



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:50 PM UTC

PDB ID : 2WL6 / pdb_00002wl6
Title : BIOSYNTHETIC THIOLASE FROM Z. RAMIGERA. THE N316H-H348N MUTANT.
Authors : Merilainen, G.; Poikela, V.; Kursula, P.; Wierenga, R.K.
Deposited on : 2009-06-22
Resolution : 2.98 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

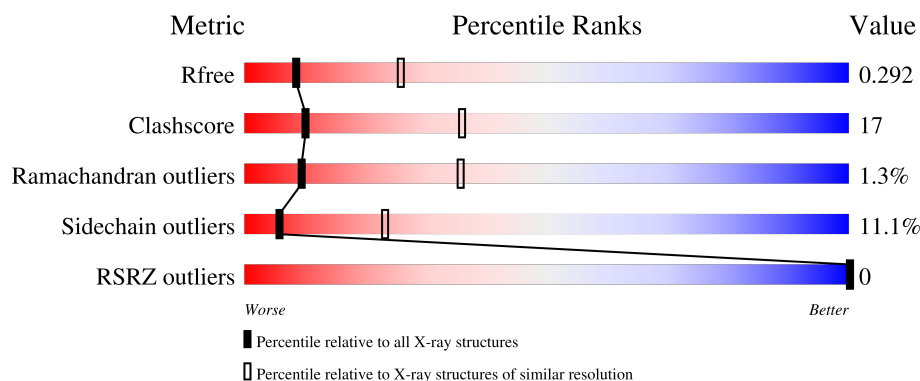
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.98 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}	180053	3580 (3.00-2.96)
Clashscore	190562	3904 (3.00-2.96)
Ramachandran outliers	187476	3761 (3.00-2.96)
Sidechain outliers	187428	3764 (3.00-2.96)
RSRZ outliers	180081	3579 (3.00-2.96)

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 $\leq 5\%$. 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	392	
1	B	392	
1	C	392	
1	D	392	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11278 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 ACETYL-COA ACETYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	389	Total	C	N	O	S	0	0	0
			2813	1746	509	537	21			
1	B	389	Total	C	N	O	S	0	0	0
			2813	1746	509	537	21			
1	C	389	Total	C	N	O	S	0	0	0
			2813	1746	509	537	21			
1	D	389	Total	C	N	O	S	0	0	0
			2813	1746	509	537	21			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	129	ARG	ALA	SEE REMARK 999	UNP P07097
A	316	HIS	ASN	engineered mutation	UNP P07097
A	348	ASN	HIS	engineered mutation	UNP P07097
B	129	ARG	ALA	SEE REMARK 999	UNP P07097
B	316	HIS	ASN	engineered mutation	UNP P07097
B	348	ASN	HIS	engineered mutation	UNP P07097
C	129	ARG	ALA	SEE REMARK 999	UNP P07097
C	316	HIS	ASN	engineered mutation	UNP P07097
C	348	ASN	HIS	engineered mutation	UNP P07097
D	129	ARG	ALA	SEE REMARK 999	UNP P07097
D	316	HIS	ASN	engineered mutation	UNP P07097
D	348	ASN	HIS	engineered mutation	UNP P07097

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	10	Total	O	0	0
			10	10		
2	B	8	Total	O	0	0
			8	8		

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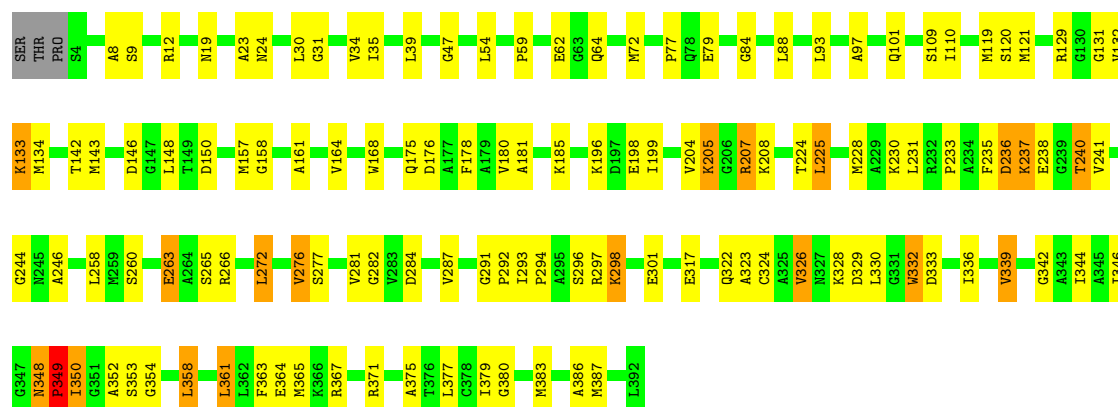
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	3	Total	O	0	0
			3	3		
2	D	5	Total	O	0	0
			5	5		

3 Residue-property plots

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.

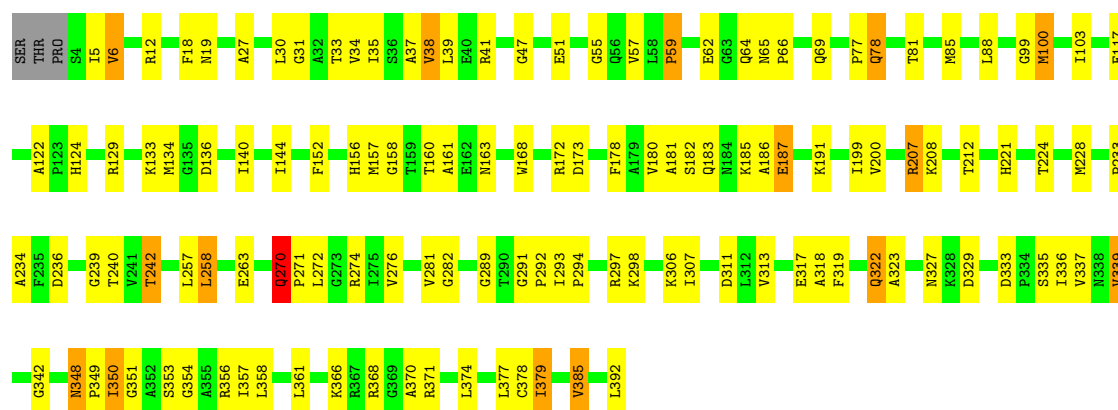
• Molecule 1: ACETYL-COA ACETYLTRANSFERASE

Chain A: 



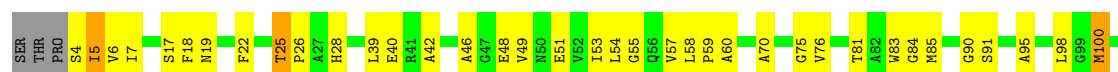
• Molecule 1: ACETYL-COA ACETYLTRANSFERASE

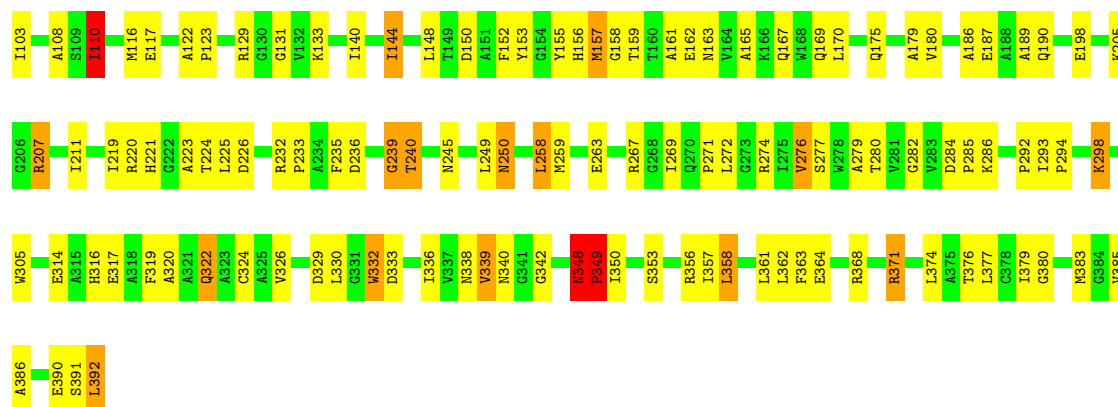
Chain B: 



• Molecule 1: ACETYL-COA ACETYLTRANSFERASE

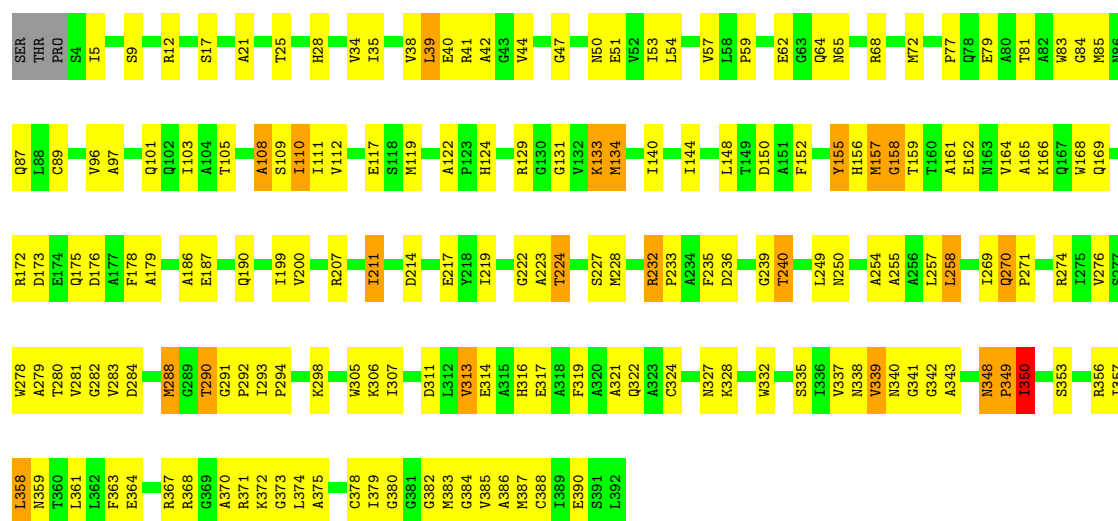
Chain C: 





• Molecule 1: ACETYL-COA ACETYLTRANSFERASE

Chain D: 55% 39% 5% •



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	84.23Å 79.87Å 149.92Å 90.00° 92.86° 90.00°	Depositor
Resolution (Å)	37.42 – 2.98 37.42 – 2.98	Depositor EDS
% Data completeness (in resolution range)	95.9 (37.42-2.98) 82.8 (37.42-2.98)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	REFMAC 5.4.0077	Depositor
R, R_{free}	0.232 , 0.291 0.233 , 0.292	Depositor DCC
R_{free} test set	1974 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	0.217	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 27.1	EDS
L-test for twinning ¹	$\langle L \rangle = 0.34$, $\langle L^2 \rangle = 0.17$	Xtriage
Estimated twinning fraction	0.176 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11278	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.83	0/2854	1.16	8/3853 (0.2%)
1	B	0.81	0/2854	1.06	10/3853 (0.3%)
1	C	0.70	0/2854	1.21	11/3853 (0.3%)
1	D	0.75	1/2854 (0.0%)	1.25	13/3853 (0.3%)
All	All	0.78	1/11416 (0.0%)	1.17	42/15412 (0.3%)

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	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	348	ASN	CA-C	-7.68	1.46	1.52

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	348	ASN	CA-C-N	-32.03	79.80	119.84
1	D	348	ASN	C-N-CA	-32.03	79.80	119.84
1	C	348	ASN	CA-C-N	-27.41	85.57	119.84
1	C	348	ASN	C-N-CA	-27.41	85.57	119.84
1	A	348	ASN	CA-C-N	-23.62	90.31	119.84
1	A	348	ASN	C-N-CA	-23.62	90.31	119.84
1	C	348	ASN	C-N-CD	-14.04	67.42	125.00
1	B	199	ILE	N-CA-C	8.18	119.53	109.30
1	D	348	ASN	C-N-CD	-7.19	95.53	125.00
1	C	5	ILE	N-CA-C	6.84	117.98	107.99
1	A	348	ASN	C-N-CD	-6.46	98.50	125.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	62	GLU	N-CA-C	6.29	119.11	111.82
1	C	284	ASP	CA-C-N	6.28	126.75	119.47
1	C	284	ASP	C-N-CA	6.28	126.75	119.47
1	A	348	ASN	N-CA-C	6.18	118.25	108.23
1	D	341	GLY	N-CA-C	6.14	117.78	111.95
1	C	349	PRO	N-CA-C	5.87	124.56	112.47
1	A	110	ILE	N-CA-C	5.81	116.37	107.77
1	B	327	ASN	N-CA-C	5.75	117.42	111.03
1	C	25	THR	CA-C-N	5.68	125.44	119.76
1	C	25	THR	C-N-CA	5.68	125.44	119.76
1	C	110	ILE	N-CA-C	5.66	116.43	108.17
1	D	108	ALA	N-CA-C	5.63	117.68	108.55
1	B	27	ALA	N-CA-C	5.50	116.95	111.07
1	D	313	VAL	N-CA-C	5.49	116.27	108.42
1	B	6	VAL	CB-CA-C	-5.43	101.13	110.30
1	B	379	ILE	N-CA-C	5.41	116.08	108.17
1	A	349	PRO	N-CA-C	-5.32	101.51	112.47
1	D	316	HIS	CA-C-N	-5.32	114.34	122.21
1	D	316	HIS	C-N-CA	-5.32	114.34	122.21
1	D	211	ILE	N-CA-C	5.22	115.64	108.12
1	D	87	GLN	N-CA-C	-5.20	104.28	110.41
1	B	99	GLY	N-CA-C	-5.19	106.48	112.50
1	D	155	TYR	N-CA-C	5.19	118.53	110.17
1	D	44	VAL	N-CA-C	5.17	116.25	109.21
1	C	42	ALA	N-CA-C	-5.05	107.11	113.28
1	B	348	ASN	N-CA-C	5.05	116.03	108.76
1	B	270	GLN	CA-C-N	5.04	125.02	120.03
1	B	270	GLN	C-N-CA	5.04	125.02	120.03
1	A	284	ASP	CA-C-N	5.03	126.13	119.84
1	A	284	ASP	C-N-CA	5.03	126.13	119.84
1	D	21	ALA	N-CA-C	5.02	117.41	111.33

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	349	PRO	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	2813	0	2819	91	0
1	B	2813	0	2819	81	0
1	C	2813	0	2819	104	0
1	D	2813	0	2819	128	0
2	A	10	0	0	0	0
2	B	8	0	0	1	0
2	C	3	0	0	0	0
2	D	5	0	0	1	0
All	All	11278	0	11276	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 17.

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)
1:C:348:ASN:O	1:C:349:PRO:C	1.65	1.21
1:C:207:ARG:HG2	1:C:207:ARG:HH11	0.88	1.04
1:D:17:SER:OG	1:D:217:GLU:HG3	1.63	0.99
1:C:207:ARG:HG2	1:C:207:ARG:NH1	1.68	0.95
1:D:9:SER:HB3	1:D:42:ALA:HB2	1.48	0.94
1:C:348:ASN:OD1	1:C:349:PRO:N	2.03	0.93
1:C:250:ASN:HB2	1:C:349:PRO:HD3	1.54	0.89
1:C:207:ARG:HH11	1:C:207:ARG:CG	1.82	0.89
1:B:180:VAL:HG22	1:B:228:MET:HE3	1.53	0.88
1:D:236:ASP:HB3	1:D:239:GLY:HA3	1.58	0.84
1:D:41:ARG:HH21	1:D:200:VAL:HB	1.43	0.84
1:A:358:LEU:HD11	1:A:387:MET:HE1	1.58	0.84
1:B:157:MET:HE2	1:B:157:MET:HA	1.62	0.82
1:B:348:ASN:OD1	1:B:353:SER:OG	1.98	0.82
1:D:175:GLN:HE22	1:D:240:THR:CG2	1.94	0.81
1:D:47:GLY:HA2	1:D:77:PRO:HG3	1.63	0.80
1:C:379:ILE:HG21	1:C:383:MET:HE3	1.63	0.78
1:A:132:VAL:O	1:C:129:ARG:HA	1.83	0.78
1:C:348:ASN:O	1:C:349:PRO:O	2.02	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:288:MET:SD	1:D:380:GLY:HA2	2.25	0.75
1:B:354:GLY:HA2	1:B:377:LEU:HD11	1.69	0.75
1:B:207:ARG:HD3	1:B:207:ARG:H	1.52	0.74
1:B:339:VAL:HG11	1:B:368:ARG:NH2	2.03	0.74
1:B:293:ILE:HB	1:B:294:PRO:CD	2.17	0.74
1:B:293:ILE:HB	1:B:294:PRO:HD3	1.68	0.74
1:A:348:ASN:C	1:A:348:ASN:OD1	2.30	0.73
1:A:64:GLN:HG2	1:B:88:LEU:HD21	1.69	0.73
1:B:281:VAL:HG12	1:B:282:GLY:N	2.06	0.71
1:C:236:ASP:HB3	1:C:239:GLY:HA3	1.71	0.71
1:D:348:ASN:OD1	1:D:348:ASN:C	2.34	0.70
1:D:236:ASP:HB3	1:D:239:GLY:CA	2.20	0.70
1:A:131:GLY:HA2	1:C:131:GLY:HA3	1.75	0.69
1:A:348:ASN:OD1	1:A:348:ASN:O	2.11	0.69
1:B:168:TRP:CH2	1:B:329:ASP:HB2	2.28	0.68
1:C:317:GLU:CD	1:C:342:GLY:HA3	2.18	0.68
1:C:179:ALA:HB2	1:C:245:ASN:HB3	1.76	0.68
1:C:333:ASP:O	1:C:336:ILE:HG12	1.94	0.68
1:B:51:GLU:OE2	1:B:81:THR:OG1	2.13	0.67
1:D:17:SER:HG	1:D:217:GLU:HG3	1.59	0.67
1:C:348:ASN:OD1	1:C:348:ASN:C	2.38	0.66
1:C:95:ALA:HA	1:C:98:LEU:HD12	1.78	0.66
1:D:175:GLN:HE22	1:D:240:THR:HG21	1.58	0.66
1:D:175:GLN:HE22	1:D:240:THR:HG23	1.60	0.65
1:D:5:ILE:HD12	1:D:103:ILE:CG2	2.27	0.65
1:C:54:LEU:O	1:C:84:GLY:HA2	1.96	0.65
1:A:93:LEU:HD11	1:A:387:MET:HE2	1.79	0.65
1:C:157:MET:HE2	1:C:157:MET:HA	1.78	0.65
1:D:97:ALA:O	1:D:101:GLN:HG3	1.97	0.65
1:C:198:GLU:HB3	1:C:363:PHE:CD2	2.32	0.64
1:C:162:GLU:OE2	1:C:240:THR:N	2.29	0.64
1:A:133:LYS:O	1:A:134:MET:HB3	1.96	0.64
1:D:101:GLN:O	1:D:105:THR:HG23	1.97	0.63
1:A:180:VAL:HG21	1:A:225:LEU:HA	1.80	0.63
1:A:292:PRO:O	1:A:296:SER:OG	2.12	0.63
1:D:150:ASP:OD2	1:D:152:PHE:HB2	1.99	0.62
1:A:348:ASN:HB2	1:A:353:SER:OG	1.99	0.62
1:C:358:LEU:HD22	1:C:362:LEU:HG	1.82	0.62
1:C:148:LEU:O	1:C:157:MET:HG2	1.99	0.62
1:D:5:ILE:HD12	1:D:103:ILE:HG21	1.82	0.62
1:B:356:ARG:C	1:B:356:ARG:HD2	2.25	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:190:GLN:HE21	1:D:219:ILE:HD12	1.64	0.61
1:A:148:LEU:O	1:A:157:MET:HG2	2.01	0.61
1:D:41:ARG:NH2	1:D:200:VAL:HB	2.15	0.61
1:A:88:LEU:HD12	1:A:380:GLY:O	2.01	0.61
1:D:119:MET:HE3	1:D:350:ILE:CG1	2.30	0.61
1:A:129:ARG:HH21	1:B:122:ALA:HB3	1.66	0.60
1:B:168:TRP:HH2	1:B:329:ASP:HB2	1.64	0.60
1:C:280:THR:HG23	1:D:81:THR:HG21	1.83	0.60
1:C:364:GLU:O	1:C:368:ARG:HG2	2.02	0.60
1:A:317:GLU:CD	1:A:342:GLY:HA3	2.27	0.59
1:B:293:ILE:O	1:B:297:ARG:HG3	2.03	0.59
1:C:282:GLY:O	1:D:79:GLU:HA	2.02	0.59
1:D:96:VAL:HG12	1:D:387:MET:HE1	1.84	0.58
1:B:356:ARG:HD2	1:B:356:ARG:O	2.03	0.58
1:A:168:TRP:CH2	1:A:329:ASP:HB2	2.38	0.58
1:B:258:LEU:N	1:B:258:LEU:HD22	2.18	0.58
1:A:142:THR:O	1:A:146:ASP:HB2	2.04	0.58
1:C:162:GLU:OE1	1:C:240:THR:HG22	2.04	0.57
1:B:371:ARG:NH2	2:B:2008:HOH:O	2.37	0.57
1:A:281:VAL:HG12	1:A:282:GLY:N	2.19	0.57
1:A:236:ASP:C	1:A:236:ASP:OD1	2.47	0.57
1:B:313:VAL:HB	1:B:337:VAL:HG22	1.87	0.56
1:D:53:ILE:HD13	1:D:83:TRP:CZ2	2.40	0.56
1:D:179:ALA:HB3	1:D:228:MET:SD	2.45	0.56
1:D:119:MET:HE3	1:D:350:ILE:HD11	1.88	0.56
1:A:228:MET:HE1	1:A:244:GLY:C	2.31	0.56
1:A:258:LEU:N	1:A:258:LEU:HD22	2.22	0.55
1:B:12:ARG:O	1:B:200:VAL:HG12	2.07	0.55
1:C:53:ILE:HD13	1:C:83:TRP:CZ2	2.42	0.55
1:D:165:ALA:O	1:D:169:GLN:N	2.39	0.55
1:D:269:ILE:O	1:D:271:PRO:HD3	2.06	0.55
1:D:317:GLU:CD	1:D:342:GLY:HA3	2.32	0.55
1:A:361:LEU:O	1:A:365:MET:HG3	2.07	0.55
1:C:57:VAL:O	1:C:59:PRO:HD3	2.08	0.54
1:B:158:GLY:O	1:B:161:ALA:HB3	2.08	0.54
1:C:339:VAL:HG11	1:C:368:ARG:NH2	2.22	0.54
1:D:313:VAL:CG1	1:D:314:GLU:N	2.69	0.54
1:A:263:GLU:HA	1:A:266:ARG:NH1	2.23	0.54
1:C:348:ASN:OD1	1:C:349:PRO:CA	2.56	0.54
1:D:155:TYR:CD1	1:D:159:THR:HB	2.43	0.54
1:D:57:VAL:O	1:D:59:PRO:HD3	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:34:VAL:CG1	1:D:255:ALA:HB3	2.38	0.54
1:D:367:ARG:O	1:D:367:ARG:HG2	2.08	0.54
1:A:157:MET:HE2	1:A:157:MET:HA	1.89	0.54
1:D:258:LEU:HD22	1:D:258:LEU:N	2.22	0.54
1:A:30:LEU:O	1:A:34:VAL:HG23	2.08	0.53
1:B:134:MET:HE3	1:C:144:ILE:HD11	1.90	0.53
1:D:313:VAL:HG12	1:D:337:VAL:HG13	1.89	0.53
1:B:281:VAL:CG1	1:B:282:GLY:N	2.71	0.53
1:B:374:LEU:HD23	1:B:374:LEU:C	2.33	0.53
1:C:54:LEU:HD13	1:C:116:MET:CE	2.37	0.53
1:A:291:GLY:O	1:A:294:PRO:HD2	2.07	0.53
1:D:109:SER:O	1:D:110:ILE:HG12	2.09	0.53
1:B:233:PRO:HA	1:B:242:THR:HG22	1.91	0.53
1:D:379:ILE:HD11	1:D:385:VAL:HG12	1.90	0.52
1:C:207:ARG:NH1	1:C:207:ARG:CG	2.53	0.52
1:D:158:GLY:HA3	1:D:235:PHE:CG	2.44	0.52
1:B:281:VAL:HG12	1:B:282:GLY:H	1.71	0.52
1:D:156:HIS:ND1	1:D:157:MET:N	2.57	0.52
1:A:164:VAL:O	1:A:168:TRP:HB2	2.09	0.52
1:A:77:PRO:HB2	1:A:79:GLU:OE1	2.09	0.51
1:C:322:GLN:O	1:C:326:VAL:HG23	2.09	0.51
1:D:348:ASN:OD1	1:D:349:PRO:N	2.43	0.51
1:A:175:GLN:HE22	1:A:240:THR:CG2	2.23	0.51
1:D:162:GLU:HB3	1:D:166:LYS:HE3	1.92	0.51
1:A:47:GLY:HA2	1:A:77:PRO:HD3	1.92	0.51
1:B:236:ASP:O	1:B:239:GLY:N	2.44	0.51
1:A:158:GLY:O	1:A:161:ALA:HB3	2.11	0.51
1:B:172:ARG:NH2	1:B:242:THR:HG21	2.25	0.51
1:C:122:ALA:HB3	1:D:129:ARG:HH21	1.75	0.51
1:D:348:ASN:OD1	1:D:349:PRO:CA	2.59	0.51
1:D:290:THR:O	1:D:294:PRO:HD2	2.11	0.51
1:D:292:PRO:HD3	1:D:378:CYS:HA	1.91	0.51
1:A:12:ARG:O	1:A:199:ILE:HA	2.10	0.51
1:B:117:GLU:OE1	1:B:351:GLY:N	2.43	0.51
1:D:307:ILE:HD12	1:D:307:ILE:H	1.76	0.51
1:B:183:GLN:OE1	1:B:183:GLN:HA	2.10	0.51
1:D:51:GLU:HB3	1:D:111:ILE:HD12	1.93	0.51
1:D:65:ASN:O	1:D:68:ARG:HB3	2.10	0.51
1:A:9:SER:HA	1:A:272:LEU:HD22	1.93	0.50
1:D:54:LEU:O	1:D:84:GLY:HA2	2.11	0.50
1:C:269:ILE:O	1:C:271:PRO:HD3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:316:HIS:CG	1:C:377:LEU:HD23	2.46	0.50
1:C:385:VAL:HG22	1:C:386:ALA:N	2.25	0.50
1:A:276:VAL:O	1:A:277:SER:HB3	2.11	0.50
1:B:64:GLN:O	1:B:65:ASN:C	2.50	0.50
1:B:333:ASP:C	1:B:335:SER:H	2.20	0.50
1:D:12:ARG:O	1:D:199:ILE:HA	2.12	0.50
1:D:150:ASP:HB2	1:D:157:MET:HE3	1.92	0.50
1:D:348:ASN:OD1	1:D:349:PRO:C	2.54	0.50
1:A:298:LYS:NZ	1:A:301:GLU:OE1	2.44	0.50
1:B:37:ALA:O	1:B:41:ARG:HG3	2.10	0.50
1:D:148:LEU:O	1:D:157:MET:HG2	2.11	0.50
1:A:8:ALA:O	1:A:9:SER:HB3	2.11	0.50
1:D:12:ARG:HB2	1:D:254:ALA:HB2	1.94	0.50
1:D:161:ALA:HB2	1:D:319:PHE:CD1	2.47	0.50
1:B:333:ASP:O	1:B:336:ILE:HG12	2.12	0.49
1:B:317:GLU:CD	1:B:342:GLY:HA3	2.37	0.49
1:D:378:CYS:C	1:D:379:ILE:HG13	2.36	0.49
1:D:357:ILE:CD1	1:D:375:ALA:HB1	2.43	0.49
1:C:190:GLN:NE2	1:C:219:ILE:HD12	2.28	0.49
1:C:314:GLU:OE2	1:C:338:ASN:HA	2.13	0.49
1:C:348:ASN:CG	1:C:349:PRO:N	2.69	0.49
1:D:364:GLU:O	1:D:368:ARG:HG2	2.13	0.49
1:D:293:ILE:HB	1:D:294:PRO:HD3	1.95	0.49
1:A:150:ASP:HB2	1:A:157:MET:HE1	1.94	0.49
1:C:5:ILE:HG13	1:C:100:MET:HG2	1.95	0.49
1:D:9:SER:CB	1:D:42:ALA:HB2	2.32	0.48
1:D:158:GLY:HA2	1:D:319:PHE:CE2	2.47	0.48
1:C:293:ILE:HB	1:C:294:PRO:CD	2.44	0.48
1:A:326:VAL:HG13	1:A:330:LEU:HD12	1.94	0.48
1:B:31:GLY:O	1:B:35:ILE:HG13	2.13	0.48
1:B:47:GLY:HA2	1:B:77:PRO:HD3	1.95	0.48
1:B:57:VAL:O	1:B:59:PRO:HD3	2.12	0.48
1:C:103:ILE:HA	1:C:108:ALA:O	2.13	0.48
1:D:314:GLU:OE2	1:D:338:ASN:HA	2.14	0.48
1:A:207:ARG:H	1:A:207:ARG:HD3	1.79	0.48
1:A:342:GLY:O	1:A:346:ILE:HD12	2.14	0.48
1:C:250:ASN:OD1	1:C:348:ASN:N	2.44	0.48
1:C:263:GLU:CD	1:C:267:ARG:HE	2.22	0.48
1:C:330:LEU:HD12	1:C:332:TRP:CZ2	2.49	0.48
1:C:153:TYR:CE2	1:C:286:LYS:HD3	2.49	0.47
1:C:175:GLN:HE22	1:C:240:THR:HG21	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:VAL:CG2	1:A:228:MET:HG3	2.44	0.47
1:A:168:TRP:HH2	1:A:329:ASP:HB2	1.76	0.47
1:A:349:PRO:O	1:A:350:ILE:C	2.58	0.47
1:B:257:LEU:HD23	1:B:257:LEU:C	2.39	0.47
1:B:293:ILE:CB	1:B:294:PRO:CD	2.86	0.47
1:A:354:GLY:HA2	1:A:377:LEU:HD21	1.96	0.47
1:B:18:PHE:O	1:B:19:ASN:C	2.54	0.47
1:C:90:GLY:O	1:C:91:SER:C	2.58	0.47
1:A:317:GLU:OE1	1:A:344:ILE:HG13	2.15	0.47
1:C:152:PHE:CE1	1:D:72:MET:HG3	2.49	0.47
1:D:12:ARG:HD2	1:D:356:ARG:HG2	1.96	0.47
1:A:97:ALA:O	1:A:101:GLN:HG3	2.15	0.47
1:A:204:VAL:O	1:A:205:LYS:C	2.58	0.47
1:A:282:GLY:HA2	1:A:383:MET:HA	1.97	0.47
1:B:156:HIS:ND1	1:B:157:MET:N	2.63	0.47
1:A:297:ARG:HG2	1:A:297:ARG:HH11	1.80	0.47
1:B:274:ARG:HE	1:B:392:LEU:HD21	1.78	0.47
1:D:178:PHE:HE1	1:D:317:GLU:CD	2.23	0.47
1:A:59:PRO:O	1:A:62:GLU:HB2	2.15	0.47
1:A:292:PRO:HB2	1:A:326:VAL:HG21	1.97	0.47
1:D:51:GLU:OE2	1:D:81:THR:OG1	2.30	0.47
1:B:12:ARG:HD2	1:B:356:ARG:HG2	1.97	0.46
1:C:329:ASP:C	1:C:329:ASP:OD1	2.58	0.46
1:D:133:LYS:HA	1:D:133:LYS:HD3	1.60	0.46
1:D:233:PRO:HB2	1:D:236:ASP:O	2.14	0.46
1:D:278:TRP:HA	1:D:386:ALA:O	2.14	0.46
1:B:234:ALA:H	1:B:242:THR:HA	1.79	0.46
1:B:291:GLY:N	1:B:292:PRO:CD	2.79	0.46
1:C:51:GLU:CD	1:C:83:TRP:HE1	2.24	0.46
1:C:156:HIS:ND1	1:C:157:MET:N	2.63	0.46
1:C:371:ARG:HA	1:C:391:SER:OG	2.15	0.46
1:D:39:LEU:O	1:D:40:GLU:C	2.57	0.46
1:D:186:ALA:HA	1:D:340:ASN:O	2.15	0.46
1:A:198:GLU:HB3	1:A:363:PHE:CD2	2.50	0.46
1:C:75:GLY:O	1:C:76:VAL:C	2.59	0.46
1:D:119:MET:HE3	1:D:350:ILE:CD1	2.45	0.46
1:A:330:LEU:HD13	1:A:332:TRP:CH2	2.50	0.46
1:A:348:ASN:HA	1:A:349:PRO:HD3	1.20	0.46
1:B:207:ARG:HG2	1:B:208:LYS:H	1.81	0.46
1:A:54:LEU:O	1:A:84:GLY:HA2	2.15	0.46
1:D:124:HIS:HA	1:D:140:ILE:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:120:SER:O	1:B:129:ARG:HD2	2.16	0.46
1:D:317:GLU:HB2	1:D:343:ALA:H	1.81	0.46
1:A:31:GLY:O	1:A:35:ILE:HG13	2.16	0.46
1:B:289:GLY:HA2	1:B:378:CYS:SG	2.55	0.46
1:C:22:PHE:HB3	1:C:25:THR:HB	1.98	0.46
1:D:28:HIS:ND1	1:D:62:GLU:OE2	2.34	0.46
1:D:57:VAL:C	1:D:59:PRO:HD3	2.41	0.46
1:D:117:GLU:HA	1:D:117:GLU:OE1	2.15	0.46
1:D:317:GLU:HB2	1:D:343:ALA:N	2.31	0.46
1:A:204:VAL:O	1:A:204:VAL:HG12	2.16	0.45
1:B:160:THR:HA	1:B:163:ASN:HD22	1.81	0.45
1:B:281:VAL:CG1	1:B:282:GLY:H	2.29	0.45
1:C:155:TYR:CD2	1:C:159:THR:HG21	2.51	0.45
1:D:199:ILE:HG22	1:D:200:VAL:N	2.31	0.45
1:A:333:ASP:O	1:A:336:ILE:HG12	2.16	0.45
1:B:318:ALA:N	1:B:322:GLN:OE1	2.47	0.45
1:C:356:ARG:NH2	1:C:357:ILE:HG22	2.31	0.45
1:D:305:TRP:CE2	1:D:372:LYS:HD3	2.50	0.45
1:A:176:ASP:O	1:A:180:VAL:HG23	2.15	0.45
1:A:236:ASP:O	1:A:237:LYS:C	2.58	0.45
1:B:34:VAL:O	1:B:38:VAL:HG13	2.16	0.45
1:C:7:ILE:HA	1:C:258:LEU:HD12	1.98	0.45
1:C:186:ALA:HA	1:C:340:ASN:O	2.17	0.45
1:C:279:ALA:CB	1:C:298:LYS:HB3	2.45	0.45
1:D:232:ARG:H	1:D:232:ARG:HE	1.65	0.45
1:A:364:GLU:HA	1:A:367:ARG:HG2	1.98	0.45
1:C:57:VAL:HB	1:C:117:GLU:OE1	2.15	0.45
1:C:205:LYS:HA	1:C:205:LYS:HD3	1.43	0.45
1:C:129:ARG:HH21	1:D:122:ALA:HB3	1.81	0.45
1:C:293:ILE:HB	1:C:294:PRO:HD2	1.98	0.45
1:D:35:ILE:O	1:D:38:VAL:HG22	2.17	0.45
1:C:276:VAL:HG21	1:C:305:TRP:CZ2	2.52	0.45
1:D:358:LEU:HD23	1:D:358:LEU:HA	1.86	0.45
1:B:157:MET:HA	1:B:157:MET:CE	2.41	0.45
1:B:207:ARG:HG2	1:B:207:ARG:HH11	1.82	0.45
1:D:357:ILE:HD12	1:D:375:ALA:HB1	1.99	0.45
1:C:60:ALA:O	1:C:123:PRO:HG3	2.17	0.45
1:B:55:GLY:HA2	1:B:85:MET:O	2.18	0.44
1:B:178:PHE:CZ	1:B:323:ALA:HB1	2.52	0.44
1:D:359:ASN:O	1:D:363:PHE:HD1	2.00	0.44
1:A:158:GLY:HA3	1:A:235:PHE:CD2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:PRO:HA	1:A:241:VAL:O	2.17	0.44
1:A:379:ILE:HB	1:A:383:MET:HB2	1.99	0.44
1:C:6:VAL:HA	1:C:274:ARG:HA	1.98	0.44
1:D:279:ALA:O	1:D:385:VAL:HA	2.16	0.44
1:C:165:ALA:O	1:C:169:GLN:N	2.50	0.44
1:C:348:ASN:O	1:C:349:PRO:CB	2.22	0.44
1:A:244:GLY:C	1:A:246:ALA:H	2.25	0.44
1:B:30:LEU:O	1:B:34:VAL:HG23	2.18	0.44
1:D:51:GLU:HA	1:D:81:THR:O	2.16	0.44
1:D:379:ILE:HB	1:D:383:MET:HB2	2.00	0.44
1:A:129:ARG:HH11	1:A:129:ARG:HG2	1.83	0.44
1:A:228:MET:HA	1:A:231:LEU:HD12	2.00	0.44
1:C:319:PHE:O	1:C:320:ALA:C	2.60	0.44
1:C:18:PHE:O	1:C:19:ASN:HB2	2.18	0.44
1:B:349:PRO:O	1:B:350:ILE:C	2.61	0.44
1:D:129:ARG:C	1:D:131:GLY:H	2.26	0.44
1:A:352:ALA:O	1:A:353:SER:C	2.61	0.44
1:B:306:LYS:O	1:B:307:ILE:C	2.61	0.44
1:B:379:ILE:HD11	1:B:385:VAL:HG12	2.00	0.44
1:A:358:LEU:HD11	1:A:387:MET:CE	2.40	0.43
1:C:57:VAL:HG12	1:C:58:LEU:HG	2.00	0.43
1:D:119:MET:HE3	1:D:350:ILE:HG13	1.99	0.43
1:D:236:ASP:O	1:D:239:GLY:N	2.51	0.43
1:C:158:GLY:O	1:C:161:ALA:HB3	2.18	0.43
1:C:326:VAL:O	1:C:330:LEU:HG	2.18	0.43
1:D:175:GLN:NE2	1:D:240:THR:HG23	2.30	0.43
1:D:305:TRP:CZ3	1:D:388:CYS:HB3	2.53	0.43
1:C:170:LEU:HD11	1:C:324:CYS:HB2	1.98	0.43
1:C:271:PRO:HG2	1:C:392:LEU:HD12	2.01	0.43
1:C:385:VAL:CG2	1:C:386:ALA:N	2.80	0.43
1:D:214:ASP:OD1	1:D:214:ASP:C	2.61	0.43
1:A:181:ALA:O	1:A:185:LYS:HG3	2.19	0.43
1:C:46:ALA:HB1	1:C:76:VAL:HA	1.99	0.43
1:C:292:PRO:HB3	1:C:376:THR:OG1	2.18	0.43
1:D:199:ILE:CG2	1:D:200:VAL:N	2.81	0.43
1:D:222:GLY:O	1:D:224:THR:N	2.51	0.43
1:D:368:ARG:HB3	2:D:2004:HOH:O	2.17	0.43
1:C:274:ARG:NH2	1:C:390:GLU:OE1	2.51	0.43
1:D:311:ASP:HB2	1:D:370:ALA:HB1	2.01	0.43
1:D:349:PRO:O	1:D:350:ILE:C	2.62	0.43
1:A:175:GLN:HE22	1:A:240:THR:HG23	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:274:ARG:NH2	1:D:390:GLU:OE1	2.49	0.43
1:A:131:GLY:HA2	1:C:131:GLY:CA	2.46	0.43
1:A:328:LYS:HE3	1:A:328:LYS:HB2	1.80	0.43
1:D:150:ASP:HB2	1:D:157:MET:CE	2.48	0.43
1:B:66:PRO:HA	1:B:69:GLN:HG3	2.01	0.43
1:B:181:ALA:O	1:B:185:LYS:HG3	2.18	0.43
1:D:172:ARG:HE	1:D:176:ASP:CG	2.26	0.43
1:A:88:LEU:HB2	1:A:379:ILE:HG23	1.99	0.42
1:A:361:LEU:HD22	1:A:365:MET:HG3	2.00	0.42
1:C:187:GLU:OE2	1:C:221:HIS:HA	2.19	0.42
1:C:150:ASP:C	1:C:152:PHE:H	2.27	0.42
1:C:28:HIS:HB2	1:C:70:ALA:HB2	2.00	0.42
1:C:156:HIS:CG	1:C:235:PHE:HE1	2.37	0.42
1:C:163:ASN:O	1:C:167:GLN:HB2	2.20	0.42
1:B:158:GLY:HA2	1:B:319:PHE:CZ	2.54	0.42
1:D:281:VAL:HG12	1:D:282:GLY:N	2.33	0.42
1:A:293:ILE:HG21	1:A:329:ASP:OD1	2.20	0.42
1:C:7:ILE:HD12	1:C:362:LEU:HD11	2.01	0.42
1:D:85:MET:HE2	1:D:85:MET:HB2	1.82	0.42
1:A:281:VAL:HG12	1:A:282:GLY:H	1.84	0.42
1:D:161:ALA:HB1	1:D:321:ALA:HB3	2.01	0.42
1:D:284:ASP:C	1:D:284:ASP:OD1	2.62	0.42
1:A:133:LYS:H	1:A:133:LYS:HG2	1.42	0.42
1:A:375:ALA:O	1:A:386:ALA:HA	2.19	0.42
1:D:53:ILE:HD13	1:D:83:TRP:HZ2	1.83	0.42
1:D:134:MET:HE2	1:D:134:MET:HB3	1.92	0.42
1:B:57:VAL:C	1:B:59:PRO:HD3	2.44	0.42
1:C:379:ILE:HG22	1:C:380:GLY:O	2.20	0.42
1:D:280:THR:HA	1:D:384:GLY:O	2.20	0.42
1:D:288:MET:HA	1:D:382:GLY:HA2	2.02	0.42
1:D:306:LYS:O	1:D:307:ILE:C	2.62	0.42
1:D:313:VAL:HG13	1:D:314:GLU:N	2.35	0.42
1:A:19:ASN:C	1:A:23:ALA:HB2	2.45	0.41
1:A:265:SER:O	1:A:266:ARG:C	2.60	0.41
1:D:339:VAL:HG11	1:D:368:ARG:HH21	1.85	0.41
1:A:291:GLY:N	1:A:292:PRO:CD	2.82	0.41
1:B:6:VAL:O	1:B:6:VAL:HG23	2.18	0.41
1:B:78:GLN:HE21	1:B:78:GLN:HB3	1.55	0.41
1:B:187:GLU:HG3	1:B:221:HIS:HA	2.02	0.41
1:C:55:GLY:HA2	1:C:85:MET:O	2.20	0.41
1:C:293:ILE:CB	1:C:294:PRO:CD	2.98	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:257:LEU:HD23	1:D:258:LEU:N	2.35	0.41
1:B:156:HIS:ND1	1:B:158:GLY:N	2.52	0.41
1:A:24:ASN:HA	1:A:121:MET:SD	2.60	0.41
1:A:119:MET:HE2	1:A:143:MET:SD	2.61	0.41
1:B:124:HIS:HA	1:B:140:ILE:O	2.21	0.41
1:B:291:GLY:N	1:B:292:PRO:HD2	2.35	0.41
1:D:164:VAL:O	1:D:168:TRP:HD1	2.04	0.41
1:B:270:GLN:HA	1:B:271:PRO:HD3	1.86	0.41
1:A:178:PHE:CZ	1:A:323:ALA:HB1	2.56	0.41
1:C:81:THR:HG22	1:D:383:MET:HG2	2.03	0.41
1:C:189:ALA:HB1	1:C:340:ASN:HB3	2.02	0.41
1:D:108:ALA:CB	1:D:111:ILE:HD11	2.51	0.41
1:A:72:MET:HG3	1:B:152:PHE:CZ	2.56	0.41
1:A:148:LEU:C	1:A:157:MET:HG2	2.45	0.41
1:A:350:ILE:HD13	1:A:350:ILE:HG21	1.73	0.41
1:B:207:ARG:HG2	1:B:207:ARG:NH1	2.36	0.41
1:C:25:THR:HA	1:C:26:PRO:HD3	1.72	0.41
1:A:293:ILE:HG13	1:A:326:VAL:HG22	2.03	0.41
1:B:5:ILE:HG13	1:B:100:MET:HG2	2.02	0.41
1:B:5:ILE:HD12	1:B:103:ILE:HB	2.03	0.41
1:B:186:ALA:O	1:B:187:GLU:C	2.62	0.41
1:B:333:ASP:C	1:B:335:SER:N	2.78	0.41
1:C:189:ALA:CB	1:C:340:ASN:HB3	2.50	0.41
1:D:158:GLY:HA3	1:D:235:PHE:CD2	2.56	0.41
1:D:291:GLY:N	1:D:292:PRO:HD2	2.35	0.41
1:D:313:VAL:HB	1:D:337:VAL:HG22	2.03	0.41
1:D:348:ASN:N	1:D:349:PRO:HD3	2.28	0.41
1:C:110:ILE:CD1	1:C:259:MET:HE2	2.51	0.41
1:D:64:GLN:O	1:D:65:ASN:C	2.64	0.41
1:D:291:GLY:N	1:D:292:PRO:CD	2.84	0.41
1:B:136:ASP:OD1	1:C:140:ILE:HA	2.20	0.40
1:C:180:VAL:HG21	1:C:225:LEU:HA	2.03	0.40
1:C:233:PRO:HB2	1:C:236:ASP:O	2.20	0.40
1:D:361:LEU:HD21	1:D:373:GLY:HA3	2.02	0.40
1:A:158:GLY:HA3	1:A:235:PHE:CE2	2.56	0.40
1:D:50:ASN:OD1	1:D:109:SER:N	2.53	0.40
1:D:250:ASN:HB2	1:D:348:ASN:O	2.21	0.40
1:B:311:ASP:HB2	1:B:370:ALA:HB1	2.02	0.40
1:C:153:TYR:HE1	1:C:285:PRO:HG2	1.86	0.40
1:C:7:ILE:CD1	1:C:362:LEU:HD11	2.52	0.40
1:C:153:TYR:CZ	1:C:286:LYS:CG	3.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:276:VAL:O	1:C:277:SER:HB3	2.22	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	387/392 (99%)	345 (89%)	37 (10%)	5 (1%)	9	36
1	B	387/392 (99%)	354 (92%)	31 (8%)	2 (0%)	24	58
1	C	387/392 (99%)	337 (87%)	44 (11%)	6 (2%)	7	31
1	D	387/392 (99%)	330 (85%)	50 (13%)	7 (2%)	6	28
All	All	1548/1568 (99%)	1366 (88%)	162 (10%)	20 (1%)	9	36

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	349	PRO
1	C	348	ASN
1	C	349	PRO
1	C	350	ILE
1	D	223	ALA
1	D	349	PRO
1	D	350	ILE
1	B	350	ILE
1	C	223	ALA
1	D	240	THR
1	A	205	LYS
1	C	240	THR
1	D	335	SER
1	A	350	ILE

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Mol	Chain	Res	Type
1	A	339	VAL
1	D	158	GLY
1	C	239	GLY
1	D	270	GLN
1	A	287	VAL
1	B	59	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	276/279 (99%)	250 (91%)	26 (9%)	8	30
1	B	276/279 (99%)	247 (90%)	29 (10%)	6	26
1	C	276/279 (99%)	244 (88%)	32 (12%)	5	21
1	D	276/279 (99%)	241 (87%)	35 (13%)	4	18
All	All	1104/1116 (99%)	982 (89%)	122 (11%)	6	23

All (122) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	LEU
1	A	109	SER
1	A	133	LYS
1	A	196	LYS
1	A	207	ARG
1	A	208	LYS
1	A	224	THR
1	A	225	LEU
1	A	230	LYS
1	A	236	ASP
1	A	237	LYS
1	A	238	GLU
1	A	240	THR
1	A	260	SER
1	A	263	GLU

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Mol	Chain	Res	Type
1	A	272	LEU
1	A	276	VAL
1	A	298	LYS
1	A	322	GLN
1	A	324	CYS
1	A	326	VAL
1	A	332	TRP
1	A	339	VAL
1	A	358	LEU
1	A	361	LEU
1	A	371	ARG
1	B	33	THR
1	B	38	VAL
1	B	39	LEU
1	B	78	GLN
1	B	100	MET
1	B	133	LYS
1	B	144	ILE
1	B	173	ASP
1	B	182	SER
1	B	187	GLU
1	B	191	LYS
1	B	207	ARG
1	B	212	THR
1	B	224	THR
1	B	240	THR
1	B	242	THR
1	B	258	LEU
1	B	263	GLU
1	B	270	GLN
1	B	272	LEU
1	B	276	VAL
1	B	298	LYS
1	B	322	GLN
1	B	339	VAL
1	B	357	ILE
1	B	358	LEU
1	B	361	LEU
1	B	366	LYS
1	B	385	VAL
1	C	4	SER
1	C	17	SER

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Mol	Chain	Res	Type
1	C	39	LEU
1	C	40	GLU
1	C	48	GLU
1	C	49	VAL
1	C	100	MET
1	C	110	ILE
1	C	133	LYS
1	C	144	ILE
1	C	157	MET
1	C	207	ARG
1	C	211	ILE
1	C	220	ARG
1	C	224	THR
1	C	226	ASP
1	C	232	ARG
1	C	249	LEU
1	C	250	ASN
1	C	258	LEU
1	C	272	LEU
1	C	276	VAL
1	C	298	LYS
1	C	322	GLN
1	C	332	TRP
1	C	339	VAL
1	C	353	SER
1	C	358	LEU
1	C	361	LEU
1	C	371	ARG
1	C	374	LEU
1	C	392	LEU
1	D	25	THR
1	D	39	LEU
1	D	89	CYS
1	D	110	ILE
1	D	112	VAL
1	D	133	LYS
1	D	134	MET
1	D	144	ILE
1	D	157	MET
1	D	173	ASP
1	D	187	GLU
1	D	207	ARG

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Mol	Chain	Res	Type
1	D	211	ILE
1	D	224	THR
1	D	227	SER
1	D	232	ARG
1	D	249	LEU
1	D	258	LEU
1	D	270	GLN
1	D	276	VAL
1	D	283	VAL
1	D	288	MET
1	D	290	THR
1	D	298	LYS
1	D	322	GLN
1	D	324	CYS
1	D	327	ASN
1	D	328	LYS
1	D	332	TRP
1	D	339	VAL
1	D	350	ILE
1	D	353	SER
1	D	358	LEU
1	D	371	ARG
1	D	374	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	GLN
1	A	102	GLN
1	A	167	GLN
1	A	175	GLN
1	A	190	GLN
1	A	245	ASN
1	B	78	GLN
1	B	163	ASN
1	B	175	GLN
1	B	184	ASN
1	C	78	GLN
1	C	101	GLN
1	C	175	GLN
1	C	184	ASN
1	C	190	GLN

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Mol	Chain	Res	Type
1	C	316	HIS
1	D	19	ASN
1	D	64	GLN
1	D	78	GLN
1	D	175	GLN
1	D	184	ASN
1	D	190	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 oligosaccharides 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	389/392 (99%)	-1.74	0 100 100	2, 12, 41, 48	0
1	B	389/392 (99%)	-1.72	0 100 100	2, 11, 45, 50	0
1	C	389/392 (99%)	-1.58	0 100 100	19, 32, 49, 57	0
1	D	389/392 (99%)	-1.48	0 100 100	16, 39, 66, 71	0
All	All	1556/1568 (99%)	-1.63	0 100 100	2, 27, 54, 71	0

There are no RSRZ outliers to report.

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 oligosaccharides 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.