



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:31 AM UTC

PDB ID : 1YIA / pdb_00001yia
Title : Crystal structure of tryptophanyl tRNA synthetase II from *Deinococcus radiodurans* in complex with 5-Hydroxy tryptophan.
Authors : Buddha, M.R.; Crane, B.R.
Deposited on : 2005-01-11
Resolution : 3.70 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
Mogul : 2022.3.0, CSD as543be (2022)
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

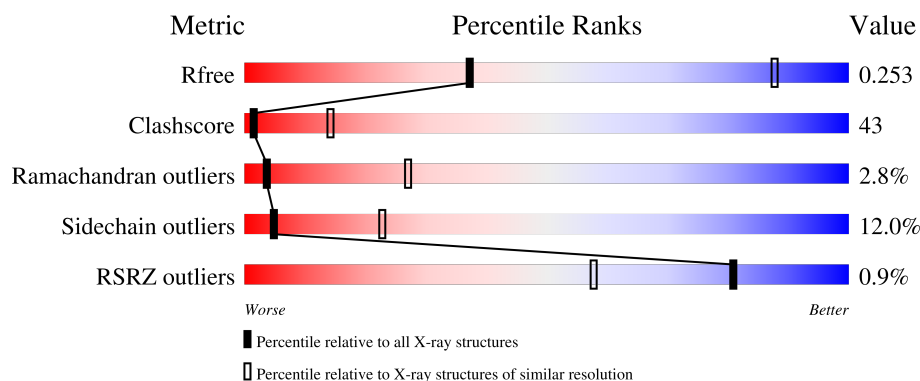
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.70 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	1131 (3.80-3.60)
Clashscore	190562	1171 (3.80-3.60)
Ramachandran outliers	187476	1129 (3.80-3.60)
Sidechain outliers	187428	1126 (3.80-3.60)
RSRZ outliers	180081	1130 (3.80-3.60)

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	351	2% 29% 51% 13% 6%
1	B	351	41% 43% 9% 6%
1	C	351	% 50% 35% 6% 6%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	HRP	B	1154	-	-	X	-

2 Entry composition [i](#)

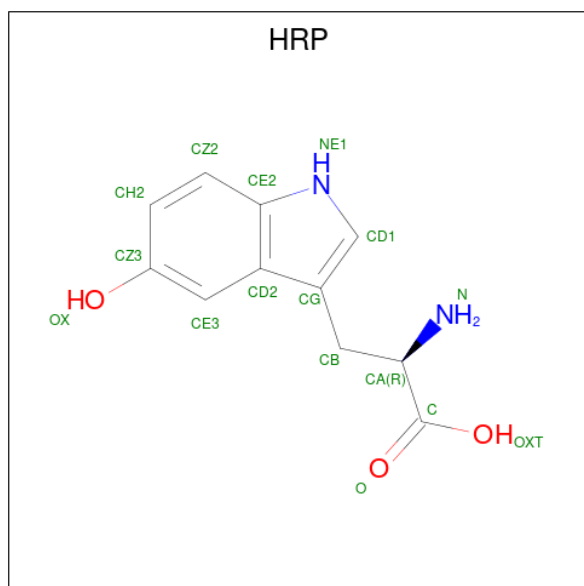
There are 2 unique types of molecules in this entry. The entry contains 7603 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 tryptophanyl-tRNA synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	331	Total	C	N	O	S	2	0	0
			2544	1599	468	471	6			
1	A	331	Total	C	N	O	S	0	0	0
			2511	1581	457	467	6			
1	C	331	Total	C	N	O	S	0	0	0
			2532	1593	462	471	6			

- Molecule 2 is 5-HYDROXY-L-TRYPTOPHAN (CCD ID: HRP) (formula: C₁₁H₁₂N₂O₃).

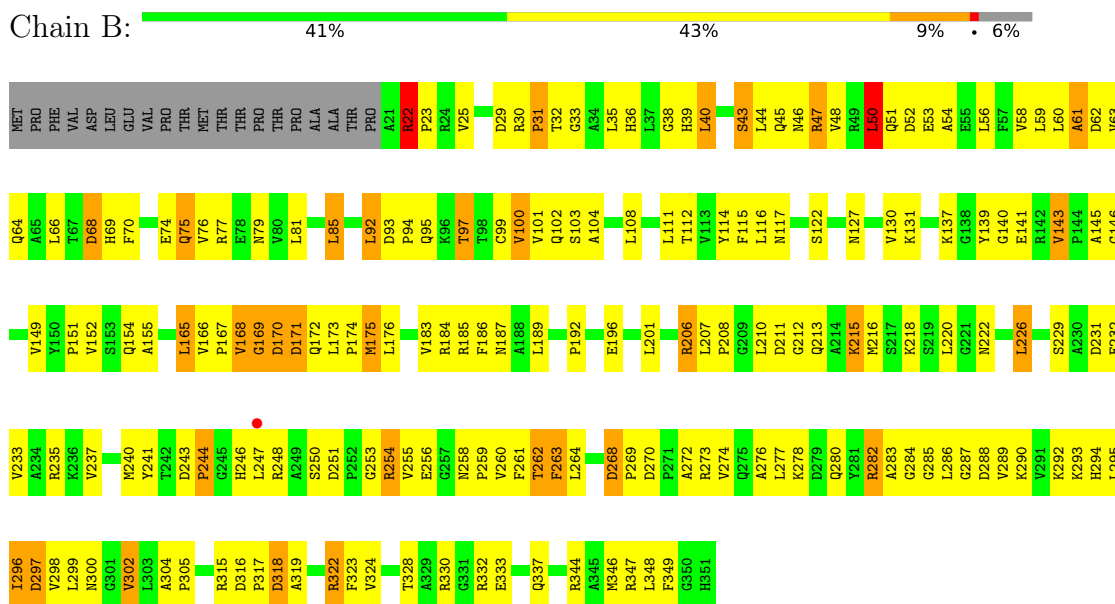


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			16	11	2	3		

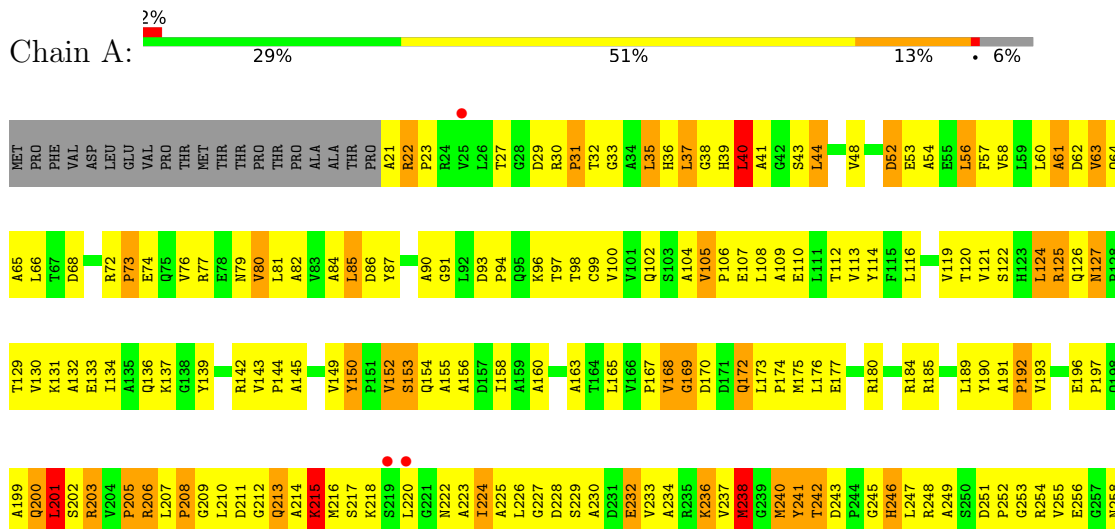
3 Residue-property plots

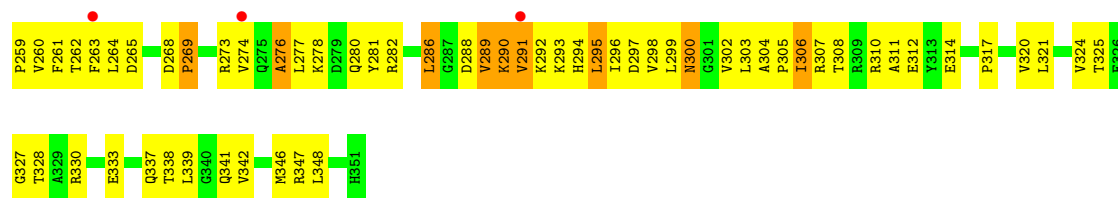
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: tryptophanyl-tRNA synthetase

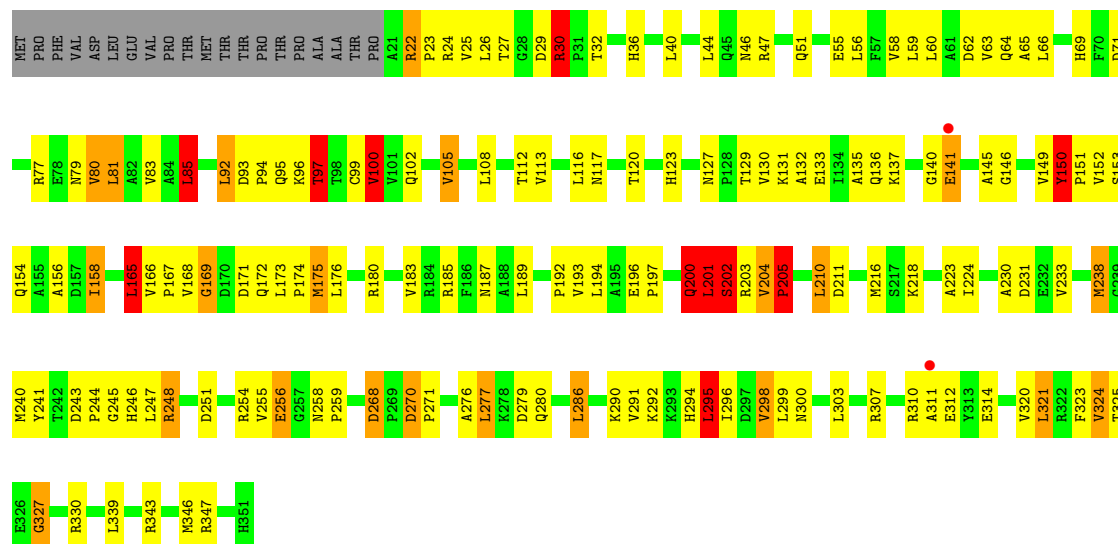


• Molecule 1: tryptophanyl-tRNA synthetase





• Molecule 1: tryptophanyl-tRNA synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	212.20Å 58.40Å 86.80Å 90.00° 96.50° 90.00°	Depositor
Resolution (Å)	30.00 – 3.70 30.00 – 3.73	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-3.70) 80.2 (30.00-3.73)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.42 (at 3.78Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.280 , 0.290 0.254 , 0.253	Depositor DCC
R_{free} test set	439 reflections (3.89%)	wwPDB-VP
Wilson B-factor (Å ²)	74.3	Xtriage
Anisotropy	0.748	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 66.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	7603	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.85% 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 ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: HRP

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	0.90	4/2558 (0.2%)	1.23	24/3478 (0.7%)
1	B	0.88	3/2591 (0.1%)	1.20	19/3519 (0.5%)
1	C	0.91	1/2579 (0.0%)	1.25	30/3505 (0.9%)
All	All	0.90	8/7728 (0.1%)	1.23	73/10502 (0.7%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	238	MET	SD-CE	10.13	2.04	1.79
1	A	291	VAL	CA-CB	8.55	1.63	1.54
1	A	238	MET	SD-CE	8.36	2.00	1.79
1	A	240	MET	SD-CE	6.64	1.96	1.79
1	A	342	VAL	CA-CB	-6.53	1.47	1.54
1	B	183	VAL	CA-CB	-6.08	1.47	1.54
1	B	324	VAL	CA-CB	-5.47	1.47	1.54
1	B	175	MET	SD-CE	5.04	1.92	1.79

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	215	LYS	N-CA-C	9.55	121.25	108.07
1	C	327	GLY	N-CA-C	-8.94	102.13	112.50
1	C	201	LEU	N-CA-C	8.70	122.51	110.24
1	B	268	ASP	CA-C-N	7.87	127.51	119.56
1	B	268	ASP	C-N-CA	7.87	127.51	119.56
1	A	43	SER	N-CA-C	7.76	119.74	111.28
1	B	272	ALA	N-CA-C	-7.53	103.01	111.07
1	B	332	ARG	N-CA-C	-7.47	103.14	111.28
1	A	105	VAL	CA-C-N	7.21	127.22	119.28
1	A	105	VAL	C-N-CA	7.21	127.22	119.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	263	PHE	N-CA-C	-7.14	104.20	113.12
1	A	243	ASP	CA-C-N	7.11	126.81	119.56
1	A	243	ASP	C-N-CA	7.11	126.81	119.56
1	A	127	ASN	CA-C-N	6.97	126.22	118.97
1	A	127	ASN	C-N-CA	6.97	126.22	118.97
1	A	44	LEU	N-CA-C	6.95	119.70	111.71
1	B	85	LEU	N-CA-C	-6.93	103.65	111.07
1	A	153	SER	N-CA-C	-6.90	104.73	113.01
1	B	100	VAL	N-CA-C	6.76	117.77	107.77
1	C	153	SER	N-CA-C	-6.46	104.32	111.82
1	B	244	PRO	N-CA-C	-6.44	102.20	111.22
1	B	171	ASP	N-CA-C	6.41	119.08	111.71
1	C	140	GLY	N-CA-C	-6.39	105.79	111.67
1	C	268	ASP	CA-C-N	6.34	126.09	119.05
1	C	268	ASP	C-N-CA	6.34	126.09	119.05
1	B	210	LEU	N-CA-C	-6.23	104.49	111.28
1	A	201	LEU	N-CA-C	6.23	119.74	110.52
1	B	155	ALA	N-CA-C	-6.22	104.58	111.36
1	B	212	GLY	N-CA-C	-6.15	106.14	115.00
1	C	295	LEU	N-CA-C	-6.15	104.50	111.14
1	A	80	VAL	N-CA-C	-6.14	103.84	110.23
1	A	276	ALA	N-CA-C	-6.09	103.38	111.24
1	C	158	ILE	N-CA-C	6.09	116.56	110.23
1	C	270	ASP	CA-C-N	6.08	125.78	119.82
1	C	270	ASP	C-N-CA	6.08	125.78	119.82
1	B	215	LYS	N-CA-C	6.03	118.77	109.07
1	C	150	TYR	CA-C-N	-6.02	112.66	119.28
1	C	150	TYR	C-N-CA	-6.02	112.66	119.28
1	C	100	VAL	N-CA-C	6.02	117.42	108.46
1	A	236	LYS	N-CA-C	-6.00	104.66	111.14
1	A	169	GLY	N-CA-C	-5.94	103.50	112.41
1	C	141	GLU	CA-C-O	5.85	125.22	118.97
1	C	324	VAL	N-CA-C	5.77	117.70	111.58
1	A	314	GLU	N-CA-C	-5.76	105.00	111.28
1	B	143	VAL	CB-CA-C	-5.62	105.72	110.94
1	C	200	GLN	N-CA-C	-5.59	97.09	107.71
1	C	149	VAL	CB-CA-C	-5.58	105.78	112.19
1	A	241	TYR	N-CA-C	5.55	117.76	107.99
1	A	100	VAL	N-CA-C	5.51	115.83	108.11
1	A	155	ALA	N-CA-C	-5.46	104.72	111.33
1	C	169	GLY	N-CA-C	-5.43	101.07	112.04
1	B	63	VAL	N-CA-C	5.34	116.11	110.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	85	LEU	N-CA-C	-5.33	105.47	111.28
1	A	40	LEU	N-CA-C	-5.33	105.12	111.03
1	A	168	VAL	N-CA-C	5.25	115.92	108.42
1	C	165	LEU	CA-C-N	-5.25	118.29	122.85
1	C	165	LEU	C-N-CA	-5.25	118.29	122.85
1	A	90	ALA	CA-C-N	-5.24	117.21	123.08
1	A	90	ALA	C-N-CA	-5.24	117.21	123.08
1	B	269	PRO	N-CA-C	5.21	120.37	113.86
1	C	202	SER	N-CA-C	5.20	121.88	110.80
1	C	30	ARG	CA-C-N	5.17	126.30	119.84
1	C	30	ARG	C-N-CA	5.17	126.30	119.84
1	C	175	MET	N-CA-C	-5.16	105.56	111.14
1	B	50	LEU	N-CA-C	-5.16	106.95	113.20
1	B	169	GLY	N-CA-C	-5.08	106.04	112.54
1	C	97	THR	N-CA-C	5.08	117.77	109.59
1	A	91	GLY	N-CA-C	5.07	120.51	114.48
1	C	218	LYS	N-CA-C	-5.07	105.65	111.07
1	C	303	LEU	N-CA-C	5.04	118.59	112.23
1	B	112	THR	N-CA-C	-5.03	105.71	111.14
1	C	105	VAL	CA-C-N	5.00	124.72	119.82
1	C	105	VAL	C-N-CA	5.00	124.72	119.82

There are no chirality outliers.

There are no planarity outliers.

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	2511	0	2507	284	0
1	B	2544	0	2568	205	0
1	C	2532	0	2546	172	0
2	B	16	0	9	15	0
All	All	7603	0	7630	654	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 43.

All (654) 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:238:MET:CE	1:C:238:MET:SD	2.04	1.43
2:B:1154:HRP:HE3	2:B:1154:HRP:N	1.52	1.24
1:B:51:GLN:HE21	1:B:97:THR:HG22	1.09	1.16
1:A:168:VAL:HG23	1:A:172:GLN:HB2	1.15	1.12
1:B:154:GLN:NE2	1:B:175:MET:HE2	1.68	1.08
1:A:35:LEU:HD11	1:A:86:ASP:HB3	1.38	1.03
2:B:1154:HRP:HE3	2:B:1154:HRP:H	1.10	1.02
1:B:47:ARG:HA	1:B:50:LEU:HG	1.39	1.02
1:B:280:GLN:HB2	1:B:286:LEU:HD12	1.44	1.00
1:C:204:VAL:HG12	1:C:205:PRO:HD2	1.43	1.00
1:A:226:LEU:HB3	1:A:306:ILE:HD11	1.43	0.99
1:C:51:GLN:HE22	1:C:93:ASP:H	1.07	0.99
1:B:154:GLN:HG2	2:B:1154:HRP:NE1	1.79	0.98
1:B:240:MET:HE3	1:B:260:VAL:HG12	1.48	0.96
2:B:1154:HRP:H	2:B:1154:HRP:CE3	1.78	0.96
1:C:120:THR:HG23	1:C:123:HIS:H	1.31	0.95
2:B:1154:HRP:N	2:B:1154:HRP:CE3	2.29	0.95
1:A:207:LEU:HD12	1:A:208:PRO:HD2	1.49	0.94
1:B:154:GLN:CG	2:B:1154:HRP:NE1	2.30	0.94
1:A:73:PRO:O	1:A:76:VAL:HG12	1.67	0.94
1:B:254:ARG:HD3	1:B:256:GLU:HG2	1.49	0.93
1:C:254:ARG:HD3	1:C:256:GLU:HG2	1.49	0.93
1:A:288:ASP:HA	1:A:291:VAL:HG12	1.48	0.91
1:B:51:GLN:NE2	1:B:97:THR:HG22	1.86	0.91
1:C:180:ARG:HG3	1:C:180:ARG:HH11	1.36	0.91
1:B:51:GLN:HE22	1:B:93:ASP:H	0.94	0.90
1:C:216:MET:HE3	1:C:223:ALA:HB1	1.53	0.89
1:B:51:GLN:HE22	1:B:93:ASP:N	1.72	0.88
1:B:295:LEU:O	1:B:298:VAL:HG12	1.72	0.88
1:B:154:GLN:HG2	2:B:1154:HRP:HE1	1.34	0.87
1:B:154:GLN:CD	1:B:175:MET:HE2	1.98	0.87
1:C:311:ALA:HA	1:C:314:GLU:HG3	1.56	0.87
1:A:207:LEU:HD23	1:A:216:MET:HE1	1.54	0.87
1:A:35:LEU:CD1	1:A:86:ASP:HB3	2.05	0.86
1:A:193:VAL:HG21	1:A:341:GLN:HB3	1.56	0.86
1:C:93:ASP:HB3	1:C:96:LYS:HB2	1.57	0.85
1:A:127:ASN:HB3	1:A:130:VAL:HG12	1.58	0.85
1:B:68:ASP:HB3	1:B:69:HIS:HD2	1.41	0.84
1:A:30:ARG:HH22	1:A:68:ASP:HB3	1.43	0.82
1:B:154:GLN:NE2	2:B:1154:HRP:CE2	2.43	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:92:LEU:HG	1:C:97:THR:HG21	1.62	0.82
1:B:154:GLN:HE21	2:B:1154:HRP:CE2	1.93	0.82
1:B:168:VAL:HG23	1:B:172:GLN:HB2	1.63	0.81
1:C:51:GLN:HE21	1:C:97:THR:HG22	1.44	0.81
1:A:207:LEU:CD1	1:A:208:PRO:HD2	2.09	0.81
1:A:254:ARG:HB2	1:A:256:GLU:OE1	1.82	0.80
1:C:77:ARG:O	1:C:80:VAL:HG12	1.82	0.80
1:A:303:LEU:O	1:A:307:ARG:HB2	1.81	0.80
1:A:60:LEU:O	1:A:62:ASP:N	2.15	0.79
1:B:154:GLN:NE2	1:B:175:MET:CE	2.45	0.79
1:A:85:LEU:HD13	1:A:310:ARG:NH2	1.98	0.78
1:A:168:VAL:CG2	1:A:172:GLN:HB2	2.05	0.78
1:A:218:LYS:HA	1:A:218:LYS:HE2	1.63	0.78
1:C:233:VAL:HG21	1:C:307:ARG:HH21	1.47	0.78
1:A:27:THR:O	1:A:58:VAL:HA	1.82	0.78
1:A:127:ASN:HB3	1:A:130:VAL:CG1	2.13	0.78
1:A:288:ASP:CA	1:A:291:VAL:HG12	2.14	0.78
1:B:146:GLY:HA3	1:C:117:ASN:ND2	1.99	0.78
1:B:68:ASP:OD1	1:B:137:LYS:HE2	1.82	0.78
1:B:36:HIS:HD2	1:B:38:GLY:H	1.33	0.77
1:A:273:ARG:O	1:A:276:ALA:HB3	1.85	0.77
1:A:288:ASP:HA	1:A:291:VAL:CG1	2.14	0.77
1:B:316:ASP:OD2	1:B:319:ALA:HB2	1.85	0.76
1:A:254:ARG:HH11	1:A:256:GLU:HB2	1.48	0.76
1:B:44:LEU:O	1:B:48:VAL:HG23	1.85	0.76
1:B:50:LEU:HD11	1:B:56:LEU:HD13	1.68	0.75
1:B:277:LEU:HD12	1:B:286:LEU:HD11	1.68	0.75
1:A:168:VAL:HG22	1:A:169:GLY:O	1.86	0.75
1:A:40:LEU:HA	1:A:44:LEU:HB3	1.68	0.75
1:A:260:VAL:HG13	1:A:295:LEU:CD2	2.16	0.75
1:A:269:PRO:HD2	1:A:273:ARG:NH1	2.02	0.74
1:A:169:GLY:H	1:A:172:GLN:HG3	1.53	0.74
1:B:51:GLN:HE21	1:B:97:THR:CG2	1.95	0.74
1:A:173:LEU:N	1:A:174:PRO:HD2	2.03	0.74
1:C:290:LYS:HE2	1:C:290:LYS:HA	1.70	0.74
1:A:260:VAL:HG13	1:A:295:LEU:HD22	1.69	0.73
1:C:201:LEU:HD23	1:C:202:SER:H	1.52	0.73
1:C:22:ARG:HB2	1:C:23:PRO:CD	2.17	0.73
1:A:168:VAL:HG11	1:A:199:ALA:HB1	1.70	0.73
1:B:68:ASP:HB3	1:B:69:HIS:CD2	2.24	0.73
1:C:254:ARG:CD	1:C:256:GLU:HG2	2.18	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:56:LEU:HD12	1:C:97:THR:HB	1.72	0.72
1:B:349:PHE:CD2	1:C:80:VAL:HG11	2.24	0.72
1:B:154:GLN:HG3	2:B:1154:HRP:CD1	2.21	0.71
1:A:125:ARG:HG3	1:A:125:ARG:HH21	1.53	0.71
1:A:259:PRO:HA	1:A:262:THR:CG2	2.20	0.71
1:C:51:GLN:NE2	1:C:97:THR:HG22	2.06	0.70
1:B:237:VAL:HG13	1:B:240:MET:HE1	1.73	0.70
1:C:40:LEU:HA	1:C:44:LEU:HD12	1.74	0.70
1:A:112:THR:HG22	1:A:153:SER:HA	1.72	0.70
1:C:26:LEU:HD23	1:C:158:ILE:HD13	1.74	0.69
1:C:175:MET:SD	1:C:175:MET:C	2.75	0.69
1:B:273:ARG:NH2	1:B:294:HIS:NE2	2.40	0.69
1:B:51:GLN:NE2	1:B:93:ASP:H	1.79	0.69
1:A:52:ASP:HB3	1:A:96:LYS:NZ	2.08	0.68
1:B:207:LEU:HD12	1:B:208:PRO:HD2	1.75	0.68
1:A:112:THR:O	1:A:116:LEU:HG	1.92	0.68
1:C:22:ARG:HB2	1:C:23:PRO:HD3	1.75	0.68
1:B:255:VAL:HG21	1:B:278:LYS:HG3	1.76	0.68
1:A:248:ARG:HD3	1:A:248:ARG:N	2.09	0.68
1:A:110:GLU:O	1:A:113:VAL:HG12	1.93	0.68
1:A:208:PRO:HG3	1:A:262:THR:HG23	1.76	0.67
1:B:33:GLY:HA2	1:B:79:ASN:OD1	1.94	0.67
1:A:98:THR:CG2	1:A:330:ARG:HD2	2.25	0.67
1:B:22:ARG:CG	1:B:23:PRO:HD3	2.25	0.67
1:A:333:GLU:O	1:A:337:GLN:HG2	1.94	0.67
1:A:184:ARG:HD2	1:A:196:GLU:OE1	1.93	0.67
1:A:216:MET:HG3	1:A:223:ALA:HA	1.76	0.67
1:A:139:TYR:CD2	1:A:143:VAL:HG13	2.30	0.67
1:A:296:ILE:HG13	1:A:297:ASP:N	2.10	0.67
1:C:180:ARG:HG3	1:C:180:ARG:NH1	2.08	0.67
1:C:166:VAL:HG21	1:C:176:LEU:HD23	1.77	0.67
1:C:216:MET:HE3	1:C:223:ALA:CB	2.24	0.67
1:C:258:ASN:C	1:C:258:ASN:HD22	2.02	0.67
1:B:40:LEU:HD23	1:B:226:LEU:HD21	1.76	0.67
1:A:180:ARG:HG2	1:A:196:GLU:HG2	1.77	0.67
1:C:127:ASN:OD1	1:C:129:THR:HG22	1.95	0.67
1:B:154:GLN:CG	2:B:1154:HRP:CD1	2.72	0.66
1:B:259:PRO:HA	1:B:262:THR:CG2	2.26	0.66
1:A:85:LEU:HD13	1:A:310:ARG:HH22	1.57	0.66
1:A:139:TYR:O	1:A:142:ARG:HG3	1.96	0.66
1:A:226:LEU:HD23	1:A:306:ILE:CD1	2.25	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:SER:OG	1:A:232:GLU:HB2	1.95	0.66
1:B:175:MET:SD	1:B:175:MET:C	2.78	0.66
1:A:98:THR:OG1	1:A:330:ARG:NH1	2.29	0.66
1:A:305:PRO:HG2	1:A:306:ILE:H	1.60	0.65
1:A:317:PRO:O	1:A:320:VAL:HG12	1.96	0.65
1:B:146:GLY:HA3	1:C:117:ASN:HD22	1.60	0.65
1:B:22:ARG:HH21	1:B:53:GLU:HG3	1.60	0.65
1:A:202:SER:OG	1:A:205:PRO:HB3	1.97	0.65
1:C:171:ASP:HB2	1:C:172:GLN:NE2	2.10	0.65
1:B:22:ARG:CB	1:B:23:PRO:HD3	2.26	0.65
1:B:145:ALA:HB3	1:C:116:LEU:O	1.95	0.65
1:A:160:ALA:HA	1:A:338:THR:HG21	1.78	0.65
1:A:226:LEU:HD12	1:A:226:LEU:H	1.60	0.65
1:A:63:VAL:HG13	1:A:102:GLN:HG2	1.78	0.65
1:B:154:GLN:HE21	2:B:1154:HRP:CZ2	2.09	0.65
1:B:268:ASP:OD2	1:B:274:VAL:HG23	1.95	0.65
1:C:56:LEU:HD11	1:C:58:VAL:HG13	1.78	0.65
1:A:36:HIS:H	1:A:39:HIS:HD2	1.44	0.65
1:A:242:THR:HB	1:A:258:ASN:HD21	1.62	0.64
1:A:304:ALA:N	1:A:305:PRO:HD2	2.13	0.64
1:B:36:HIS:CE1	1:B:39:HIS:CE1	2.86	0.64
1:A:175:MET:SD	1:A:176:LEU:N	2.71	0.64
1:A:254:ARG:HD2	1:A:256:GLU:HB2	1.80	0.64
1:C:180:ARG:HD3	1:C:197:PRO:O	1.97	0.64
1:A:175:MET:SD	1:A:175:MET:C	2.81	0.64
1:A:230:ALA:HB1	1:A:300:ASN:HD21	1.61	0.64
1:A:294:HIS:HA	1:A:297:ASP:OD2	1.98	0.64
1:B:40:LEU:HA	1:B:44:LEU:HB2	1.80	0.63
1:A:265:ASP:HB3	1:A:274:VAL:HG11	1.81	0.63
1:C:183:VAL:HG13	1:C:194:LEU:HB2	1.80	0.63
1:A:30:ARG:O	1:A:32:THR:N	2.32	0.63
1:B:43:SER:O	1:B:46:ASN:N	2.32	0.63
1:A:277:LEU:HD12	1:A:286:LEU:HD11	1.81	0.63
1:B:254:ARG:CD	1:B:256:GLU:HG2	2.26	0.63
1:A:30:ARG:HH22	1:A:68:ASP:CB	2.12	0.63
1:C:166:VAL:HG23	1:C:166:VAL:O	1.99	0.63
1:C:204:VAL:O	1:C:205:PRO:O	2.17	0.63
1:B:322:ARG:HG2	1:B:322:ARG:HH21	1.64	0.62
1:C:171:ASP:HB2	1:C:172:GLN:HE22	1.62	0.62
1:A:33:GLY:HA2	1:A:79:ASN:CG	2.24	0.62
1:C:59:LEU:HD12	1:C:100:VAL:HG22	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:254:ARG:CG	1:C:256:GLU:HG2	2.29	0.62
1:B:184:ARG:HG2	1:B:184:ARG:HH21	1.64	0.62
1:B:173:LEU:N	1:B:174:PRO:HD2	2.14	0.62
1:A:232:GLU:O	1:A:236:LYS:HB2	1.99	0.62
1:A:176:LEU:O	1:A:176:LEU:HD13	1.99	0.62
1:A:255:VAL:HG22	1:A:261:PHE:CE1	2.33	0.62
1:A:133:GLU:O	1:A:137:LYS:HG3	2.00	0.62
1:A:208:PRO:HG2	1:A:263:PHE:CE2	2.35	0.62
1:A:229:SER:O	1:A:233:VAL:HG23	2.00	0.61
1:A:105:VAL:HG13	1:A:105:VAL:O	2.00	0.61
1:A:220:LEU:C	1:A:222:ASN:H	2.08	0.61
1:B:255:VAL:HG22	1:B:261:PHE:CE2	2.34	0.61
1:A:158:ILE:HG23	1:A:163:ALA:HB3	1.82	0.61
1:A:293:LYS:O	1:A:296:ILE:HG12	2.00	0.61
1:A:33:GLY:HA2	1:A:79:ASN:ND2	2.15	0.61
1:A:273:ARG:O	1:A:277:LEU:HD23	2.00	0.61
1:A:288:ASP:C	1:A:291:VAL:HG12	2.25	0.61
1:C:51:GLN:HE22	1:C:93:ASP:N	1.89	0.61
1:A:79:ASN:HA	1:A:82:ALA:HB3	1.82	0.61
1:B:247:LEU:HB2	1:B:248:ARG:HH21	1.66	0.60
1:A:288:ASP:O	1:A:291:VAL:HG12	2.00	0.60
1:A:125:ARG:HG3	1:A:125:ARG:NH2	2.17	0.60
1:C:230:ALA:HB1	1:C:300:ASN:HD21	1.65	0.60
1:C:254:ARG:HD3	1:C:256:GLU:CG	2.27	0.60
1:A:259:PRO:HA	1:A:262:THR:HG22	1.82	0.60
1:C:51:GLN:NE2	1:C:93:ASP:H	1.89	0.60
1:B:255:VAL:HG13	1:B:261:PHE:CD2	2.36	0.60
1:A:149:VAL:O	1:A:149:VAL:HG12	2.01	0.60
1:A:241:TYR:O	1:A:259:PRO:HD2	2.00	0.60
1:A:254:ARG:HH11	1:A:256:GLU:CB	2.13	0.60
1:A:260:VAL:O	1:A:264:LEU:HB2	2.00	0.60
1:A:201:LEU:CD2	1:A:202:SER:H	2.14	0.60
1:A:201:LEU:HD23	1:A:202:SER:H	1.66	0.60
1:B:36:HIS:CD2	1:B:38:GLY:H	2.17	0.60
1:C:36:HIS:CG	1:C:216:MET:HE2	2.38	0.59
1:B:296:ILE:HD12	1:B:300:ASN:ND2	2.17	0.59
1:C:108:LEU:O	1:C:112:THR:HG23	2.02	0.59
1:A:56:LEU:HD22	1:A:57:PHE:N	2.17	0.59
1:C:216:MET:CE	1:C:224:ILE:H	2.14	0.59
1:A:142:ARG:O	1:A:142:ARG:HD3	2.02	0.59
1:B:66:LEU:C	1:B:68:ASP:H	2.09	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:294:HIS:O	1:C:298:VAL:HG12	2.03	0.59
1:A:168:VAL:CG1	1:A:199:ALA:HB1	2.33	0.59
1:A:261:PHE:CE1	1:A:278:LYS:HG2	2.37	0.59
1:C:216:MET:HE3	1:C:224:ILE:H	1.68	0.59
1:B:248:ARG:H	1:B:248:ARG:HD3	1.68	0.59
1:A:32:THR:HG22	1:A:32:THR:O	2.03	0.59
1:A:173:LEU:O	1:A:177:GLU:HB2	2.02	0.59
1:B:293:LYS:O	1:B:296:ILE:HG23	2.02	0.58
1:A:321:LEU:O	1:A:325:THR:HG23	2.02	0.58
1:C:311:ALA:CA	1:C:314:GLU:HG3	2.31	0.58
1:A:290:LYS:HE2	1:A:290:LYS:N	2.18	0.58
1:C:40:LEU:HA	1:C:44:LEU:HB2	1.85	0.58
1:C:36:HIS:ND1	1:C:216:MET:HE2	2.17	0.58
1:A:76:VAL:HG13	1:A:77:ARG:N	2.19	0.58
1:A:253:GLY:O	1:A:282:ARG:HA	2.02	0.58
1:C:56:LEU:CD1	1:C:58:VAL:HG13	2.33	0.58
1:B:254:ARG:HH11	1:B:256:GLU:CG	2.17	0.58
1:B:51:GLN:NE2	1:B:97:THR:CG2	2.59	0.58
1:C:36:HIS:HB2	1:C:216:MET:CE	2.34	0.58
1:C:65:ALA:O	1:C:69:HIS:HB2	2.04	0.58
1:B:248:ARG:HD3	1:B:248:ARG:N	2.18	0.58
1:A:220:LEU:C	1:A:222:ASN:N	2.61	0.57
1:B:22:ARG:HH21	1:B:53:GLU:CG	2.17	0.57
1:A:44:LEU:O	1:A:48:VAL:HG23	2.04	0.57
1:A:259:PRO:O	1:A:263:PHE:HB2	2.03	0.57
1:C:168:VAL:CG2	1:C:172:GLN:HB2	2.33	0.57
1:A:116:LEU:HD23	1:A:152:VAL:HG21	1.86	0.57
1:A:180:ARG:O	1:A:184:ARG:HD3	2.05	0.57
1:C:95:GLN:HA	1:C:95:GLN:NE2	2.20	0.57
1:B:296:ILE:CD1	1:B:300:ASN:ND2	2.68	0.57
1:B:60:LEU:O	1:B:62:ASP:N	2.36	0.56
1:A:295:LEU:O	1:A:299:LEU:HG	2.05	0.56
1:A:225:ALA:N	1:A:228:ASP:OD2	2.37	0.56
1:C:310:ARG:O	1:C:314:GLU:HG2	2.05	0.56
1:B:101:VAL:HG13	1:B:101:VAL:O	2.04	0.56
1:B:186:PHE:CD1	1:B:186:PHE:C	2.83	0.56
1:C:60:LEU:HD22	1:C:83:VAL:HG21	1.87	0.56
1:B:229:SER:O	1:B:233:VAL:HG12	2.06	0.56
1:B:296:ILE:HG13	1:B:297:ASP:N	2.20	0.56
1:C:112:THR:HG22	1:C:156:ALA:CB	2.35	0.56
1:A:136:GLN:O	1:A:136:GLN:HG3	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:VAL:O	1:A:152:VAL:HG13	2.06	0.56
1:A:216:MET:CG	1:A:223:ALA:HA	2.36	0.56
1:C:64:GLN:H	1:C:102:GLN:HE22	1.54	0.56
1:B:53:GLU:OE1	1:B:53:GLU:HA	2.06	0.55
1:A:130:VAL:O	1:A:134:ILE:HG13	2.05	0.55
1:C:201:LEU:HD23	1:C:202:SER:N	2.21	0.55
1:B:137:LYS:HD3	1:B:139:TYR:CE2	2.42	0.55
1:B:206:ARG:NH1	1:B:215:LYS:HE2	2.21	0.55
1:C:166:VAL:HG21	1:C:176:LEU:CD2	2.36	0.55
1:A:277:LEU:HA	1:A:280:GLN:NE2	2.22	0.55
1:C:36:HIS:HB2	1:C:216:MET:HE2	1.89	0.55
1:B:22:ARG:HB2	1:B:23:PRO:HD3	1.88	0.55
1:C:339:LEU:O	1:C:343:ARG:HG3	2.06	0.55
1:A:41:ALA:CB	1:A:207:LEU:HD22	2.37	0.55
1:A:185:ARG:HD2	1:A:189:LEU:HD22	1.89	0.55
1:C:51:GLN:NE2	1:C:97:THR:CG2	2.69	0.55
1:A:58:VAL:HG13	1:A:58:VAL:O	2.05	0.55
1:A:211:ASP:C	1:A:213:GLN:H	2.15	0.54
1:C:233:VAL:HG21	1:C:307:ARG:NH2	2.20	0.54
1:B:233:VAL:O	1:B:237:VAL:HG23	2.07	0.54
1:A:41:ALA:HB2	1:A:207:LEU:HD22	1.89	0.54
1:A:110:GLU:HG2	1:A:114:TYR:CE1	2.42	0.54
1:A:190:TYR:O	1:A:191:ALA:C	2.49	0.54
1:C:286:LEU:HD12	1:C:291:VAL:HG23	1.89	0.54
1:B:237:VAL:HG13	1:B:240:MET:CE	2.37	0.54
1:B:270:ASP:O	1:B:273:ARG:HB3	2.08	0.54
1:A:193:VAL:HG21	1:A:341:GLN:CB	2.31	0.54
1:A:205:PRO:O	1:A:206:ARG:C	2.51	0.54
1:A:232:GLU:HA	1:A:232:GLU:OE1	2.07	0.54
1:C:254:ARG:HD2	1:C:256:GLU:O	2.08	0.54
1:A:23:PRO:O	1:A:54:ALA:HB1	2.08	0.54
1:A:56:LEU:HD13	1:A:97:THR:HG23	1.89	0.54
1:A:107:GLU:HB3	1:A:339:LEU:HG	1.89	0.54
1:A:207:LEU:CG	1:A:208:PRO:HD2	2.37	0.54
1:A:237:VAL:O	1:A:240:MET:N	2.34	0.54
1:B:50:LEU:C	1:B:50:LEU:HD12	2.33	0.54
1:B:277:LEU:CD1	1:B:286:LEU:HD11	2.38	0.53
1:B:330:ARG:O	1:B:333:GLU:HB3	2.08	0.53
1:C:127:ASN:OD1	1:C:129:THR:CG2	2.56	0.53
1:C:180:ARG:HE	1:C:196:GLU:CD	2.15	0.53
1:B:99:CYS:HB2	1:B:323:PHE:CZ	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:25:VAL:HG22	1:B:165:LEU:HB3	1.90	0.53
1:C:40:LEU:CA	1:C:44:LEU:HD12	2.38	0.53
1:B:168:VAL:HG22	1:B:169:GLY:O	2.08	0.53
1:C:180:ARG:HD2	1:C:196:GLU:HG3	1.91	0.53
1:A:98:THR:OG1	1:A:330:ARG:HD2	2.09	0.53
1:C:295:LEU:CD2	1:C:299:LEU:HG	2.38	0.53
1:C:346:MET:O	1:C:347:ARG:HB2	2.08	0.53
1:A:52:ASP:HB3	1:A:96:LYS:HZ3	1.74	0.53
1:C:105:VAL:O	1:C:105:VAL:HG13	2.09	0.53
1:B:117:ASN:ND2	1:C:146:GLY:HA3	2.24	0.52
1:B:263:PHE:HB3	1:B:295:LEU:HD11	1.91	0.52
1:C:256:GLU:H	1:C:256:GLU:CD	2.16	0.52
1:B:66:LEU:C	1:B:68:ASP:N	2.65	0.52
1:B:43:SER:C	1:B:47:ARG:HD2	2.35	0.52
1:B:64:GLN:H	1:B:102:GLN:HE22	1.58	0.52
1:C:211:ASP:OD1	1:C:211:ASP:O	2.27	0.52
1:B:36:HIS:CD2	1:B:216:MET:SD	3.02	0.52
1:B:243:ASP:OD1	1:B:244:PRO:O	2.28	0.52
1:A:173:LEU:N	1:A:174:PRO:CD	2.73	0.52
1:A:22:ARG:N	1:A:23:PRO:CD	2.73	0.52
1:A:35:LEU:HD12	1:A:226:LEU:HD13	1.91	0.52
1:A:260:VAL:HG13	1:A:295:LEU:HD23	1.92	0.52
1:A:260:VAL:HA	1:A:295:LEU:HD22	1.91	0.52
1:A:261:PHE:HZ	1:A:281:TYR:CG	2.27	0.52
1:A:290:LYS:HE2	1:A:290:LYS:CA	2.40	0.52
1:B:231:ASP:O	1:B:235:ARG:HG3	2.10	0.52
1:A:170:ASP:OD1	1:A:203:ARG:NH2	2.43	0.52
1:A:237:VAL:HA	1:A:240:MET:HG3	1.91	0.52
1:A:112:THR:CG2	1:A:153:SER:HA	2.38	0.52
1:A:289:VAL:O	1:A:293:LYS:HB2	2.10	0.52
1:C:40:LEU:HD12	1:C:44:LEU:HB2	1.91	0.52
1:C:60:LEU:HD22	1:C:83:VAL:CG2	2.40	0.52
1:C:295:LEU:HD22	1:C:299:LEU:HG	1.92	0.52
1:B:259:PRO:HA	1:B:262:THR:HG23	1.92	0.51
1:A:226:LEU:HD23	1:A:306:ILE:HD12	1.92	0.51
1:C:29:ASP:OD2	1:C:47:ARG:NH2	2.41	0.51
1:B:254:ARG:HD3	1:B:256:GLU:CG	2.30	0.51
1:A:276:ALA:O	1:A:280:GLN:HG3	2.11	0.51
1:A:302:VAL:HG13	1:A:303:LEU:HD13	1.93	0.51
1:C:105:VAL:O	1:C:105:VAL:HG22	2.10	0.51
1:C:258:ASN:C	1:C:258:ASN:ND2	2.69	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:VAL:HG13	1:A:131:LYS:N	2.24	0.51
1:A:208:PRO:HG2	1:A:263:PHE:HE2	1.75	0.51
1:A:212:GLY:O	1:A:213:GLN:C	2.53	0.51
1:A:298:VAL:O	1:A:302:VAL:HG12	2.10	0.51
1:B:154:GLN:HE22	1:B:175:MET:HE2	1.68	0.51
1:B:35:LEU:HD13	1:B:44:LEU:HD11	1.92	0.51
1:B:254:ARG:NH1	1:B:256:GLU:HG3	2.25	0.51
1:B:289:VAL:O	1:B:293:LYS:HG3	2.10	0.51
1:A:209:GLY:O	1:A:210:LEU:C	2.53	0.51
1:B:333:GLU:HG3	1:B:337:GLN:HE21	1.76	0.51
1:A:237:VAL:HG13	1:A:295:LEU:HD23	1.93	0.51
1:B:232:GLU:O	1:B:233:VAL:C	2.55	0.50
1:B:292:LYS:O	1:B:296:ILE:CG2	2.60	0.50
1:A:112:THR:HG22	1:A:156:ALA:CB	2.41	0.50
1:A:130:VAL:CG1	1:A:131:LYS:N	2.73	0.50
1:A:193:VAL:HG22	1:A:193:VAL:O	2.12	0.50
1:C:56:LEU:HD11	1:C:58:VAL:CG1	2.40	0.50
1:C:171:ASP:CB	1:C:172:GLN:NE2	2.73	0.50
1:C:210:LEU:HD21	1:C:224:ILE:HG13	1.93	0.50
1:A:63:VAL:HG13	1:A:102:GLN:CG	2.41	0.50
1:A:108:LEU:O	1:A:112:THR:HG23	2.11	0.50
1:A:215:LYS:NZ	1:A:217:SER:OG	2.44	0.50
1:C:203:ARG:C	1:C:204:VAL:HG23	2.37	0.50
1:B:211:ASP:OD1	1:B:222:ASN:HB3	2.11	0.50
1:C:58:VAL:O	1:C:58:VAL:HG23	2.10	0.50
1:C:276:ALA:O	1:C:280:GLN:HG3	2.11	0.50
1:B:58:VAL:HG23	1:B:99:CYS:HA	1.93	0.50
1:C:80:VAL:HG13	1:C:81:LEU:N	2.27	0.50
1:A:68:ASP:OD2	1:A:137:LYS:NZ	2.45	0.50
1:A:99:CYS:SG	1:A:324:VAL:HG12	2.52	0.50
1:B:295:LEU:HA	1:B:298:VAL:HG12	1.94	0.50
1:B:101:VAL:HG12	1:B:328:THR:OG1	2.12	0.50
1:A:305:PRO:HG2	1:A:306:ILE:HG22	1.93	0.50
1:B:40:LEU:O	1:B:45:GLN:HG2	2.12	0.49
1:A:256:GLU:H	1:A:256:GLU:CD	2.20	0.49
1:C:56:LEU:HD13	1:C:56:LEU:C	2.37	0.49
1:A:248:ARG:HD3	1:A:248:ARG:H	1.78	0.49
1:C:295:LEU:HD22	1:C:299:LEU:CD1	2.42	0.49
1:B:140:GLY:O	1:B:143:VAL:HG23	2.13	0.49
1:B:298:VAL:O	1:B:302:VAL:HG12	2.13	0.49
1:A:265:ASP:CB	1:A:274:VAL:HG11	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:295:LEU:HA	1:C:298:VAL:CG1	2.43	0.49
1:B:102:GLN:HG3	1:B:108:LEU:HD12	1.95	0.49
1:A:241:TYR:O	1:A:259:PRO:CD	2.60	0.49
1:A:260:VAL:CG1	1:A:295:LEU:HD22	2.40	0.49
1:A:66:LEU:HD21	1:A:76:VAL:HG11	1.93	0.49
1:A:98:THR:HG23	1:A:330:ARG:HD2	1.94	0.49
1:A:293:LYS:O	1:A:296:ILE:CG1	2.60	0.49
1:C:241:TYR:OH	1:C:244:PRO:HD3	2.13	0.49
1:B:213:GLN:OE1	1:B:220:LEU:HD13	2.13	0.49
1:C:310:ARG:O	1:C:314:GLU:CG	2.61	0.49
1:A:274:VAL:HG12	1:A:278:LYS:HE3	1.95	0.48
1:A:30:ARG:HD2	1:A:65:ALA:HA	1.94	0.48
1:A:76:VAL:CG1	1:A:77:ARG:N	2.77	0.48
1:B:278:LYS:HG2	1:B:282:ARG:NH2	2.28	0.48
1:A:255:VAL:HG21	1:A:278:LYS:HD3	1.93	0.48
1:B:30:ARG:HH22	1:B:68:ASP:HB2	1.77	0.48
1:B:74:GLU:OE1	1:C:347:ARG:NH1	2.46	0.48
1:B:244:PRO:C	1:B:246:HIS:H	2.21	0.48
1:B:295:LEU:C	1:B:295:LEU:HD23	2.38	0.48
1:A:274:VAL:C	1:A:276:ALA:H	2.19	0.48
1:A:63:VAL:HG13	1:A:102:GLN:CD	2.37	0.48
1:A:288:ASP:O	1:A:292:LYS:HG3	2.14	0.48
1:C:120:THR:HG22	1:C:123:HIS:HB2	1.95	0.48
1:A:277:LEU:HD13	1:A:280:GLN:NE2	2.29	0.48
1:B:264:LEU:HG	1:B:295:LEU:HD12	1.95	0.48
1:C:204:VAL:O	1:C:205:PRO:C	2.56	0.48
1:C:295:LEU:O	1:C:298:VAL:HG13	2.12	0.48
1:B:315:ARG:O	1:B:315:ARG:HG2	2.14	0.48
1:A:37:LEU:HD21	1:A:303:LEU:HD11	1.95	0.48
1:B:101:VAL:HG13	1:B:104:ALA:HB3	1.96	0.48
1:B:273:ARG:O	1:B:276:ALA:HB3	2.13	0.48
1:A:214:ALA:C	1:A:215:LYS:HG3	2.38	0.48
1:C:62:ASP:O	1:C:66:LEU:HG	2.13	0.48
1:B:241:TYR:O	1:B:259:PRO:HG2	2.14	0.47
1:A:37:LEU:O	1:A:41:ALA:N	2.28	0.47
1:B:154:GLN:HE22	1:B:175:MET:CE	2.23	0.47
1:A:261:PHE:CD1	1:A:278:LYS:HG2	2.48	0.47
1:A:269:PRO:HD2	1:A:273:ARG:HH12	1.74	0.47
1:C:201:LEU:HD22	1:C:202:SER:HB3	1.96	0.47
1:A:21:ALA:HB3	1:A:53:GLU:O	2.14	0.47
1:A:248:ARG:NE	1:A:251:ASP:OD1	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:294:HIS:O	1:A:296:ILE:N	2.48	0.47
1:C:127:ASN:HB3	1:C:130:VAL:HG22	1.97	0.47
1:B:22:ARG:HE	1:B:53:GLU:HG3	1.79	0.47
1:B:43:SER:O	1:B:47:ARG:HD2	2.14	0.47
1:B:293:LYS:C	1:B:296:ILE:HG23	2.40	0.47
1:A:253:GLY:HA3	1:A:281:TYR:CE2	2.50	0.47
1:B:22:ARG:NH2	1:B:53:GLU:HG3	2.28	0.47
1:B:62:ASP:OD1	1:B:103:SER:OG	2.33	0.47
1:B:296:ILE:CD1	1:B:300:ASN:HD21	2.28	0.47
1:A:237:VAL:O	1:A:240:MET:HB2	2.15	0.47
1:C:154:GLN:O	1:C:158:ILE:HG12	2.15	0.47
1:C:321:LEU:HD22	1:C:325:THR:HG23	1.97	0.47
1:B:22:ARG:HG3	1:B:23:PRO:HD3	1.95	0.47
1:B:130:VAL:HG13	1:B:131:LYS:N	2.30	0.47
1:B:260:VAL:HG23	1:B:261:PHE:N	2.28	0.47
1:C:58:VAL:CG2	1:C:99:CYS:HA	2.45	0.47
1:B:29:ASP:O	1:B:31:PRO:HD3	2.15	0.47
1:A:105:VAL:HG13	1:A:108:LEU:HG	1.96	0.47
1:C:254:ARG:HG2	1:C:255:VAL:N	2.29	0.47
1:B:278:LYS:HB2	1:B:278:LYS:HE3	1.71	0.46
1:A:37:LEU:HG	1:A:224:ILE:HG12	1.97	0.46
1:A:110:GLU:O	1:A:113:VAL:CG1	2.61	0.46
1:B:127:ASN:HB3	1:B:130:VAL:HG12	1.98	0.46
1:A:36:HIS:HD2	1:A:38:GLY:H	1.62	0.46
1:C:22:ARG:CB	1:C:23:PRO:CD	2.91	0.46
1:A:52:ASP:OD2	1:A:52:ASP:N	2.36	0.46
1:C:210:LEU:HD12	1:C:240:MET:HG2	1.97	0.46
1:B:237:VAL:O	1:B:240:MET:HE2	2.15	0.46
1:B:254:ARG:NH1	1:B:256:GLU:CG	2.79	0.46
1:A:40:LEU:HA	1:A:44:LEU:CB	2.42	0.46
1:A:208:PRO:HG3	1:A:262:THR:CG2	2.43	0.46
1:A:62:ASP:H	1:A:102:GLN:HB3	1.81	0.46
1:A:127:ASN:HB3	1:A:130:VAL:HG11	1.96	0.46
1:A:299:LEU:O	1:A:303:LEU:HD13	2.16	0.46
1:B:253:GLY:O	1:B:282:ARG:HA	2.15	0.46
1:B:255:VAL:HG22	1:B:261:PHE:CD2	2.51	0.46
1:C:168:VAL:HG23	1:C:172:GLN:HB2	1.97	0.46
1:A:84:ALA:O	1:A:87:TYR:HB2	2.16	0.46
1:A:144:PRO:O	1:A:145:ALA:C	2.56	0.46
1:B:43:SER:O	1:B:46:ASN:HB3	2.17	0.45
1:B:292:LYS:O	1:B:293:LYS:C	2.59	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:172:GLN:C	1:A:174:PRO:HD2	2.39	0.45
1:A:215:LYS:O	1:A:222:ASN:ND2	2.49	0.45
1:C:32:THR:O	1:C:32:THR:HG22	2.16	0.45
1:A:222:ASN:CG	1:A:222:ASN:O	2.58	0.45
1:A:293:LYS:NZ	1:A:296:ILE:HD11	2.31	0.45
1:B:66:LEU:HD12	1:B:70:PHE:HA	1.97	0.45
1:B:127:ASN:HB3	1:B:130:VAL:CG1	2.46	0.45
1:B:211:ASP:HB3	1:B:213:GLN:HG3	1.98	0.45
1:C:63:VAL:H	1:C:102:GLN:HE21	1.63	0.45
1:A:269:PRO:HD2	1:A:273:ARG:HH11	1.78	0.45
1:A:302:VAL:O	1:A:305:PRO:HD2	2.17	0.45
1:C:112:THR:HG22	1:C:156:ALA:HB3	1.97	0.45
1:A:63:VAL:CG1	1:A:102:GLN:HG2	2.45	0.45
1:A:308:THR:O	1:A:311:ALA:HB3	2.17	0.45
1:B:173:LEU:HD23	1:B:173:LEU:HA	1.84	0.45
1:B:316:ASP:N	1:B:317:PRO:CD	2.79	0.45
1:A:79:ASN:O	1:A:80:VAL:C	2.58	0.45
1:A:230:ALA:CB	1:A:300:ASN:HD21	2.30	0.45
1:C:32:THR:O	1:C:32:THR:CG2	2.65	0.45
1:C:180:ARG:NH1	1:C:180:ARG:CG	2.77	0.45
1:A:93:ASP:HA	1:A:94:PRO:HD3	1.84	0.45
1:B:92:LEU:HD12	1:B:92:LEU:HA	1.84	0.45
1:B:274:VAL:C	1:B:276:ALA:N	2.74	0.45
1:A:136:GLN:O	1:A:136:GLN:CG	2.65	0.45
1:A:241:TYR:O	1:A:241:TYR:CD1	2.69	0.45
1:C:165:LEU:HD22	1:C:200:GLN:HB2	1.99	0.45
1:A:180:ARG:HB3	1:A:184:ARG:HH11	1.82	0.45
1:C:30:ARG:HD2	1:C:64:GLN:HE21	1.82	0.45
1:B:263:PHE:CB	1:B:295:LEU:HD11	2.48	0.44
1:B:295:LEU:C	1:B:298:VAL:HG12	2.40	0.44
1:B:74:GLU:CD	1:C:347:ARG:NH1	2.76	0.44
1:A:210:LEU:HB2	1:A:222:ASN:OD1	2.17	0.44
1:A:242:THR:CB	1:A:258:ASN:HD21	2.28	0.44
1:A:296:ILE:CG1	1:A:297:ASP:N	2.79	0.44
1:A:294:HIS:C	1:A:296:ILE:N	2.75	0.44
1:C:92:LEU:HD12	1:C:92:LEU:HA	1.81	0.44
1:A:216:MET:HA	1:A:222:ASN:O	2.17	0.44
1:C:25:VAL:HG22	1:C:165:LEU:HB3	1.98	0.44
1:C:243:ASP:HA	1:C:244:PRO:HD2	1.82	0.44
1:B:22:ARG:CB	1:B:23:PRO:CD	2.95	0.44
1:B:273:ARG:O	1:B:277:LEU:HD22	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:292:LYS:O	1:B:296:ILE:HG23	2.17	0.44
1:A:52:ASP:HB3	1:A:96:LYS:HZ1	1.82	0.44
1:A:290:LYS:HE2	1:A:290:LYS:HA	1.98	0.44
1:A:35:LEU:O	1:A:226:LEU:CD1	2.66	0.44
1:A:36:HIS:HA	1:A:226:LEU:HD12	1.99	0.44
1:A:193:VAL:HG21	1:A:341:GLN:HE21	1.82	0.44
1:A:296:ILE:HG13	1:A:297:ASP:H	1.78	0.44
1:A:296:ILE:O	1:A:300:ASN:HB2	2.17	0.44
1:C:270:ASP:O	1:C:270:ASP:CG	2.61	0.44
1:A:109:ALA:O	1:A:110:GLU:C	2.61	0.44
1:C:71:ASP:OD2	1:C:71:ASP:C	2.61	0.44
1:C:166:VAL:O	1:C:166:VAL:CG2	2.65	0.44
1:C:185:ARG:HH11	1:C:189:LEU:HD22	1.82	0.44
1:B:75:GLN:O	1:B:79:ASN:ND2	2.51	0.44
1:B:117:ASN:HD22	1:C:146:GLY:HA3	1.81	0.44
1:A:121:VAL:O	1:A:125:ARG:HB2	2.18	0.44
1:A:226:LEU:HD12	1:A:226:LEU:N	2.29	0.44
1:A:233:VAL:O	1:A:236:LYS:HB3	2.18	0.44
1:C:25:VAL:HG13	1:C:167:PRO:HD3	1.99	0.44
1:C:150:TYR:N	1:C:151:PRO:CD	2.80	0.44
1:B:268:ASP:OD1	1:B:273:ARG:NH2	2.51	0.43
1:A:35:LEU:CG	1:A:86:ASP:HB3	2.48	0.43
1:A:150:TYR:O	1:A:154:GLN:HG3	2.18	0.43
1:A:237:VAL:O	1:A:240:MET:CB	2.66	0.43
1:C:173:LEU:N	1:C:174:PRO:HD2	2.33	0.43
1:C:292:LYS:O	1:C:296:ILE:HG13	2.18	0.43
1:B:154:GLN:CD	2:B:1154:HRP:NE1	2.74	0.43
1:B:23:PRO:HD2	1:B:54:ALA:HB2	2.00	0.43
1:C:248:ARG:O	1:C:251:ASP:HB2	2.18	0.43
1:C:277:LEU:HD13	1:C:277:LEU:HA	1.83	0.43
1:B:241:TYR:OH	1:B:244:PRO:HD3	2.18	0.43
1:A:242:THR:HB	1:A:258:ASN:ND2	2.31	0.43
1:C:79:ASN:O	1:C:83:VAL:HG22	2.17	0.43
1:C:93:ASP:O	1:C:97:THR:HG23	2.18	0.43
1:C:201:LEU:CD2	1:C:202:SER:N	2.81	0.43
1:A:149:VAL:O	1:A:149:VAL:CG1	2.66	0.43
1:A:252:PRO:CB	1:A:282:ARG:O	2.66	0.43
1:A:261:PHE:CZ	1:A:281:TYR:CG	3.07	0.43
1:A:346:MET:O	1:A:347:ARG:HB2	2.17	0.43
1:B:58:VAL:HG23	1:B:58:VAL:O	2.19	0.43
1:B:187:ASN:OD1	1:B:192:PRO:HA	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:LEU:C	1:A:40:LEU:CD2	2.91	0.43
1:A:154:GLN:NE2	1:A:175:MET:HG3	2.33	0.43
1:A:300:ASN:O	1:A:304:ALA:HB2	2.19	0.43
1:B:186:PHE:CD1	1:B:186:PHE:O	2.72	0.43
1:B:316:ASP:OD1	1:B:318:ASP:OD2	2.36	0.43
1:C:241:TYR:CZ	1:C:244:PRO:HD3	2.54	0.43
1:B:70:PHE:CD1	1:B:70:PHE:C	2.96	0.43
1:B:111:LEU:HD11	1:B:115:PHE:CE1	2.53	0.43
1:A:237:VAL:O	1:A:240:MET:HG3	2.18	0.43
1:C:243:ASP:OD2	1:C:246:HIS:HB2	2.19	0.43
1:B:22:ARG:HB2	1:B:23:PRO:CD	2.49	0.43
1:A:29:ASP:O	1:A:31:PRO:HD3	2.19	0.43
1:A:245:GLY:O	1:A:247:LEU:HD13	2.19	0.43
1:A:255:VAL:HG22	1:A:261:PHE:CD1	2.53	0.43
1:C:30:ARG:H	1:C:30:ARG:HG2	1.62	0.43
1:C:99:CYS:O	1:C:327:GLY:HA3	2.19	0.43
1:B:59:LEU:HD12	1:B:100:VAL:O	2.19	0.42
1:B:122:SER:OG	1:C:141:GLU:HB3	2.18	0.42
1:B:258:ASN:HA	1:B:259:PRO:HD2	1.94	0.42
1:A:149:VAL:CG1	1:A:152:VAL:HG13	2.49	0.42
1:A:180:ARG:HG2	1:A:197:PRO:HD2	2.00	0.42
1:A:247:LEU:HD23	1:A:248:ARG:NH2	2.34	0.42
1:B:68:ASP:C	1:B:69:HIS:CD2	2.97	0.42
1:B:76:VAL:O	1:B:77:ARG:C	2.61	0.42
1:A:200:GLN:HE21	1:A:200:GLN:HB3	1.68	0.42
1:B:296:ILE:HD12	1:B:296:ILE:C	2.45	0.42
1:B:296:ILE:HD12	1:B:296:ILE:O	2.18	0.42
1:C:24:ARG:HD3	1:C:55:GLU:OE2	2.19	0.42
1:A:105:VAL:HA	1:A:106:PRO:HD2	1.86	0.42
1:A:119:VAL:HG22	1:A:120:THR:N	2.34	0.42
1:A:129:THR:O	1:A:132:ALA:HB3	2.19	0.42
1:C:132:ALA:O	1:C:135:ALA:HB3	2.19	0.42
1:C:268:ASP:OD1	1:C:294:HIS:HE1	2.02	0.42
1:B:248:ARG:NE	1:B:251:ASP:OD2	2.51	0.42
1:A:58:VAL:O	1:A:58:VAL:CG1	2.68	0.42
1:B:255:VAL:HG13	1:B:261:PHE:CG	2.55	0.42
2:B:1154:HRP:H	2:B:1154:HRP:CD2	2.31	0.42
1:C:36:HIS:CB	1:C:216:MET:HE2	2.50	0.42
1:C:63:VAL:HB	1:C:102:GLN:NE2	2.34	0.42
1:C:254:ARG:HG2	1:C:256:GLU:HG2	2.02	0.42
1:A:291:VAL:HG13	1:A:292:LYS:N	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:HIS:CD2	1:B:69:HIS:N	2.88	0.42
1:B:149:VAL:O	1:B:149:VAL:CG1	2.67	0.42
1:B:218:LYS:HE3	1:B:218:LYS:HB2	1.91	0.42
1:A:274:VAL:C	1:A:276:ALA:N	2.78	0.42
1:C:133:GLU:HA	1:C:136:GLN:HG3	2.01	0.42
1:C:270:ASP:HA	1:C:271:PRO:HD2	1.73	0.42
1:B:32:THR:O	1:B:75:GLN:NE2	2.53	0.42
1:B:154:GLN:CG	2:B:1154:HRP:HE1	2.06	0.42
1:B:232:GLU:O	1:B:235:ARG:N	2.53	0.42
1:A:30:ARG:O	1:A:31:PRO:C	2.63	0.42
1:A:234:ALA:O	1:A:238:MET:HG2	2.20	0.42
1:C:245:GLY:O	1:C:247:LEU:HD13	2.20	0.42
1:B:116:LEU:O	1:C:145:ALA:HB3	2.20	0.42
1:B:184:ARG:HG2	1:B:184:ARG:NH2	2.33	0.42
1:B:299:LEU:HA	1:B:302:VAL:CG1	2.49	0.42
1:A:246:HIS:ND1	1:A:288:ASP:OD1	2.53	0.42
1:C:169:GLY:C	1:C:171:ASP:N	2.78	0.42
1:B:44:LEU:HA	1:B:47:ARG:CD	2.49	0.41
1:B:61:ALA:HB1	1:B:64:GLN:HB3	2.02	0.41
1:A:61:ALA:HB1	1:A:64:GLN:HB3	2.01	0.41
1:C:85:LEU:CD1	1:C:310:ARG:CZ	2.98	0.41
1:C:93:ASP:HA	1:C:94:PRO:HD3	1.95	0.41
1:C:165:LEU:HA	1:C:165:LEU:HD23	1.79	0.41
1:A:58:VAL:CG1	1:A:99:CYS:HA	2.51	0.41
1:A:237:VAL:CG2	1:A:296:ILE:HG22	2.50	0.41
1:A:258:ASN:HA	1:A:259:PRO:HD2	1.90	0.41
1:C:58:VAL:HG23	1:C:99:CYS:HA	2.01	0.41
1:C:324:VAL:HG23	1:C:325:THR:N	2.36	0.41
1:B:58:VAL:CG2	1:B:99:CYS:HA	2.49	0.41
1:B:243:ASP:OD2	1:B:254:ARG:HG3	2.20	0.41
1:B:260:VAL:CG2	1:B:261:PHE:N	2.82	0.41
1:A:30:ARG:NH2	1:A:68:ASP:CB	2.82	0.41
1:A:327:GLY:O	1:A:330:ARG:HB3	2.20	0.41
1:C:295:LEU:HD22	1:C:299:LEU:CG	2.51	0.41
1:B:173:LEU:N	1:B:174:PRO:CD	2.83	0.41
1:B:304:ALA:N	1:B:305:PRO:HD2	2.35	0.41
1:A:104:ALA:CB	1:A:328:THR:HG21	2.50	0.41
1:A:106:PRO:C	1:A:108:LEU:N	2.76	0.41
1:A:211:ASP:C	1:A:213:GLN:N	2.79	0.41
1:A:242:THR:HG23	1:A:292:LYS:CE	2.51	0.41
1:C:133:GLU:O	1:C:136:GLN:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:211:ASP:CB	1:B:213:GLN:HG3	2.51	0.41
1:C:51:GLN:HB2	1:C:96:LYS:HB3	2.02	0.41
1:C:127:ASN:O	1:C:131:LYS:HB2	2.20	0.41
1:B:287:GLY:O	1:B:288:ASP:C	2.61	0.41
1:A:58:VAL:HG12	1:A:99:CYS:HA	2.03	0.41
1:A:293:LYS:C	1:A:296:ILE:HG12	2.45	0.41
1:C:27:THR:HG22	1:C:167:PRO:HD2	2.01	0.41
1:C:187:ASN:CG	1:C:192:PRO:HA	2.44	0.41
1:A:225:ALA:C	1:A:227:GLY:H	2.28	0.41
1:A:308:THR:HG22	1:A:312:GLU:OE2	2.20	0.41
1:C:238:MET:CE	1:C:238:MET:CG	2.96	0.41
1:B:77:ARG:NE	1:C:347:ARG:NH2	2.68	0.41
1:B:333:GLU:HA	1:B:333:GLU:OE2	2.20	0.41
1:B:349:PHE:CD2	1:C:80:VAL:CG1	3.00	0.41
1:A:116:LEU:HD21	1:A:149:VAL:HG11	2.02	0.41
1:B:23:PRO:HD2	1:B:54:ALA:CB	2.51	0.41
1:A:32:THR:O	1:A:32:THR:CG2	2.67	0.41
1:A:97:THR:HG22	1:A:98:THR:N	2.35	0.41
1:A:122:SER:C	1:A:124:LEU:N	2.77	0.41
1:A:337:GLN:HG2	1:A:337:GLN:H	1.65	0.41
1:C:204:VAL:HG12	1:C:205:PRO:CD	2.32	0.41
1:C:320:VAL:O	1:C:323:PHE:HB3	2.20	0.41
1:B:93:ASP:OD2	1:B:94:PRO:HD2	2.21	0.41
1:B:277:LEU:HD22	1:B:277:LEU:H	1.85	0.41
1:A:218:LYS:HA	1:A:218:LYS:CE	2.42	0.41
1:A:254:ARG:CD	1:A:256:GLU:HB2	2.49	0.41
1:C:254:ARG:CD	1:C:256:GLU:O	2.69	0.41
1:C:258:ASN:HA	1:C:259:PRO:HD2	1.94	0.41
1:B:169:GLY:O	1:B:170:ASP:C	2.64	0.40
1:B:274:VAL:C	1:B:276:ALA:H	2.29	0.40
1:B:348:LEU:O	1:B:349:PHE:C	2.63	0.40
1:A:191:ALA:HA	1:A:192:PRO:HD3	1.87	0.40
1:A:255:VAL:HG13	1:A:261:PHE:CD1	2.57	0.40
1:B:114:TYR:CE1	1:B:346:MET:HE1	2.56	0.40
1:B:166:VAL:O	1:B:167:PRO:C	2.65	0.40
1:B:184:ARG:HE	1:B:196:GLU:CD	2.29	0.40
1:A:234:ALA:HA	1:A:296:ILE:CG2	2.51	0.40
1:A:167:PRO:HA	1:A:200:GLN:HB3	2.03	0.40
1:C:216:MET:HE3	1:C:224:ILE:N	2.36	0.40
1:A:72:ARG:N	1:A:73:PRO:CD	2.84	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	329/351 (94%)	282 (86%)	32 (10%)	15 (5%)	2	18
1	B	329/351 (94%)	299 (91%)	21 (6%)	9 (3%)	4	27
1	C	329/351 (94%)	307 (93%)	18 (6%)	4 (1%)	10	40
All	All	987/1053 (94%)	888 (90%)	71 (7%)	28 (3%)	4	27

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	22	ARG
1	B	61	ALA
1	A	61	ALA
1	C	202	SER
1	C	205	PRO
1	B	284	GLY
1	A	31	PRO
1	A	249	ALA
1	A	295	LEU
1	C	22	ARG
1	B	43	SER
1	A	205	PRO
1	A	213	GLN
1	A	269	PRO
1	A	37	LEU
1	C	279	ASP
1	B	283	ALA
1	A	242	THR
1	B	206	ARG
1	B	250	SER
1	A	73	PRO
1	A	206	ARG
1	A	238	MET
1	B	285	GLY

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Mol	Chain	Res	Type
1	A	192	PRO
1	A	208	PRO
1	A	22	ARG
1	B	31	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	256/282 (91%)	227 (89%)	29 (11%)	5	25
1	B	264/282 (94%)	229 (87%)	35 (13%)	4	20
1	C	262/282 (93%)	232 (88%)	30 (12%)	5	24
All	All	782/846 (92%)	688 (88%)	94 (12%)	5	23

All (94) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	22	ARG
1	B	40	LEU
1	B	47	ARG
1	B	50	LEU
1	B	52	ASP
1	B	68	ASP
1	B	75	GLN
1	B	81	LEU
1	B	85	LEU
1	B	92	LEU
1	B	95	GLN
1	B	97	THR
1	B	141	GLU
1	B	151	PRO
1	B	152	VAL
1	B	165	LEU
1	B	168	VAL
1	B	170	ASP

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Mol	Chain	Res	Type
1	B	171	ASP
1	B	176	LEU
1	B	185	ARG
1	B	189	LEU
1	B	201	LEU
1	B	226	LEU
1	B	254	ARG
1	B	262	THR
1	B	282	ARG
1	B	290	LYS
1	B	296	ILE
1	B	297	ASP
1	B	302	VAL
1	B	318	ASP
1	B	322	ARG
1	B	344	ARG
1	B	347	ARG
1	A	35	LEU
1	A	40	LEU
1	A	52	ASP
1	A	56	LEU
1	A	63	VAL
1	A	74	GLU
1	A	81	LEU
1	A	85	LEU
1	A	124	LEU
1	A	125	ARG
1	A	126	GLN
1	A	150	TYR
1	A	152	VAL
1	A	165	LEU
1	A	172	GLN
1	A	200	GLN
1	A	201	LEU
1	A	203	ARG
1	A	215	LYS
1	A	224	ILE
1	A	232	GLU
1	A	246	HIS
1	A	268	ASP
1	A	286	LEU
1	A	289	VAL

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Mol	Chain	Res	Type
1	A	290	LYS
1	A	300	ASN
1	A	306	ILE
1	A	348	LEU
1	C	30	ARG
1	C	46	ASN
1	C	80	VAL
1	C	81	LEU
1	C	85	LEU
1	C	92	LEU
1	C	97	THR
1	C	100	VAL
1	C	113	VAL
1	C	137	LYS
1	C	150	TYR
1	C	152	VAL
1	C	165	LEU
1	C	193	VAL
1	C	200	GLN
1	C	201	LEU
1	C	202	SER
1	C	204	VAL
1	C	205	PRO
1	C	210	LEU
1	C	231	ASP
1	C	248	ARG
1	C	256	GLU
1	C	277	LEU
1	C	286	LEU
1	C	295	LEU
1	C	298	VAL
1	C	312	GLU
1	C	321	LEU
1	C	330	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	36	HIS
1	B	39	HIS
1	B	46	ASN
1	B	51	GLN

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Mol	Chain	Res	Type
1	B	69	HIS
1	B	75	GLN
1	B	102	GLN
1	B	117	ASN
1	B	136	GLN
1	B	154	GLN
1	B	178	GLN
1	B	200	GLN
1	B	300	ASN
1	B	337	GLN
1	A	36	HIS
1	A	46	ASN
1	A	69	HIS
1	A	79	ASN
1	A	123	HIS
1	A	154	GLN
1	A	172	GLN
1	A	200	GLN
1	A	280	GLN
1	A	300	ASN
1	A	341	GLN
1	C	46	ASN
1	C	51	GLN
1	C	69	HIS
1	C	95	GLN
1	C	102	GLN
1	C	117	ASN
1	C	126	GLN
1	C	200	GLN
1	C	213	GLN
1	C	258	ASN
1	C	294	HIS
1	C	300	ASN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HRP	B	1154	-	16,17,17	2.02	5 (31%)	20,24,24	1.46	3 (15%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HRP	B	1154	-	-	4/8/8/8	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1154	HRP	OX-CZ3	-5.32	1.25	1.37
2	B	1154	HRP	CE3-CD2	3.31	1.45	1.39
2	B	1154	HRP	CH2-CZ3	2.47	1.43	1.39
2	B	1154	HRP	CB-CG	2.07	1.54	1.50
2	B	1154	HRP	CB-CA	-2.00	1.44	1.53

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1154	HRP	CA-CB-CG	-4.30	105.33	113.49
2	B	1154	HRP	CE3-CD2-CE2	-2.12	117.32	119.39
2	B	1154	HRP	OXT-C-O	2.05	128.73	124.08

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1154	HRP	CA-CB-CG-CD2
2	B	1154	HRP	CA-CB-CG-CD1
2	B	1154	HRP	O-C-CA-CB
2	B	1154	HRP	OXT-C-CA-CB

There are no ring outliers.

1 monomer is involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1154	HRP	15	0

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	331/351 (94%)	0.12	6 (1%) 67 42	27, 96, 156, 167	0
1	B	331/351 (94%)	-0.12	1 (0%) 90 76	19, 67, 126, 170	1 (0%)
1	C	331/351 (94%)	-0.15	2 (0%) 85 66	20, 64, 95, 123	0
All	All	993/1053 (94%)	-0.05	9 (0%) 81 58	19, 72, 140, 170	1 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	219	SER	3.8
1	A	274	VAL	3.7
1	C	311	ALA	3.3
1	A	220	LEU	3.1
1	A	291	VAL	2.7
1	A	263	PHE	2.5
1	A	25	VAL	2.5
1	B	247	LEU	2.2
1	C	141	GLU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	HRP	B	1154	16/16	0.88	0.15	88,92,107,109	0

6.5 Other polymers

There are no such residues in this entry.