
1:A:19:LEU:HD22	1:A:298:ALA:CB	2.36	0.54
1:A:320:MET:HB2	1:A:445:PHE:CE1	2.43	0.54
1:A:183:LEU:O	1:A:186:LEU:HB2	2.08	0.54
1:A:59:ALA:CB	1:A:140:LEU:HD13	2.21	0.54
1:A:181:ILE:N	1:A:181:ILE:CD1	2.71	0.54
1:A:404:ILE:HG13	1:A:405:PRO:HD3	0.82	0.54
1:A:98:GLY:HA2	1:A:101:LEU:HG	1.90	0.54
1:A:233:MET:HG3	1:A:234:LYS:CG	2.38	0.54
1:A:160:ILE:HD13	1:A:168:PHE:CE1	2.43	0.53
1:A:63:MET:O	1:A:66:MET:HG3	2.07	0.53
1:A:97:PHE:C	1:A:97:PHE:CD1	2.82	0.53
1:A:174:ILE:O	1:A:178:VAL:HG22	2.08	0.53
1:A:292:MET:HE1	1:A:330:SER:HA	1.90	0.53
1:A:25:PRO:HA	1:A:170:MET:CE	2.38	0.53
1:A:76:GLY:HA2	1:A:244:VAL:HB	1.90	0.53
1:A:323:PHE:HE1	1:A:377:VAL:HG12	1.74	0.53
1:A:180:ASN:ND2	1:A:201:ALA:O	2.41	0.53
1:A:118:LEU:HA	1:A:121:PRO:HG2	1.89	0.53
1:A:181:ILE:H	1:A:181:ILE:CD1	2.22	0.53
1:A:215:ARG:O	1:A:217:PHE:N	2.42	0.53
1:A:404:ILE:O	1:A:406:LEU:N	2.41	0.53
1:A:18:LEU:HB3	1:A:302:ILE:HD13	1.90	0.53
1:A:36:TYR:HD1	1:A:181:ILE:HD13	1.71	0.53
1:A:15:ILE:HA	1:A:18:LEU:HB2	1.91	0.52
1:A:383:GLN:C	1:A:385:LEU:N	2.62	0.52
1:A:75:ILE:HG22	1:A:76:GLY:N	2.24	0.52
1:A:293:MET:CE	1:A:376:ILE:HD11	2.39	0.52
1:A:254:SER:HB3	1:A:399:GLN:CG	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:THR:O	1:A:322:GLY:N	2.39	0.52
1:A:96:VAL:O	1:A:100:ILE:HG12	2.10	0.52
1:A:101:LEU:HA	1:A:104:ILE:HD12	1.91	0.52
1:A:60:PHE:HB3	1:A:61:PRO:CD	2.35	0.52
1:A:396:LEU:O	1:A:401:ILE:CG1	2.55	0.52
1:A:40:ASN:ND2	1:A:184:ASP:OD2	2.42	0.52
1:A:145:LEU:N	1:A:146:GLY:HA3	2.24	0.52
1:A:188:ILE:HG23	1:A:189:PHE:H	1.74	0.52
1:A:178:VAL:O	1:A:181:ILE:HB	2.09	0.52
1:A:250:VAL:O	1:A:254:SER:N	2.42	0.52
1:A:220:GLY:O	1:A:222:SER:N	2.42	0.52
1:A:383:GLN:O	1:A:385:LEU:N	2.42	0.52
1:A:365:LEU:N	1:A:420:TRP:HE1	2.08	0.52
1:A:340:PHE:N	1:A:341:PRO:HD3	2.25	0.52
1:A:149:PHE:HB3	1:A:206:GLN:O	2.09	0.52
1:A:94:ASN:ND2	1:A:231:LEU:HD13	2.25	0.52
1:A:72:ALA:HB2	1:A:249:PHE:HD2	1.75	0.52
1:A:8:GLU:O	1:A:9:ARG:C	2.47	0.52
1:A:303:VAL:HG23	1:A:318:THR:HG21	1.92	0.51
1:A:112:PHE:CD1	1:A:145:LEU:O	2.61	0.51
1:A:160:ILE:CD1	1:A:168:PHE:CD1	2.94	0.51
1:A:295:ILE:O	1:A:298:ALA:HB3	2.09	0.51
1:A:411:PRO:O	1:A:412:HIS:C	2.46	0.51
1:A:400:ILE:HG13	1:A:401:ILE:H	1.76	0.51
1:A:166:ALA:O	1:A:169:ALA:HB3	2.11	0.51
1:A:170:MET:O	1:A:174:ILE:N	2.39	0.51
1:A:188:ILE:CG2	1:A:189:PHE:H	2.24	0.51
1:A:442:ARG:O	1:A:446:PHE:HB3	2.11	0.51
1:A:141:PHE:N	1:A:142:PRO:HD2	2.26	0.51
1:A:25:PRO:HA	1:A:170:MET:SD	2.50	0.51
1:A:247:PRO:CG	1:A:388:PRO:HB2	2.23	0.51
1:A:411:PRO:O	1:A:412:HIS:O	2.29	0.51
1:A:397:SER:HB2	1:A:402:PHE:HD1	1.75	0.51
1:A:227:HIS:HB3	1:A:231:LEU:HD13	1.93	0.51
1:A:35:LEU:O	1:A:36:TYR:C	2.48	0.50
1:A:105:LEU:HA	1:A:108:SER:HB2	1.93	0.50
1:A:163:GLU:HG3	1:A:164:GLY:N	2.21	0.50
1:A:315:LEU:HD12	1:A:316:ARG:H	1.76	0.50
1:A:369:THR:O	1:A:372:ILE:N	2.36	0.50
1:A:54:ALA:O	1:A:57:THR:N	2.44	0.50
1:A:19:LEU:HD23	1:A:298:ALA:CB	2.40	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:VAL:HA	1:A:240:GLU:HG3	1.93	0.50
1:A:131:VAL:HG23	1:A:132:THR:H	1.76	0.50
1:A:357:ALA:HA	1:A:360:SER:HB2	1.92	0.50
1:A:65:ILE:HG22	1:A:66:MET:H	1.77	0.50
1:A:46:TYR:CB	1:A:189:PHE:HZ	2.24	0.50
1:A:78:ALA:HA	1:A:162:SER:HB3	1.82	0.50
1:A:410:LEU:CD2	1:A:421:TRP:HB2	2.41	0.50
1:A:209:VAL:O	1:A:212:LEU:N	2.45	0.49
1:A:210:THR:HA	1:A:213:VAL:HG23	1.94	0.49
1:A:10:LEU:HB3	1:A:12:THR:H	1.76	0.49
1:A:148:ILE:CG1	1:A:149:PHE:CD1	2.95	0.49
1:A:306:ASN:HB3	1:A:315:LEU:HB3	1.93	0.49
1:A:337:LEU:HD21	1:A:366:PHE:CB	2.41	0.49
1:A:107:LEU:HA	1:A:110:ILE:HD13	1.95	0.49
1:A:148:ILE:CG1	1:A:149:PHE:N	2.71	0.49
1:A:297:GLN:O	1:A:301:PRO:HD3	2.12	0.49
1:A:325:VAL:HG22	1:A:326:ALA:N	2.28	0.49
1:A:7:SER:CB	1:A:223:THR:O	2.56	0.49
1:A:184:ASP:O	1:A:185:VAL:C	2.51	0.49
1:A:169:ALA:HA	1:A:172:THR:OG1	2.12	0.49
1:A:24:ILE:N	1:A:25:PRO:CD	2.75	0.49
1:A:129:THR:OG1	1:A:267:LEU:HD11	2.13	0.49
1:A:97:PHE:C	1:A:97:PHE:HD1	2.16	0.49
1:A:323:PHE:CE1	1:A:377:VAL:HG11	2.48	0.49
1:A:45:SER:HA	1:A:49:GLY:O	2.13	0.49
1:A:157:ASN:ND2	1:A:172:THR:OG1	2.45	0.49
1:A:31:PHE:CB	1:A:290:MET:SD	2.99	0.49
1:A:100:ILE:C	1:A:104:ILE:HD12	2.33	0.49
1:A:331:ILE:O	1:A:331:ILE:HD12	2.13	0.49
1:A:120:GLY:N	1:A:121:PRO:HD2	2.28	0.49
1:A:215:ARG:C	1:A:217:PHE:H	2.16	0.49
1:A:28:ILE:HB	1:A:170:MET:SD	2.53	0.48
1:A:163:GLU:CD	1:A:223:THR:HB	2.33	0.48
1:A:333:ILE:O	1:A:337:LEU:N	2.41	0.48
1:A:161:ARG:O	1:A:163:GLU:O	2.32	0.48
1:A:292:MET:CE	1:A:330:SER:HA	2.44	0.48
1:A:337:LEU:HD21	1:A:366:PHE:CG	2.49	0.48
1:A:303:VAL:CG1	1:A:303:VAL:O	2.60	0.48
1:A:358:GLY:O	1:A:362:MET:HG2	2.13	0.48
1:A:375:GLN:HG3	1:A:431:PHE:HA	1.96	0.48
1:A:433:LEU:O	1:A:436:VAL:HG12	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:ASP:OD1	1:A:201:ALA:HB3	2.12	0.48
1:A:18:LEU:HB3	1:A:302:ILE:CD1	2.44	0.48
1:A:283:ARG:O	1:A:287:PHE:HB2	2.14	0.48
1:A:302:ILE:HG22	1:A:303:VAL:N	2.27	0.48
1:A:436:VAL:O	1:A:440:ARG:HB2	2.14	0.48
1:A:178:VAL:O	1:A:182:LEU:HG	2.13	0.48
1:A:410:LEU:O	1:A:414:PHE:HB2	2.13	0.48
1:A:330:SER:OG	1:A:373:GLY:N	2.45	0.48
1:A:36:TYR:O	1:A:37:ASN:C	2.52	0.48
1:A:19:LEU:CD2	1:A:298:ALA:HB3	2.40	0.48
1:A:320:MET:HG3	1:A:445:PHE:CZ	2.49	0.48
1:A:152:PHE:CD2	1:A:210:THR:HG23	2.49	0.47
1:A:302:ILE:C	1:A:304:GLY:N	2.67	0.47
1:A:98:GLY:C	1:A:232:ARG:HB3	2.34	0.47
1:A:16:PRO:CD	1:A:17:LYS:N	2.75	0.47
1:A:362:MET:C	1:A:363:HIS:O	2.53	0.47
1:A:115:ALA:HA	1:A:144:LEU:HD22	1.95	0.47
1:A:227:HIS:CB	1:A:231:LEU:HB2	2.44	0.47
1:A:187:PHE:HA	1:A:191:LEU:CD1	2.35	0.47
1:A:148:ILE:HG13	1:A:149:PHE:CD1	2.50	0.47
1:A:401:ILE:O	1:A:402:PHE:CD2	2.68	0.47
1:A:262:ILE:HB	1:A:278:PHE:CZ	2.50	0.47
1:A:221:LYS:HB3	1:A:226:LEU:HD12	1.97	0.47
1:A:31:PHE:CA	1:A:290:MET:SD	3.03	0.47
1:A:322:GLY:O	1:A:325:VAL:HG12	2.14	0.47
1:A:365:LEU:HA	1:A:424:PRO:HG3	1.97	0.47
1:A:151:PHE:O	1:A:152:PHE:C	2.53	0.46
1:A:34:ALA:O	1:A:38:VAL:HG23	2.15	0.46
1:A:300:GLN:CB	1:A:301:PRO:HD3	2.31	0.46
1:A:371:LEU:HD13	1:A:428:VAL:HG23	1.97	0.46
1:A:183:LEU:O	1:A:187:PHE:CD1	2.69	0.46
1:A:160:ILE:HD13	1:A:168:PHE:CD1	2.50	0.46
1:A:195:VAL:HG12	1:A:196:LEU:N	2.30	0.46
1:A:125:LEU:O	1:A:126:PHE:CB	2.62	0.46
1:A:184:ASP:O	1:A:186:LEU:N	2.48	0.46
1:A:224:LEU:HB3	1:A:225:SER:H	1.55	0.46
1:A:31:PHE:C	1:A:31:PHE:CD2	2.89	0.46
1:A:173:MET:O	1:A:174:ILE:C	2.52	0.46
1:A:286:MET:O	1:A:289:MET:HB2	2.16	0.46
1:A:139:TYR:O	1:A:143:ILE:HD12	2.16	0.46
1:A:338:MET:HA	1:A:341:PRO:HG3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:340:PHE:N	1:A:341:PRO:CD	2.79	0.46
1:A:15:ILE:HD13	1:A:15:ILE:C	2.31	0.46
1:A:188:ILE:CG2	1:A:189:PHE:N	2.79	0.46
1:A:179:LEU:HD22	1:A:183:LEU:HB2	1.97	0.46
1:A:160:ILE:HG22	1:A:217:PHE:CZ	2.50	0.46
1:A:122:ALA:O	1:A:125:LEU:HB2	2.16	0.46
1:A:156:ALA:CB	1:A:213:VAL:HG11	2.46	0.46
1:A:362:MET:SD	1:A:362:MET:N	2.89	0.45
1:A:157:ASN:ND2	1:A:169:ALA:HA	2.31	0.45
1:A:101:LEU:HD23	1:A:101:LEU:H	1.81	0.45
1:A:396:LEU:C	1:A:401:ILE:HG12	2.35	0.45
1:A:434:THR:CG2	1:A:435:VAL:N	2.79	0.45
1:A:36:TYR:CE1	1:A:180:ASN:OD1	2.69	0.45
1:A:148:ILE:HG12	1:A:149:PHE:CD1	2.52	0.45
1:A:277:VAL:HA	1:A:280:LEU:HD12	1.99	0.45
1:A:187:PHE:CD1	1:A:201:ALA:HB2	2.51	0.45
1:A:94:ASN:HD21	1:A:227:HIS:CB	2.29	0.45
1:A:110:ILE:O	1:A:113:ILE:HB	2.15	0.45
1:A:7:SER:O	1:A:8:GLU:C	2.56	0.45
1:A:245:GLY:O	1:A:246:LEU:C	2.55	0.45
1:A:302:ILE:CG2	1:A:318:THR:HG21	2.47	0.45
1:A:368:VAL:HG11	1:A:427:ASP:CB	2.47	0.45
1:A:252:GLN:O	1:A:256:SER:N	2.37	0.45
1:A:365:LEU:HD11	1:A:423:PHE:CD1	2.52	0.44
1:A:48:VAL:HG11	1:A:135:TYR:CE2	2.53	0.44
1:A:260:ILE:HG13	1:A:260:ILE:H	1.63	0.44
1:A:420:TRP:CZ3	1:A:421:TRP:CE3	2.97	0.44
1:A:160:ILE:HD11	1:A:169:ALA:N	2.32	0.44
1:A:227:HIS:HD2	1:A:231:LEU:HB2	1.81	0.44
1:A:339:LEU:C	1:A:341:PRO:HD3	2.38	0.44
1:A:177:ALA:O	1:A:180:ASN:N	2.51	0.44
1:A:300:GLN:O	1:A:384:SER:HB3	2.18	0.44
1:A:356:GLN:O	1:A:360:SER:N	2.50	0.44
1:A:219:THR:O	1:A:220:GLY:C	2.56	0.44
1:A:270:GLY:HA3	1:A:274:TYR:CD1	2.53	0.44
1:A:406:LEU:O	1:A:406:LEU:HD23	2.18	0.44
1:A:23:SER:C	1:A:25:PRO:HD2	2.37	0.44
1:A:87:GLU:N	1:A:308:GLY:CA	2.80	0.44
1:A:107:LEU:C	1:A:109:VAL:N	2.71	0.44
1:A:157:ASN:HD21	1:A:169:ALA:HA	1.83	0.44
1:A:285:MET:O	1:A:289:MET:N	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:15:ILE:HG21	1:A:314:ARG:HB3	1.99	0.44
1:A:75:ILE:O	1:A:76:GLY:C	2.56	0.44
1:A:237:VAL:O	1:A:240:GLU:N	2.48	0.44
1:A:292:MET:HE3	1:A:330:SER:N	2.33	0.44
1:A:377:VAL:O	1:A:381:LEU:HG	2.18	0.44
1:A:333:ILE:HG22	1:A:369:THR:HG21	2.00	0.43
1:A:104:ILE:O	1:A:107:LEU:N	2.51	0.43
1:A:283:ARG:O	1:A:287:PHE:CB	2.67	0.43
1:A:288:VAL:O	1:A:288:VAL:CG1	2.65	0.43
1:A:120:GLY:O	1:A:121:PRO:C	2.55	0.43
1:A:365:LEU:O	1:A:365:LEU:HG	2.17	0.43
1:A:357:ALA:HA	1:A:360:SER:CB	2.47	0.43
1:A:368:VAL:CG2	1:A:424:PRO:HA	2.48	0.43
1:A:180:ASN:OD1	1:A:202:THR:CG2	2.63	0.43
1:A:57:THR:H	1:A:57:THR:HG1	1.51	0.43
1:A:303:VAL:O	1:A:385:LEU:HD21	2.18	0.43
1:A:95:GLN:HA	1:A:95:GLN:HE21	1.84	0.43
1:A:227:HIS:CG	1:A:231:LEU:HB2	2.53	0.43
1:A:292:MET:CE	1:A:330:SER:CA	2.97	0.43
1:A:339:LEU:N	1:A:341:PRO:HD3	2.33	0.43
1:A:127:GLY:HA3	1:A:267:LEU:HG	2.00	0.43
1:A:254:SER:HB2	1:A:399:GLN:HG2	2.01	0.43
1:A:169:ALA:C	1:A:171:VAL:N	2.72	0.43
1:A:32:VAL:O	1:A:35:LEU:N	2.52	0.43
1:A:400:ILE:O	1:A:403:LEU:N	2.51	0.43
1:A:230:ASP:O	1:A:231:LEU:C	2.57	0.43
1:A:374:ALA:C	1:A:376:ILE:H	2.23	0.43
1:A:279:GLY:O	1:A:283:ARG:CD	2.65	0.43
1:A:25:PRO:HB2	1:A:166:ALA:HB1	2.00	0.42
1:A:428:VAL:O	1:A:432:ILE:HD12	2.19	0.42
1:A:287:PHE:C	1:A:287:PHE:CD2	2.92	0.42
1:A:173:MET:C	1:A:175:VAL:N	2.60	0.42
1:A:183:LEU:HA	1:A:183:LEU:HD22	1.50	0.42
1:A:12:THR:O	1:A:305:TYR:OH	2.29	0.42
1:A:446:PHE:HD2	1:A:446:PHE:H	1.67	0.42
1:A:78:ALA:CA	1:A:162:SER:HB3	2.44	0.42
1:A:301:PRO:HD2	1:A:302:ILE:H	1.84	0.42
1:A:76:GLY:O	1:A:77:GLY:C	2.57	0.42
1:A:176:PRO:HD3	1:A:209:VAL:CG1	2.50	0.42
1:A:12:THR:HG22	1:A:14:ALA:H	1.84	0.42
1:A:252:GLN:O	1:A:254:SER:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:ASN:O	1:A:181:ILE:C	2.58	0.42
1:A:246:LEU:N	1:A:247:PRO:HD2	2.35	0.42
1:A:292:MET:CE	1:A:330:SER:N	2.83	0.42
1:A:340:PHE:O	1:A:344:LEU:HD12	2.20	0.42
1:A:174:ILE:O	1:A:177:ALA:HB3	2.20	0.42
1:A:184:ASP:C	1:A:186:LEU:N	2.73	0.42
1:A:363:HIS:HB3	1:A:364:ILE:H	1.75	0.42
1:A:178:VAL:HG23	1:A:179:LEU:H	1.85	0.42
1:A:237:VAL:HA	1:A:240:GLU:CG	2.49	0.42
1:A:119:LEU:O	1:A:123:LEU:HB2	2.20	0.42
1:A:180:ASN:HD21	1:A:202:THR:CA	2.31	0.42
1:A:19:LEU:HD21	1:A:299:MET:HG2	2.01	0.42
1:A:27:MET:SD	1:A:294:GLY:HA3	2.60	0.41
1:A:157:ASN:HD21	1:A:169:ALA:CB	2.33	0.41
1:A:173:MET:CE	1:A:173:MET:HA	2.50	0.41
1:A:13:GLU:CD	1:A:306:ASN:HD21	2.18	0.41
1:A:244:VAL:HA	1:A:388:PRO:CG	2.51	0.41
1:A:374:ALA:C	1:A:376:ILE:N	2.74	0.41
1:A:180:ASN:HA	1:A:180:ASN:HD22	1.58	0.41
1:A:321:LEU:O	1:A:322:GLY:C	2.58	0.41
1:A:331:ILE:HG22	1:A:370:PHE:HB2	2.02	0.41
1:A:183:LEU:O	1:A:186:LEU:CB	2.68	0.41
1:A:187:PHE:CA	1:A:191:LEU:HD12	2.34	0.41
1:A:12:THR:HG22	1:A:13:GLU:N	2.36	0.41
1:A:354:VAL:O	1:A:358:GLY:HA3	2.21	0.41
1:A:292:MET:O	1:A:293:MET:C	2.59	0.41
1:A:20:ARG:HD3	1:A:20:ARG:HA	1.72	0.41
1:A:404:ILE:C	1:A:406:LEU:N	2.72	0.41
1:A:247:PRO:HG2	1:A:383:GLN:HE22	1.83	0.41
1:A:301:PRO:CD	1:A:302:ILE:H	2.34	0.41
1:A:312:TYR:O	1:A:316:ARG:HG3	2.21	0.41
1:A:320:MET:HG3	1:A:445:PHE:HE1	1.81	0.41
1:A:323:PHE:HE1	1:A:377:VAL:CG1	2.29	0.41
1:A:379:GLY:O	1:A:383:GLN:HG3	2.21	0.41
1:A:148:ILE:HG12	1:A:149:PHE:CE1	2.55	0.41
1:A:95:GLN:HE22	1:A:232:ARG:HA	1.86	0.41
1:A:390:GLN:HG2	1:A:390:GLN:H	1.61	0.41
1:A:183:LEU:HD12	1:A:201:ALA:HA	2.03	0.41
1:A:18:LEU:HD23	1:A:302:ILE:HG12	2.03	0.41
1:A:232:ARG:H	1:A:232:ARG:HG3	1.57	0.41
1:A:160:ILE:HA	1:A:163:GLU:HG2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:ILE:N	1:A:181:ILE:HD13	2.30	0.41
1:A:301:PRO:C	1:A:304:GLY:H	2.24	0.41
1:A:325:VAL:O	1:A:328:ILE:N	2.54	0.41
1:A:83:ARG:O	1:A:87:GLU:HB3	2.21	0.41
1:A:428:VAL:CG1	1:A:429:LEU:N	2.82	0.41
1:A:331:ILE:CD1	1:A:331:ILE:C	2.83	0.41
1:A:160:ILE:HD12	1:A:168:PHE:CD1	2.55	0.40
1:A:35:LEU:O	1:A:39:VAL:HG22	2.21	0.40
1:A:78:ALA:O	1:A:82:SER:OG	2.36	0.40
1:A:413:ILE:O	1:A:414:PHE:CD1	2.74	0.40
1:A:214:LEU:N	1:A:214:LEU:HD23	2.35	0.40
1:A:36:TYR:CD2	1:A:37:ASN:N	2.89	0.40
1:A:96:VAL:O	1:A:99:ASN:N	2.54	0.40
1:A:320:MET:CB	1:A:445:PHE:HE1	2.30	0.40
1:A:6:GLN:HG3	1:A:6:GLN:O	2.22	0.40
1:A:100:ILE:O	1:A:104:ILE:HD12	2.21	0.40
1:A:96:VAL:HG22	1:A:237:VAL:HG11	2.04	0.40
1:A:341:PRO:HB2	1:A:359:VAL:CG2	2.52	0.40
1:A:266:LEU:HD22	1:A:278:PHE:CE1	2.56	0.40
1:A:302:ILE:HG22	1:A:318:THR:HG21	2.03	0.40
1:A:252:GLN:C	1:A:254:SER:H	2.24	0.40
1:A:434:THR:O	1:A:435:VAL:C	2.60	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	444/463 (96%)	252 (57%)	129 (29%)	63 (14%)	<b>0</b> <b>1</b>

All (63) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	8	GLU
1	A	9	ARG
1	A	15	ILE
1	A	16	PRO
1	A	38	VAL
1	A	89	ARG
1	A	170	MET
1	A	174	ILE
1	A	177	ALA
1	A	181	ILE
1	A	185	VAL
1	A	196	LEU
1	A	216	TYR
1	A	221	LYS
1	A	223	THR
1	A	290	MET
1	A	300	GLN
1	A	309	ALA
1	A	311	GLN
1	A	323	PHE
1	A	324	LYS
1	A	363	HIS
1	A	369	THR
1	A	375	GLN
1	A	386	GLY
1	A	401	ILE
1	A	77	GLY
1	A	85	LEU
1	A	90	GLY
1	A	96	VAL
1	A	97	PHE
1	A	151	PHE
1	A	163	GLU
1	A	194	GLY
1	A	222	SER
1	A	234	LYS
1	A	253	SER
1	A	322	GLY
1	A	326	ALA
1	A	364	ILE
1	A	368	VAL
1	A	373	GLY
1	A	126	PHE

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Mol	Chain	Res	Type
1	A	215	ARG
1	A	293	MET
1	A	301	PRO
1	A	325	VAL
1	A	367	CYS
1	A	384	SER
1	A	412	HIS
1	A	60	PHE
1	A	76	GLY
1	A	229	SER
1	A	355	ILE
1	A	285	MET
1	A	327	THR
1	A	423	PHE
1	A	56	VAL
1	A	188	ILE
1	A	303	VAL
1	A	80	VAL
1	A	410	LEU
1	A	104	ILE

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	366/382 (96%)	271 (74%)	95 (26%)	<b>0</b> <b>3</b>

All (95) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ARG
1	A	15	ILE
1	A	21	SER
1	A	22	LEU
1	A	24	ILE
1	A	27	MET

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Mol	Chain	Res	Type
1	A	31	PHE
1	A	33	MET
1	A	48	VAL
1	A	51	GLU
1	A	57	THR
1	A	63	MET
1	A	65	ILE
1	A	66	MET
1	A	75	ILE
1	A	79	SER
1	A	97	PHE
1	A	99	ASN
1	A	101	LEU
1	A	105	LEU
1	A	119	LEU
1	A	126	PHE
1	A	129	THR
1	A	141	PHE
1	A	147	SER
1	A	148	ILE
1	A	150	PHE
1	A	159	ILE
1	A	160	ILE
1	A	167	THR
1	A	168	PHE
1	A	174	ILE
1	A	175	VAL
1	A	178	VAL
1	A	179	LEU
1	A	180	ASN
1	A	181	ILE
1	A	182	LEU
1	A	183	LEU
1	A	186	LEU
1	A	188	ILE
1	A	196	LEU
1	A	200	ILE
1	A	203	VAL
1	A	204	ILE
1	A	209	VAL
1	A	213	VAL
1	A	222	SER

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Mol	Chain	Res	Type
1	A	223	THR
1	A	224	LEU
1	A	226	LEU
1	A	228	TRP
1	A	230	ASP
1	A	232	ARG
1	A	234	LYS
1	A	236	SER
1	A	259	MET
1	A	260	ILE
1	A	267	LEU
1	A	268	ARG
1	A	271	SER
1	A	278	PHE
1	A	283	ARG
1	A	284	ILE
1	A	285	MET
1	A	288	VAL
1	A	289	MET
1	A	292	MET
1	A	299	MET
1	A	315	LEU
1	A	317	GLU
1	A	319	VAL
1	A	325	VAL
1	A	330	SER
1	A	331	ILE
1	A	336	LEU
1	A	339	LEU
1	A	345	LEU
1	A	348	PHE
1	A	359	VAL
1	A	363	HIS
1	A	370	PHE
1	A	375	GLN
1	A	382	TYR
1	A	390	GLN
1	A	397	SER
1	A	410	LEU
1	A	419	VAL
1	A	428	VAL
1	A	429	LEU

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Mol	Chain	Res	Type
1	A	431	PHE
1	A	432	ILE
1	A	434	THR
1	A	436	VAL
1	A	446	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	95	GLN
1	A	157	ASN
1	A	180	ASN
1	A	227	HIS

### 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<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	446/463 (96%)	0.47	63 (14%) 4 1	103, 190, 291, 366	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	447	LEU	11.7
1	A	3	GLN	11.0
1	A	271	SER	10.3
1	A	129	THR	8.3
1	A	270	GLY	7.8
1	A	269	PHE	6.7
1	A	6	GLN	6.6
1	A	127	GLY	6.3
1	A	225	SER	6.2
1	A	233	MET	6.0
1	A	133	GLN	5.8
1	A	352	ARG	5.8
1	A	132	THR	5.7
1	A	8	GLU	5.4
1	A	191	LEU	5.3
1	A	131	VAL	4.8
1	A	7	SER	4.7
1	A	51	GLU	4.6
1	A	155	ALA	4.5
1	A	308	GLY	4.5
1	A	293	MET	4.5
1	A	52	GLY	4.2
1	A	356	GLN	4.2
1	A	416	LEU	3.9
1	A	384	SER	3.9
1	A	128	ALA	3.9
1	A	116	PHE	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	87	GLU	3.7
1	A	234	LYS	3.5
1	A	230	ASP	3.5
1	A	89	ARG	3.5
1	A	13	GLU	3.5
1	A	272	ASP	3.4
1	A	9	ARG	3.4
1	A	353	GLU	3.3
1	A	312	TYR	3.3
1	A	130	SER	3.3
1	A	4	LYS	3.2
1	A	162	SER	3.1
1	A	306	ASN	3.0
1	A	324	LYS	3.0
1	A	64	MET	2.9
1	A	156	ALA	2.8
1	A	126	PHE	2.8
1	A	235	GLY	2.8
1	A	274	TYR	2.7
1	A	238	ILE	2.7
1	A	364	ILE	2.5
1	A	445	PHE	2.5
1	A	219	THR	2.4
1	A	439	TYR	2.3
1	A	200	ILE	2.3
1	A	91	GLU	2.2
1	A	50	ILE	2.2
1	A	135	TYR	2.2
1	A	190	GLY	2.2
1	A	354	VAL	2.1
1	A	314	ARG	2.1
1	A	386	GLY	2.1
1	A	42	ILE	2.0
1	A	350	ALA	2.0
1	A	90	GLY	2.0
1	A	98	GLY	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 [i](#)

There are no carbohydrates 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.