



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 07:54 PM UTC

PDB ID : 5M1B / pdb_00005m1b
Title : Crystal structure of C-terminally tagged apo-UbiD from E. coli
Authors : White, M.; Leys, D.
Deposited on : 2016-10-07
Resolution : 3.15 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

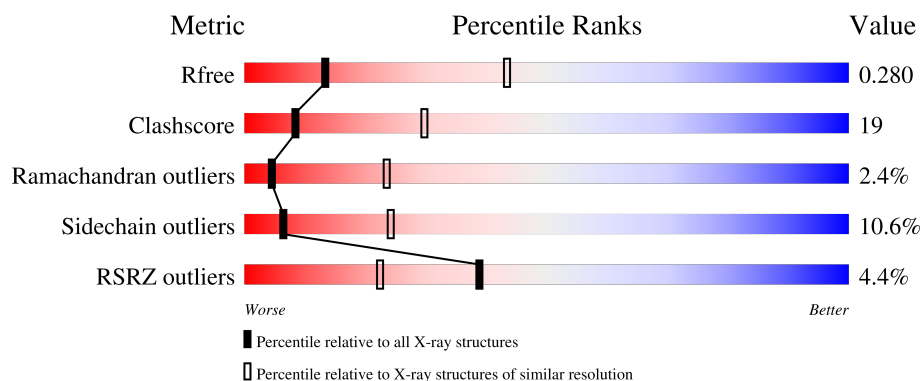
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.15 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	2361 (3.20-3.12)
Clashscore	190562	2486 (3.20-3.12)
Ramachandran outliers	187476	2405 (3.20-3.12)
Sidechain outliers	187428	2404 (3.20-3.12)
RSRZ outliers	180081	2361 (3.20-3.12)

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	505	2% 54% 28% 5% 13%
1	B	505	9% 48% 37% 6% 9%
1	C	505	% 50% 32% 6% 11%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 10593 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 3-octaprenyl-4-hydroxybenzoate carboxy-lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	0	0	0
			3451	2205	588	641	17			
1	B	461	Total	C	N	O	S	0	0	0
			3625	2318	617	673	17			
1	C	448	Total	C	N	O	S	0	0	0
			3517	2252	596	652	17			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	498	LEU	-	expression tag	UNP P0AAB5
A	499	GLU	-	expression tag	UNP P0AAB5
A	500	HIS	-	expression tag	UNP P0AAB5
A	501	HIS	-	expression tag	UNP P0AAB5
A	502	HIS	-	expression tag	UNP P0AAB5
A	503	HIS	-	expression tag	UNP P0AAB5
A	504	HIS	-	expression tag	UNP P0AAB5
A	505	HIS	-	expression tag	UNP P0AAB5
B	498	LEU	-	expression tag	UNP P0AAB5
B	499	GLU	-	expression tag	UNP P0AAB5
B	500	HIS	-	expression tag	UNP P0AAB5
B	501	HIS	-	expression tag	UNP P0AAB5
B	502	HIS	-	expression tag	UNP P0AAB5
B	503	HIS	-	expression tag	UNP P0AAB5
B	504	HIS	-	expression tag	UNP P0AAB5
B	505	HIS	-	expression tag	UNP P0AAB5
C	498	LEU	-	expression tag	UNP P0AAB5
C	499	GLU	-	expression tag	UNP P0AAB5
C	500	HIS	-	expression tag	UNP P0AAB5
C	501	HIS	-	expression tag	UNP P0AAB5
C	502	HIS	-	expression tag	UNP P0AAB5
C	503	HIS	-	expression tag	UNP P0AAB5
C	504	HIS	-	expression tag	UNP P0AAB5

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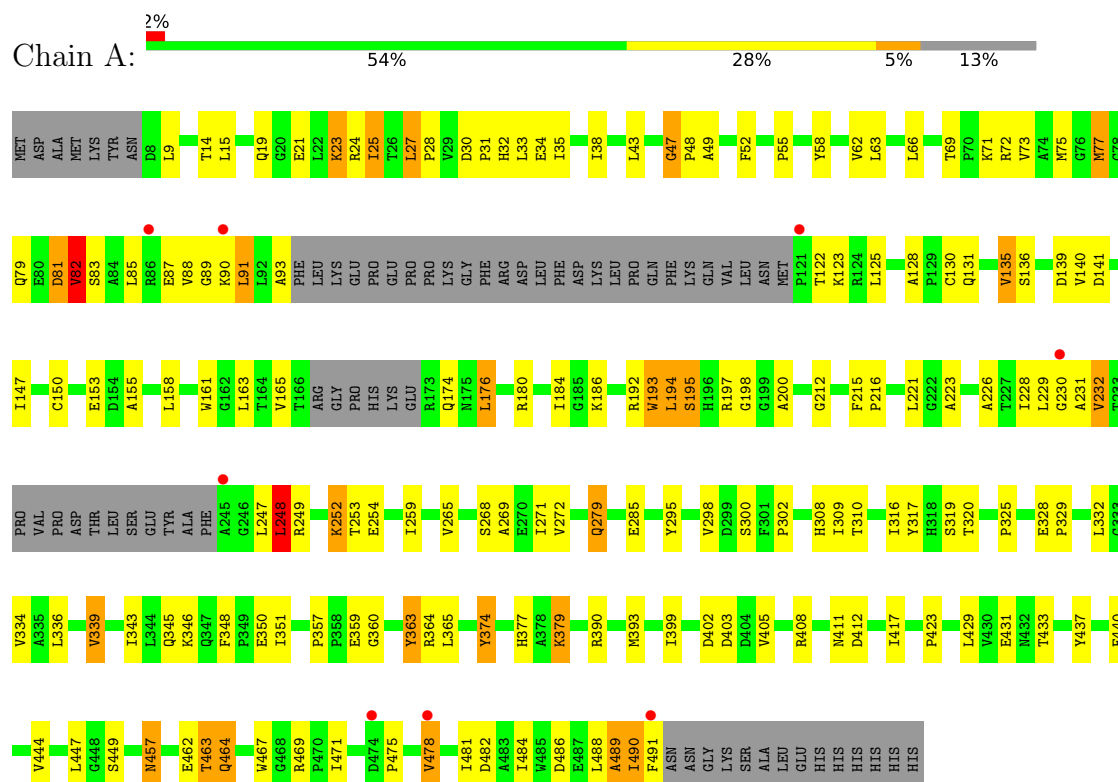
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Chain	Residue	Modelled	Actual	Comment	Reference
C	505	HIS	-	expression tag	UNP P0AAB5

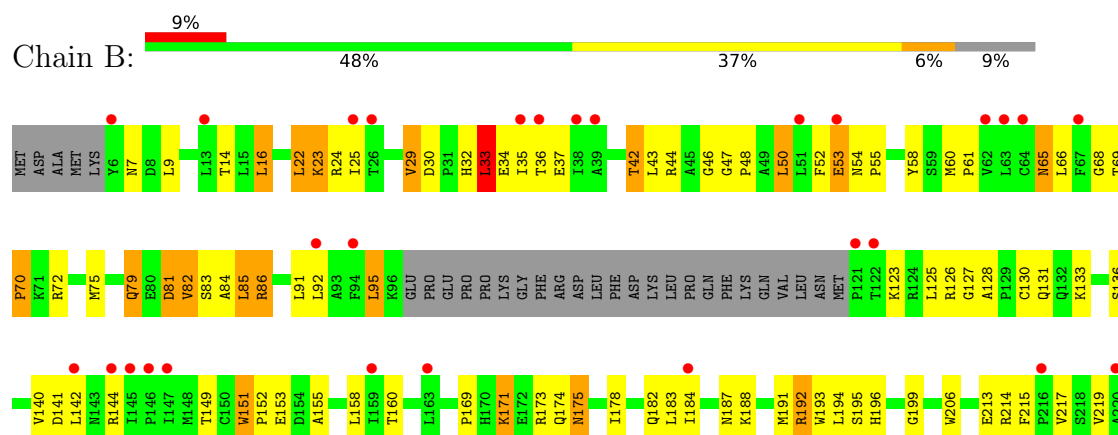
3 Residue-property plots [i](#)

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: 3-octaprenyl-4-hydroxybenzoate carboxy-lyase



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4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	210.19Å 210.19Å 109.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.90 – 3.15 29.90 – 3.15	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.90-3.15) 99.8 (29.90-3.15)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.23 (at 3.18Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.221 , 0.277 0.221 , 0.280	Depositor DCC
R_{free} test set	2151 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	74.2	Xtriage
Anisotropy	0.233	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 65.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10593	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.91% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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	0.74	0/3534	1.11	11/4812 (0.2%)
1	B	0.78	1/3717 (0.0%)	1.14	17/5065 (0.3%)
1	C	0.70	1/3606 (0.0%)	1.11	15/4914 (0.3%)
All	All	0.74	2/10857 (0.0%)	1.12	43/14791 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	428	VAL	CA-CB	-5.49	1.47	1.54
1	C	233	THR	N-CA	5.47	1.51	1.46

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	47	GLY	CA-C-N	8.35	128.38	119.78
1	A	47	GLY	C-N-CA	8.35	128.38	119.78
1	B	399	ILE	N-CA-C	-8.10	96.34	108.17
1	C	47	GLY	CA-C-N	7.75	127.70	120.03
1	C	47	GLY	C-N-CA	7.75	127.70	120.03
1	C	286	GLY	CA-C-N	6.88	126.92	119.90
1	C	286	GLY	C-N-CA	6.88	126.92	119.90
1	B	474	ASP	CA-C-N	-6.59	112.90	119.56
1	B	474	ASP	C-N-CA	-6.59	112.90	119.56
1	B	69	THR	CA-C-N	6.44	127.89	119.84
1	B	69	THR	C-N-CA	6.44	127.89	119.84
1	C	199	GLY	N-CA-C	-6.32	102.19	111.03
1	A	457	ASN	CA-C-N	-6.07	114.34	122.72
1	A	457	ASN	C-N-CA	-6.07	114.34	122.72
1	C	90	LYS	N-CA-C	-5.91	104.92	111.36
1	B	151	TRP	CA-C-N	5.90	127.22	119.84
1	B	151	TRP	C-N-CA	5.90	127.22	119.84
1	B	298	VAL	N-CA-C	5.85	116.97	109.30
1	A	462	GLU	N-CA-C	-5.78	106.06	113.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	348	PHE	CA-C-N	5.68	125.36	119.56
1	B	348	PHE	C-N-CA	5.68	125.36	119.56
1	B	384	GLY	N-CA-C	-5.65	105.65	112.49
1	C	323	GLY	N-CA-C	5.52	119.26	111.42
1	C	248	LEU	N-CA-C	-5.50	106.61	113.55
1	A	195	SER	N-CA-C	5.47	122.46	110.80
1	C	261	ASN	N-CA-C	5.46	117.10	108.96
1	C	368	VAL	CB-CA-C	-5.44	102.58	110.63
1	A	62	VAL	N-CA-C	5.40	115.74	108.17
1	B	303	VAL	N-CA-C	5.39	114.51	106.85
1	C	147	ILE	N-CA-C	-5.25	102.80	109.58
1	B	381	VAL	N-CA-C	5.25	115.39	110.30
1	B	7	ASN	N-CA-C	5.24	116.27	108.31
1	B	485	TRP	N-CA-C	5.18	116.61	111.07
1	C	62	VAL	N-CA-C	5.15	115.32	108.11
1	B	311	GLN	N-CA-C	5.11	116.06	108.86
1	C	433	THR	CA-C-N	-5.10	114.54	119.90
1	C	433	THR	C-N-CA	-5.10	114.54	119.90
1	A	198	GLY	N-CA-C	5.08	118.39	112.50
1	B	174	GLN	N-CA-C	5.06	117.95	110.46
1	A	91	LEU	N-CA-C	-5.04	105.97	111.82
1	C	92	LEU	N-CA-C	5.01	118.55	112.23
1	A	30	ASP	CA-C-N	5.00	125.27	119.47
1	A	30	ASP	C-N-CA	5.00	125.27	119.47

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	3451	0	3425	130	0
1	B	3625	0	3594	162	0
1	C	3517	0	3475	130	0
All	All	10593	0	10494	403	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 19.

All (403) 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:90:LYS:HD2	1:A:339:VAL:HG13	1.27	1.12
1:A:25:ILE:HD12	1:A:52:PHE:CD1	1.87	1.08
1:C:93:ALA:HB3	1:C:346:LYS:NZ	1.70	1.06
1:C:93:ALA:HB3	1:C:346:LYS:HZ3	1.18	1.02
1:C:135:VAL:HG13	1:C:139:ASP:HB3	1.43	0.98
1:A:27:LEU:HD23	1:A:28:PRO:HD2	1.49	0.95
1:B:227:THR:HA	1:B:245:ALA:HB1	1.48	0.94
1:A:247:LEU:HD12	1:A:248:LEU:H	1.32	0.93
1:B:234:PRO:HG2	1:B:391:GLN:OE1	1.69	0.91
1:A:82:VAL:H	1:A:85:LEU:HD12	1.35	0.90
1:A:131:GLN:HA	1:A:310:THR:HG23	1.51	0.90
1:C:25:ILE:HG22	1:C:52:PHE:HA	1.54	0.88
1:A:131:GLN:HA	1:A:310:THR:CG2	2.02	0.88
1:B:32:HIS:HA	1:B:149:THR:HB	1.59	0.85
1:C:355:TYR:CE2	1:C:357:PRO:HG3	2.10	0.84
1:A:90:LYS:HD2	1:A:339:VAL:CG1	2.08	0.83
1:A:158:LEU:HD11	1:A:180:ARG:HG3	1.60	0.82
1:C:69:THR:HG21	1:C:71:LYS:HE2	1.60	0.82
1:C:93:ALA:CB	1:C:346:LYS:NZ	2.42	0.82
1:A:25:ILE:HD12	1:A:52:PHE:HD1	1.41	0.81
1:B:313:GLU:C	1:B:315:ALA:H	1.85	0.81
1:C:135:VAL:CG1	1:C:139:ASP:HB3	2.13	0.79
1:B:222:GLY:O	1:B:270:GLU:HG2	1.83	0.79
1:C:188:LYS:HD3	1:C:305:THR:HG23	1.64	0.78
1:A:163:LEU:HB3	1:A:176:LEU:HD21	1.64	0.78
1:A:158:LEU:HD11	1:A:180:ARG:CG	2.14	0.78
1:B:219:VAL:HB	1:B:273:LEU:HD12	1.66	0.77
1:A:47:GLY:O	1:A:72:ARG:NH2	2.19	0.76
1:A:130:CYS:O	1:A:310:THR:HG22	1.85	0.76
1:A:247:LEU:HD12	1:A:248:LEU:N	2.01	0.76
1:B:126:ARG:C	1:B:128:ALA:H	1.92	0.76
1:B:23:LYS:HB2	1:B:48:PRO:HG2	1.69	0.74
1:C:393:MET:HE3	1:C:394:TYR:CE2	2.23	0.73
1:C:7:ASN:O	1:C:8:ASP:HB2	1.87	0.73
1:A:249:ARG:NH2	1:A:253:THR:HA	2.03	0.73
1:B:234:PRO:CG	1:B:391:GLN:OE1	2.37	0.73
1:B:242:TYR:HE1	1:B:266:PRO:HG3	1.54	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:93:ALA:CB	1:C:346:LYS:HZ3	1.96	0.73
1:B:24:ARG:NH1	1:B:53:GLU:OE2	2.21	0.72
1:B:486:ASP:C	1:B:488:LEU:H	1.97	0.72
1:A:194:LEU:H	1:A:197:ARG:NH1	1.88	0.72
1:B:382:MET:SD	1:B:428:VAL:HG21	2.30	0.72
1:B:9:LEU:HD13	1:B:223:ALA:O	1.90	0.71
1:B:192:ARG:HG3	1:B:285:GLU:OE2	1.91	0.70
1:A:25:ILE:CD1	1:A:52:PHE:HD1	2.04	0.70
1:A:317:TYR:CE2	1:A:319:SER:HB2	2.27	0.70
1:B:55:PRO:HB2	1:B:58:TYR:HB2	1.74	0.70
1:A:249:ARG:HH21	1:A:253:THR:HA	1.56	0.70
1:B:125:LEU:HB2	1:B:128:ALA:HB2	1.73	0.69
1:B:356:LEU:HD22	1:B:363:TYR:HA	1.74	0.69
1:C:35:ILE:HD13	1:C:62:VAL:HG11	1.75	0.69
1:A:136:SER:O	1:A:139:ASP:HB2	1.92	0.68
1:B:290:ASP:HB2	1:B:296:ASN:HD21	1.58	0.68
1:A:489:ALA:HB1	1:A:491:PHE:CE2	2.28	0.68
1:C:230:GLY:HA2	1:C:244:PHE:CD2	2.29	0.68
1:B:408:ARG:HH11	1:B:408:ARG:HG3	1.59	0.68
1:C:435:ILE:O	1:C:446:GLY:HA2	1.93	0.67
1:A:405:VAL:HG13	1:A:412:ASP:HB3	1.76	0.67
1:A:226:ALA:HB3	1:A:249:ARG:HH11	1.59	0.67
1:B:391:GLN:H	1:B:391:GLN:HE21	1.43	0.67
1:A:471:ILE:HG21	1:B:151:TRP:CE3	2.30	0.66
1:A:25:ILE:CD1	1:A:52:PHE:CD1	2.73	0.66
1:C:230:GLY:HA3	1:C:245:ALA:HB2	1.77	0.66
1:A:63:LEU:HD21	1:A:66:LEU:HD13	1.78	0.65
1:B:47:GLY:O	1:B:72:ARG:NH2	2.29	0.65
1:C:374:TYR:O	1:C:377:HIS:HB2	1.97	0.65
1:A:226:ALA:HB3	1:A:249:ARG:NH1	2.10	0.65
1:B:188:LYS:HD3	1:B:305:THR:HG23	1.79	0.65
1:C:74:ALA:HA	1:C:77:MET:HE3	1.79	0.65
1:C:130:CYS:O	1:C:310:THR:HG22	1.95	0.65
1:A:194:LEU:H	1:A:197:ARG:HH11	1.43	0.64
1:B:214:ARG:HB3	1:B:276:TYR:CD2	2.32	0.64
1:A:231:ALA:O	1:A:232:VAL:O	2.16	0.64
1:B:91:LEU:O	1:B:95:LEU:HG	1.97	0.64
1:B:313:GLU:C	1:B:315:ALA:N	2.55	0.64
1:B:52:PHE:O	1:B:55:PRO:HD3	1.97	0.63
1:A:71:LYS:HE2	1:A:75:MET:HE3	1.80	0.63
1:A:295:TYR:CE2	1:B:471:ILE:HA	2.34	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:167:ARG:HH11	1:C:263:LEU:HD23	1.64	0.63
1:A:247:LEU:CD1	1:A:248:LEU:H	2.09	0.63
1:C:136:SER:O	1:C:139:ASP:HB2	1.98	0.63
1:C:354:PHE:CE2	1:C:356:LEU:HD11	2.33	0.63
1:A:140:VAL:HG13	1:A:309:ILE:HD11	1.79	0.62
1:B:191:MET:HE1	1:B:217:VAL:HG11	1.80	0.62
1:B:258:CYS:SG	1:B:265:VAL:HG13	2.39	0.62
1:B:33:LEU:N	1:B:149:THR:O	2.25	0.62
1:C:153:GLU:HB2	1:C:295:TYR:CZ	2.35	0.61
1:B:29:VAL:HG21	1:B:52:PHE:CD1	2.36	0.61
1:B:259:ILE:HD13	1:B:310:THR:HG21	1.82	0.61
1:C:69:THR:HG21	1:C:71:LYS:CE	2.30	0.61
1:C:337:ASN:HD22	1:C:356:LEU:HD13	1.66	0.61
1:B:474:ASP:HB3	1:B:477:VAL:HG23	1.83	0.61
1:B:484:ILE:HG23	1:B:488:LEU:HG	1.81	0.61
1:A:9:LEU:HB2	1:A:223:ALA:O	2.01	0.61
1:A:82:VAL:H	1:A:85:LEU:CD1	2.11	0.61
1:C:85:LEU:HD23	1:C:88:VAL:HG21	1.83	0.61
1:B:42:THR:HG21	1:B:50:LEU:HD21	1.84	0.60
1:B:259:ILE:HD12	1:B:259:ILE:H	1.66	0.60
1:C:145:ILE:HG22	1:C:146:PRO:HD2	1.81	0.60
1:B:222:GLY:HA2	1:B:318:HIS:HB2	1.83	0.60
1:B:290:ASP:HB2	1:B:296:ASN:ND2	2.16	0.60
1:B:320:THR:HG23	1:B:321:TYR:N	2.17	0.60
1:C:10:ARG:HA	1:C:13:LEU:HD12	1.81	0.60
1:C:85:LEU:HA	1:C:88:VAL:HG23	1.82	0.60
1:C:290:ASP:HB3	1:C:292:THR:H	1.66	0.60
1:C:474:ASP:OD1	1:C:476:ASP:HB2	2.02	0.60
1:B:226:ALA:HB3	1:B:249:ARG:HE	1.67	0.59
1:C:158:LEU:HD11	1:C:180:ARG:HB2	1.84	0.59
1:A:77:MET:HB2	1:A:79:GLN:OE1	2.03	0.59
1:C:12:PHE:CD2	1:C:225:PRO:HG3	2.38	0.59
1:B:81:ASP:C	1:B:83:SER:H	2.11	0.59
1:B:131:GLN:HG2	1:B:259:ILE:HD11	1.85	0.59
1:C:337:ASN:ND2	1:C:356:LEU:HD13	2.18	0.59
1:A:82:VAL:N	1:A:85:LEU:HD12	2.12	0.58
1:C:304:PHE:CD1	1:C:304:PHE:C	2.81	0.58
1:C:28:PRO:HA	1:C:54:ASN:O	2.03	0.58
1:C:272:VAL:HB	1:C:310:THR:HB	1.85	0.58
1:B:224:ASP:HB2	1:B:225:PRO:HD2	1.85	0.58
1:C:182:GLN:OE1	1:C:285:GLU:HG2	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:ARG:C	1:B:128:ALA:N	2.62	0.58
1:C:82:VAL:HG12	1:C:85:LEU:HD12	1.85	0.57
1:B:158:LEU:O	1:B:160:THR:HG23	2.05	0.57
1:A:212:GLY:H	1:A:279:GLN:HE22	1.52	0.57
1:C:199:GLY:O	1:C:200:ALA:CB	2.52	0.57
1:A:463:THR:OG1	1:A:464:GLN:N	2.36	0.56
1:B:222:GLY:HA2	1:B:318:HIS:CB	2.35	0.56
1:C:378:ALA:O	1:C:382:MET:HG3	2.05	0.56
1:A:360:GLY:O	1:A:364:ARG:HB2	2.04	0.56
1:B:142:LEU:HD12	1:B:187:ASN:HA	1.86	0.56
1:B:486:ASP:O	1:B:488:LEU:N	2.38	0.56
1:C:85:LEU:HA	1:C:88:VAL:CG2	2.35	0.56
1:C:369:THR:HA	1:C:401:CYS:O	2.05	0.56
1:B:249:ARG:HH11	1:B:251:THR:HG23	1.71	0.56
1:A:165:VAL:HG21	1:A:265:VAL:HG12	1.87	0.56
1:A:379:LYS:HD3	1:C:432:ASN:O	2.05	0.56
1:C:31:PRO:HA	1:C:35:ILE:HD12	1.88	0.56
1:A:194:LEU:N	1:A:197:ARG:NH1	2.53	0.56
1:B:353:ASP:CG	1:B:408:ARG:HH12	2.13	0.55
1:A:32:HIS:CE1	1:A:33:LEU:HD12	2.41	0.55
1:A:81:ASP:H	1:A:85:LEU:HD11	1.71	0.55
1:A:161:TRP:CD1	1:A:320:THR:HB	2.42	0.55
1:C:69:THR:CG2	1:C:71:LYS:HE2	2.36	0.55
1:A:25:ILE:HD13	1:A:25:ILE:O	2.07	0.55
1:A:193:TRP:HB2	1:A:200:ALA:HB2	1.89	0.55
1:C:153:GLU:HB2	1:C:295:TYR:CE1	2.42	0.55
1:B:65:ASN:HD22	1:B:72:ARG:HH12	1.53	0.55
1:B:30:ASP:HB3	1:B:34:GLU:HG3	1.89	0.55
1:B:386:TRP:NE1	1:B:450:LYS:HD3	2.23	0.54
1:B:459:TRP:O	1:B:462:GLU:HB2	2.07	0.54
1:C:66:LEU:HD23	1:C:318:HIS:CE1	2.42	0.54
1:A:228:ILE:C	1:A:230:GLY:H	2.15	0.54
1:C:405:VAL:HG13	1:C:412:ASP:HB3	1.89	0.54
1:B:258:CYS:HB3	1:B:274:GLU:CD	2.33	0.54
1:B:408:ARG:HG3	1:B:408:ARG:NH1	2.22	0.54
1:B:42:THR:O	1:B:46:GLY:N	2.32	0.54
1:A:79:GLN:CB	1:A:85:LEU:HD13	2.38	0.53
1:C:87:GLU:HA	1:C:90:LYS:HD2	1.91	0.53
1:B:61:PRO:HB2	1:B:316:ILE:HG23	1.90	0.53
1:A:79:GLN:HB2	1:A:85:LEU:HD13	1.90	0.53
1:A:325:PRO:HB2	1:A:359:GLU:HA	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:363:TYR:CE1	1:B:364:ARG:HG2	2.44	0.53
1:C:365:LEU:HD23	1:C:366:ALA:N	2.24	0.53
1:C:274:GLU:HB2	1:C:308:HIS:HB2	1.91	0.53
1:A:475:PRO:O	1:A:478:VAL:HB	2.09	0.52
1:B:390:ARG:O	1:B:393:MET:HG2	2.09	0.52
1:C:136:SER:HB3	1:C:308:HIS:CD2	2.43	0.52
1:B:486:ASP:C	1:B:488:LEU:N	2.67	0.52
1:B:182:GLN:HG3	1:B:288:TYR:OH	2.08	0.52
1:B:291:HIS:O	1:B:325:PRO:HD3	2.08	0.52
1:B:313:GLU:O	1:B:315:ALA:N	2.38	0.52
1:A:90:LYS:CG	1:A:93:ALA:HB3	2.38	0.52
1:A:128:ALA:HB3	1:A:131:GLN:HG3	1.92	0.52
1:A:212:GLY:H	1:A:279:GLN:NE2	2.07	0.52
1:A:429:LEU:HD22	1:B:424:ALA:HA	1.90	0.52
1:C:54:ASN:OD1	1:C:54:ASN:N	2.42	0.52
1:B:65:ASN:ND2	1:B:72:ARG:HH12	2.07	0.52
1:B:193:TRP:HH2	1:B:215:PHE:HE2	1.57	0.52
1:A:193:TRP:HA	1:A:197:ARG:HH11	1.75	0.52
1:B:352:VAL:HG21	1:B:403:ASP:HB3	1.92	0.52
1:B:474:ASP:O	1:B:478:VAL:HG23	2.10	0.52
1:B:42:THR:HG22	1:B:47:GLY:HA3	1.91	0.51
1:C:130:CYS:CB	1:C:269:ALA:HB3	2.39	0.51
1:C:131:GLN:HA	1:C:310:THR:HG23	1.92	0.51
1:A:23:LYS:HD2	1:B:489:ALA:O	2.10	0.51
1:B:42:THR:CG2	1:B:47:GLY:HA3	2.40	0.51
1:B:75:MET:HA	1:B:79:GLN:O	2.11	0.51
1:C:130:CYS:O	1:C:310:THR:CG2	2.59	0.51
1:A:90:LYS:CD	1:A:339:VAL:HG13	2.19	0.51
1:C:421:MET:HE1	1:C:453:LEU:HG	1.93	0.51
1:B:417:ILE:O	1:B:421:MET:HG2	2.11	0.51
1:A:399:ILE:HD13	1:A:417:ILE:HD11	1.93	0.51
1:A:437:TYR:HD1	1:A:437:TYR:H	1.58	0.51
1:B:42:THR:HG22	1:B:47:GLY:CA	2.41	0.51
1:C:339:VAL:O	1:C:342:PRO:HD2	2.10	0.51
1:C:341:VAL:HB	1:C:342:PRO:HD3	1.92	0.51
1:A:193:TRP:HE1	1:A:302:PRO:HG2	1.75	0.51
1:B:329:PRO:O	1:B:333:GLY:N	2.42	0.51
1:C:354:PHE:HE2	1:C:356:LEU:HD11	1.75	0.51
1:B:82:VAL:HG12	1:B:85:LEU:HD13	1.92	0.50
1:B:328:GLU:HB3	1:B:329:PRO:HD3	1.93	0.50
1:C:351:ILE:HA	1:C:370:ILE:HG22	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:GLU:HB2	1:A:295:TYR:CZ	2.46	0.50
1:C:25:ILE:HG21	1:C:52:PHE:CD1	2.46	0.50
1:C:170:HIS:C	1:C:171:LYS:HG2	2.37	0.50
1:A:21:GLU:HB3	1:A:72:ARG:HD3	1.94	0.50
1:B:55:PRO:HG3	1:B:60:MET:O	2.12	0.50
1:B:70:PRO:HD3	1:B:331:VAL:HG13	1.94	0.50
1:C:320:THR:HG23	1:C:321:TYR:N	2.27	0.50
1:C:145:ILE:CG2	1:C:146:PRO:HD2	2.42	0.50
1:A:69:THR:CG2	1:A:71:LYS:HB3	2.42	0.50
1:A:69:THR:HG22	1:A:71:LYS:HB3	1.94	0.49
1:B:320:THR:HG23	1:B:321:TYR:H	1.77	0.49
1:B:329:PRO:HA	1:B:332:LEU:HB2	1.94	0.49
1:C:233:THR:HB	1:C:234:PRO:HD2	1.94	0.49
1:C:373:GLN:O	1:C:459:TRP:CH2	2.65	0.49
1:A:192:ARG:HD2	1:A:285:GLU:CD	2.38	0.49
1:C:231:ALA:O	1:C:232:VAL:HG13	2.11	0.49
1:B:66:LEU:O	1:B:72:ARG:HD2	2.13	0.49
1:B:413:VAL:O	1:B:416:ALA:HB3	2.13	0.48
1:A:136:SER:HB3	1:A:308:HIS:CD2	2.48	0.48
1:B:459:TRP:CD1	1:B:459:TRP:N	2.80	0.48
1:C:10:ARG:HG3	1:C:312:ARG:HH22	1.78	0.48
1:C:484:ILE:HD12	1:C:488:LEU:HD11	1.95	0.48
1:C:194:LEU:O	1:C:196:HIS:N	2.46	0.48
1:C:422:ASP:O	1:C:423:PRO:C	2.56	0.48
1:A:490:ILE:HG21	1:B:42:THR:HG23	1.95	0.48
1:B:32:HIS:CD2	1:B:33:LEU:HG	2.49	0.48
1:A:83:SER:O	1:A:87:GLU:HB2	2.14	0.48
1:A:374:TYR:HE1	1:A:377:HIS:HA	1.77	0.48
1:B:365:LEU:HD23	1:B:366:ALA:N	2.28	0.48
1:C:35:ILE:CG2	1:C:148:MET:HE3	2.43	0.48
1:C:132:GLN:HB3	1:C:313:GLU:HG3	1.94	0.48
1:A:390:ARG:O	1:A:393:MET:HG2	2.13	0.48
1:B:374:TYR:O	1:B:377:HIS:HB2	2.13	0.48
1:C:165:VAL:HA	1:C:175:ASN:O	2.14	0.48
1:A:55:PRO:HG2	1:A:58:TYR:HB2	1.96	0.48
1:A:141:ASP:OD1	1:A:186:LYS:HG3	2.13	0.48
1:B:257:LYS:NZ	1:B:262:ASP:OD1	2.36	0.48
1:C:233:THR:CB	1:C:234:PRO:HD2	2.44	0.48
1:C:456:THR:O	1:C:458:LYS:HD3	2.13	0.48
1:A:15:LEU:O	1:A:19:GLN:HG2	2.14	0.48
1:A:348:PHE:HB2	1:A:351:ILE:CD1	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:411:ASN:HA	1:B:359:GLU:HG2	1.95	0.48
1:B:409:ASP:OD1	1:B:411:ASN:N	2.43	0.47
1:B:123:LYS:O	1:B:254:GLU:HA	2.13	0.47
1:B:130:CYS:HB3	1:B:267:ALA:O	2.14	0.47
1:B:194:LEU:O	1:B:196:HIS:N	2.48	0.47
1:C:288:TYR:O	1:C:295:TYR:HA	2.14	0.47
1:A:125:LEU:HD11	1:A:254:GLU:OE1	2.14	0.47
1:A:221:LEU:HB2	1:A:271:ILE:HB	1.95	0.47
1:B:16:LEU:HB2	1:B:22:LEU:HB2	1.95	0.47
1:C:130:CYS:HB3	1:C:269:ALA:HB3	1.96	0.47
1:C:337:ASN:C	1:C:339:VAL:N	2.72	0.47
1:A:34:GLU:O	1:A:38:ILE:HG13	2.15	0.47
1:A:135:VAL:HG13	1:A:139:ASP:HB3	1.97	0.47
1:A:449:SER:OG	1:B:422:ASP:OD1	2.31	0.47
1:A:469:ARG:NH2	1:B:297:GLU:OE2	2.48	0.47
1:B:29:VAL:HG21	1:B:52:PHE:HD1	1.78	0.47
1:B:226:ALA:CB	1:B:249:ARG:HE	2.28	0.47
1:C:91:LEU:O	1:C:94:PHE:HB2	2.14	0.47
1:B:386:TRP:CE2	1:B:450:LYS:HD3	2.50	0.47
1:A:32:HIS:O	1:A:33:LEU:HB2	2.13	0.47
1:A:272:VAL:N	1:A:310:THR:O	2.47	0.47
1:A:328:GLU:HB3	1:A:329:PRO:HD3	1.97	0.47
1:A:363:TYR:CZ	1:A:364:ARG:HG3	2.49	0.47
1:C:356:LEU:HD23	1:C:363:TYR:HA	1.96	0.47
1:A:194:LEU:N	1:A:197:ARG:HH11	2.12	0.47
1:A:488:LEU:HD21	1:B:25:ILE:HG23	1.96	0.46
1:B:79:GLN:HG3	1:B:84:ALA:HB3	1.97	0.46
1:C:183:LEU:HA	1:C:189:LEU:HD23	1.96	0.46
1:B:82:VAL:HA	1:B:85:LEU:HD12	1.97	0.46
1:B:169:PRO:HB3	1:B:206:TRP:CE3	2.51	0.46
1:C:334:VAL:CG2	1:C:358:PRO:HD3	2.45	0.46
1:A:440:PHE:CD1	1:A:440:PHE:C	2.94	0.46
1:B:65:ASN:ND2	1:B:68:GLY:HA3	2.30	0.46
1:C:232:VAL:O	1:C:336:LEU:HD13	2.15	0.46
1:B:425:ARG:HG2	1:B:426:ASP:OD1	2.15	0.46
1:B:444:VAL:CG2	1:B:447:LEU:HB3	2.45	0.46
1:C:170:HIS:C	1:C:171:LYS:CG	2.88	0.46
1:C:328:GLU:HB3	1:C:329:PRO:HD3	1.96	0.46
1:A:130:CYS:O	1:A:310:THR:CG2	2.61	0.46
1:A:165:VAL:CG1	1:A:174:GLN:NE2	2.79	0.46
1:A:23:LYS:HE3	1:A:24:ARG:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:354:PHE:HD1	1:B:368:VAL:HG22	1.81	0.46
1:B:459:TRP:HB3	1:B:460:PRO:CD	2.45	0.46
1:C:133:LYS:HB2	1:C:311:GLN:HG3	1.98	0.46
1:A:357:PRO:HB3	1:A:359:GLU:OE1	2.16	0.46
1:B:422:ASP:O	1:B:423:PRO:C	2.58	0.46
1:B:402:ASP:OD2	1:B:456:THR:HA	2.16	0.46
1:A:130:CYS:CB	1:A:269:ALA:HB3	2.45	0.46
1:B:409:ASP:OD1	1:B:409:ASP:C	2.57	0.46
1:C:230:GLY:HA2	1:C:244:PHE:HD2	1.79	0.46
1:C:421:MET:HG2	1:C:454:ASP:O	2.16	0.46
1:A:153:GLU:OE1	1:B:472:LYS:NZ	2.49	0.45
1:B:242:TYR:CE1	1:B:266:PRO:HG3	2.44	0.45
1:C:372:LYS:HE2	1:C:459:TRP:CD1	2.51	0.45
1:B:52:PHE:C	1:B:54:ASN:H	2.24	0.45
1:A:90:LYS:HA	1:A:93:ALA:HB3	1.98	0.45
1:A:252:LYS:H	1:A:252:LYS:HG3	1.57	0.45
1:B:82:VAL:CG1	1:B:85:LEU:CD1	2.94	0.45
1:C:47:GLY:O	1:C:72:ARG:NH2	2.49	0.45
1:B:35:ILE:HG23	1:B:52:PHE:CE2	2.52	0.45
1:B:82:VAL:HG13	1:B:85:LEU:CD1	2.47	0.45
1:A:43:LEU:C	1:A:43:LEU:HD23	2.41	0.45
1:C:62:VAL:HG23	1:C:146:PRO:HG2	1.99	0.45
1:C:169:PRO:HG3	1:C:206:TRP:HA	1.98	0.45
1:A:490:ILE:HG13	1:B:23:LYS:HG3	1.99	0.45
1:C:337:ASN:O	1:C:339:VAL:N	2.50	0.45
1:A:345:GLN:O	1:A:346:LYS:C	2.60	0.44
1:A:228:ILE:O	1:A:230:GLY:N	2.46	0.44
1:A:457:ASN:OD1	1:A:467:TRP:NE1	2.49	0.44
1:B:136:SER:HA	1:B:307:THR:O	2.18	0.44
1:A:440:PHE:HA	1:B:465:ARG:NH2	2.31	0.44
1:C:388:PHE:CD2	1:C:389:LEU:HG	2.52	0.44
1:C:484:ILE:CD1	1:C:488:LEU:HD11	2.48	0.44
1:C:304:PHE:HD1	1:C:305:THR:N	2.16	0.44
1:A:31:PRO:HA	1:A:35:ILE:HD12	1.99	0.44
1:C:288:TYR:H	1:C:296:ASN:HB2	1.82	0.44
1:C:391:GLN:H	1:C:391:GLN:HE21	1.66	0.44
1:A:123:LYS:O	1:A:254:GLU:HA	2.17	0.44
1:C:199:GLY:O	1:C:200:ALA:HB3	2.17	0.44
1:B:324:ARG:O	1:B:324:ARG:HG3	2.18	0.43
1:B:363:TYR:CD1	1:B:363:TYR:C	2.96	0.43
1:B:267:ALA:C	1:B:269:ALA:H	2.26	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:LYS:HD3	1:A:343:ILE:CG1	2.49	0.43
1:B:65:ASN:HD22	1:B:72:ARG:NH1	2.17	0.43
1:B:330:ALA:O	1:B:334:VAL:HG23	2.18	0.43
1:C:68:GLY:H	1:C:72:ARG:HH12	1.65	0.43
1:C:184:ILE:HG21	1:C:190:ILE:HD11	1.99	0.43
1:A:69:THR:HG22	1:A:71:LYS:H	1.84	0.43
1:A:215:PHE:HA	1:A:216:PRO:HD3	1.92	0.43
1:C:297:GLU:OE2	1:C:297:GLU:HA	2.17	0.43
1:C:409:ASP:O	1:C:410:TRP:C	2.61	0.43
1:A:81:ASP:O	1:A:82:VAL:HB	2.17	0.43
1:C:474:ASP:O	1:C:478:VAL:HG23	2.18	0.43
1:A:481:ILE:HG13	1:B:37:GLU:HG2	1.99	0.43
1:B:52:PHE:C	1:B:54:ASN:N	2.76	0.43
1:A:43:LEU:HD23	1:A:43:LEU:O	2.19	0.43
1:B:60:MET:SD	1:B:271:ILE:HD11	2.59	0.43
1:A:147:ILE:HG23	1:A:317:TYR:OH	2.18	0.43
1:B:140:VAL:HA	1:B:309:ILE:HD11	2.00	0.43
1:B:141:ASP:O	1:B:144:ARG:HB2	2.18	0.43
1:C:15:LEU:O	1:C:19:GLN:HG2	2.19	0.43
1:B:290:ASP:C	1:B:292:THR:N	2.77	0.42
1:B:338:GLU:OE2	1:B:408:ARG:NH2	2.52	0.42
1:A:444:VAL:CG2	1:A:447:LEU:HB3	2.50	0.42
1:A:471:ILE:CG2	1:B:151:TRP:CE3	3.02	0.42
1:C:85:LEU:O	1:C:88:VAL:HB	2.19	0.42
1:B:175:ASN:ND2	1:B:199:GLY:HA2	2.34	0.42
1:C:70:PRO:HA	1:C:73:VAL:HG23	2.00	0.42
1:B:367:VAL:HG22	1:B:399:ILE:HB	2.02	0.42
1:B:391:GLN:H	1:B:391:GLN:NE2	2.15	0.42
1:C:77:MET:HE1	1:C:85:LEU:HD11	2.01	0.42
1:B:125:LEU:HD21	1:B:254:GLU:HG2	2.00	0.42
1:A:69:THR:HG22	1:A:71:LYS:N	2.35	0.42
1:C:167:ARG:O	1:C:215:PHE:HE1	2.02	0.42
1:C:361:CYS:O	1:C:362:SER:HB2	2.19	0.42
1:A:359:GLU:HG2	1:B:411:ASN:HA	2.01	0.42
1:B:433:THR:HB	1:B:434:PRO:CD	2.49	0.42
1:B:256:VAL:HG22	1:B:257:LYS:O	2.19	0.42
1:C:163:LEU:HD12	1:C:163:LEU:H	1.84	0.42
1:C:207:CYS:HA	1:C:279:GLN:HE22	1.85	0.42
1:C:332:LEU:HD23	1:C:332:LEU:HA	1.92	0.42
1:B:29:VAL:HG23	1:B:54:ASN:O	2.20	0.42
1:B:125:LEU:CB	1:B:128:ALA:HB2	2.47	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:325:PRO:HB2	1:B:358:PRO:O	2.20	0.42
1:C:393:MET:HE3	1:C:394:TYR:CZ	2.54	0.42
1:C:417:ILE:HG12	1:C:453:LEU:HD23	2.02	0.42
1:B:381:VAL:O	1:B:382:MET:C	2.61	0.41
1:C:252:LYS:HB2	1:C:252:LYS:HE2	1.85	0.41
1:C:337:ASN:C	1:C:339:VAL:H	2.28	0.41
1:A:47:GLY:HA2	1:A:48:PRO:HD2	1.72	0.41
1:A:423:PRO:HB3	1:B:451:MET:HB2	2.01	0.41
1:B:81:ASP:C	1:B:83:SER:N	2.77	0.41
1:B:389:LEU:HD23	1:B:389:LEU:HA	1.91	0.41
1:C:72:ARG:HA	1:C:75:MET:HB2	2.03	0.41
1:C:74:ALA:HB1	1:C:79:GLN:HB3	2.01	0.41
1:A:49:ALA:HB2	1:A:72:ARG:NH1	2.36	0.41
1:B:82:VAL:CG1	1:B:85:LEU:HD13	2.49	0.41
1:B:86:ARG:HH11	1:B:341:VAL:HG12	1.85	0.41
1:B:454:ASP:C	1:B:454:ASP:OD1	2.64	0.41
1:C:484:ILE:O	1:C:488:LEU:HG	2.21	0.41
1:B:83:SER:O	1:B:86:ARG:HB2	2.20	0.41
1:A:90:LYS:HD3	1:A:343:ILE:HG13	2.03	0.41
1:A:348:PHE:HB2	1:A:351:ILE:HD12	2.03	0.41
1:A:365:LEU:HD23	1:A:365:LEU:C	2.46	0.41
1:C:454:ASP:OD1	1:C:454:ASP:C	2.64	0.41
1:A:90:LYS:HG2	1:A:93:ALA:HB3	2.02	0.41
1:A:403:ASP:C	1:A:405:VAL:H	2.28	0.41
1:B:386:TRP:CD1	1:B:450:LYS:HD3	2.56	0.41
1:C:8:ASP:O	1:C:11:ASP:HB2	2.21	0.41
1:C:55:PRO:O	1:C:56:LYS:C	2.62	0.40
1:C:217:VAL:HG21	1:C:277:ILE:HD11	2.03	0.40
1:A:122:THR:HB	1:A:253:THR:O	2.20	0.40
1:B:235:VAL:HG12	1:B:236:PRO:O	2.21	0.40
1:C:442:SER:HB2	1:C:447:LEU:O	2.20	0.40
1:C:490:ILE:HG23	1:C:491:PHE:CD2	2.56	0.40
1:A:447:LEU:HD13	1:C:375:ALA:HB1	2.04	0.40
1:B:44:ARG:NH2	1:B:324:ARG:O	2.55	0.40
1:B:215:PHE:HD2	1:B:277:ILE:HD12	1.86	0.40
1:A:48:PRO:CG	1:B:490:ILE:HG22	2.52	0.40
1:B:33:LEU:HD23	1:B:149:THR:O	2.21	0.40
1:B:316:ILE:HG22	1:B:317:TYR:H	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	432/505 (86%)	381 (88%)	40 (9%)	11 (2%)	4	22
1	B	457/505 (90%)	397 (87%)	49 (11%)	11 (2%)	4	23
1	C	442/505 (88%)	401 (91%)	31 (7%)	10 (2%)	5	24
All	All	1331/1515 (88%)	1179 (89%)	120 (9%)	32 (2%)	4	23

All (32) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	82	VAL
1	A	232	VAL
1	A	489	ALA
1	B	195	SER
1	B	489	ALA
1	C	8	ASP
1	C	195	SER
1	C	200	ALA
1	C	489	ALA
1	B	155	ALA
1	B	314	ASP
1	A	155	ALA
1	A	229	LEU
1	A	484	ILE
1	B	152	PRO
1	B	487	GLU
1	A	248	LEU
1	A	363	TYR
1	B	171	LYS
1	C	81	ASP
1	A	195	SER
1	A	482	ASP
1	B	33	LEU
1	B	53	GLU

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Mol	Chain	Res	Type
1	C	488	LEU
1	A	89	GLY
1	C	234	PRO
1	C	222	GLY
1	C	423	PRO
1	B	70	PRO
1	C	169	PRO
1	B	127	GLY

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	371/430 (86%)	331 (89%)	40 (11%)	6	23
1	B	390/430 (91%)	349 (90%)	41 (10%)	6	24
1	C	377/430 (88%)	337 (89%)	40 (11%)	6	24
All	All	1138/1290 (88%)	1017 (89%)	121 (11%)	6	24

All (121) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	THR
1	A	23	LYS
1	A	25	ILE
1	A	27	LEU
1	A	73	VAL
1	A	77	MET
1	A	81	ASP
1	A	82	VAL
1	A	88	VAL
1	A	91	LEU
1	A	135	VAL
1	A	150	CYS
1	A	176	LEU
1	A	184	ILE

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Mol	Chain	Res	Type
1	A	193	TRP
1	A	194	LEU
1	A	248	LEU
1	A	252	LYS
1	A	259	ILE
1	A	268	SER
1	A	279	GLN
1	A	298	VAL
1	A	300	SER
1	A	316	ILE
1	A	332	LEU
1	A	334	VAL
1	A	336	LEU
1	A	339	VAL
1	A	350	GLU
1	A	374	TYR
1	A	379	LYS
1	A	402	ASP
1	A	408	ARG
1	A	431	GLU
1	A	433	THR
1	A	463	THR
1	A	464	GLN
1	A	478	VAL
1	A	486	ASP
1	A	490	ILE
1	B	14	THR
1	B	16	LEU
1	B	22	LEU
1	B	23	LYS
1	B	29	VAL
1	B	33	LEU
1	B	36	THR
1	B	42	THR
1	B	43	LEU
1	B	50	LEU
1	B	65	ASN
1	B	79	GLN
1	B	81	ASP
1	B	82	VAL
1	B	85	LEU
1	B	86	ARG

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Mol	Chain	Res	Type
1	B	92	LEU
1	B	95	LEU
1	B	133	LYS
1	B	153	GLU
1	B	171	LYS
1	B	173	ARG
1	B	175	ASN
1	B	178	ILE
1	B	183	LEU
1	B	184	ILE
1	B	192	ARG
1	B	213	GLU
1	B	233	THR
1	B	251	THR
1	B	309	ILE
1	B	322	THR
1	B	328	GLU
1	B	332	LEU
1	B	336	LEU
1	B	390	ARG
1	B	391	GLN
1	B	405	VAL
1	B	436	ASP
1	B	472	LYS
1	B	484	ILE
1	C	14	THR
1	C	23	LYS
1	C	25	ILE
1	C	54	ASN
1	C	56	LYS
1	C	62	VAL
1	C	71	LYS
1	C	82	VAL
1	C	92	LEU
1	C	94	PHE
1	C	95	LEU
1	C	132	GLN
1	C	165	VAL
1	C	171	LYS
1	C	213	GLU
1	C	232	VAL
1	C	233	THR

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Mol	Chain	Res	Type
1	C	242	TYR
1	C	248	LEU
1	C	251	THR
1	C	254	GLU
1	C	255	VAL
1	C	256	VAL
1	C	263	LEU
1	C	285	GLU
1	C	296	ASN
1	C	298	VAL
1	C	304	PHE
1	C	322	THR
1	C	324	ARG
1	C	345	GLN
1	C	359	GLU
1	C	368	VAL
1	C	391	GLN
1	C	393	MET
1	C	453	LEU
1	C	472	LYS
1	C	473	LYS
1	C	481	ILE
1	C	490	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	HIS
1	A	132	GLN
1	A	143	ASN
1	A	279	GLN
1	A	480	HIS
1	B	32	HIS
1	B	65	ASN
1	B	79	GLN
1	B	170	HIS
1	B	175	ASN
1	B	187	ASN
1	B	196	HIS
1	B	296	ASN
1	B	318	HIS
1	B	391	GLN

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Mol	Chain	Res	Type
1	B	432	ASN
1	C	131	GLN
1	C	143	ASN
1	C	261	ASN
1	C	318	HIS
1	C	337	ASN
1	C	347	GLN
1	C	391	GLN
1	C	411	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no 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

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, 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	440/505 (87%)	-0.29	8 (1%) 67 47	29, 75, 144, 182	0
1	B	461/505 (91%)	0.36	45 (9%) 13 7	26, 115, 166, 244	2 (0%)
1	C	448/505 (88%)	-0.30	7 (1%) 70 50	31, 76, 135, 174	0
All	All	1349/1515 (89%)	-0.07	60 (4%) 39 22	26, 84, 152, 244	2 (0%)

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	247	LEU	5.0
1	B	121	PRO	4.8
1	B	25	ILE	4.7
1	B	243	ALA	4.7
1	B	51	LEU	4.4
1	B	142	LEU	4.4
1	B	26	THR	3.9
1	B	145	ILE	3.5
1	B	271	ILE	3.5
1	B	63	LEU	3.4
1	C	94	PHE	3.3
1	C	480	HIS	3.2
1	B	244	PHE	3.2
1	B	6	TYR	3.2
1	B	309	ILE	3.2
1	A	86	ARG	3.2
1	B	163	LEU	3.1
1	A	90	LYS	3.1
1	C	234	PRO	3.1
1	B	53	GLU	3.0
1	B	216	PRO	3.0
1	B	221	LEU	3.0
1	B	36	THR	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	220	ALA	2.9
1	A	474	ASP	2.9
1	B	39	ALA	2.9
1	A	245	ALA	2.8
1	B	92	LEU	2.8
1	A	478	VAL	2.7
1	A	121	PRO	2.7
1	C	491	PHE	2.7
1	B	67	PHE	2.5
1	B	317	TYR	2.4
1	B	62	VAL	2.4
1	B	239	LEU	2.4
1	A	491	PHE	2.4
1	B	242	TYR	2.3
1	C	95	LEU	2.3
1	C	490	ILE	2.3
1	B	229	LEU	2.3
1	B	263	LEU	2.3
1	B	228	ILE	2.2
1	A	230	GLY	2.2
1	B	38	ILE	2.2
1	C	247	LEU	2.2
1	B	303	VAL	2.2
1	B	273	LEU	2.1
1	B	122	THR	2.1
1	B	13	LEU	2.1
1	B	251	THR	2.1
1	B	310	THR	2.1
1	B	146	PRO	2.1
1	B	35	ILE	2.1
1	B	184	ILE	2.1
1	B	64	CYS	2.1
1	B	159	ILE	2.0
1	B	318	HIS	2.0
1	B	144	ARG	2.0
1	B	147	ILE	2.0
1	B	94	PHE	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 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.