



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:51 AM UTC

PDB ID : 1GL9 / pdb_00001gl9
Title : Archaeoglobus fulgidus reverse gyrase complexed with ADPNP
Authors : Rodriguez, A.C.; Stock, D.
Deposited on : 2001-08-30
Resolution : 3.20 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

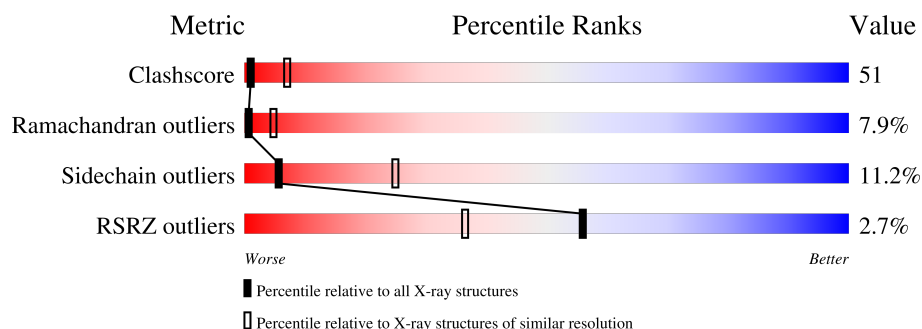
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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))
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	B	1054	2% 31% 51% 12% • •
1	C	1054	3% 31% 52% 11% • 5%

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	ANP	B	2055	X	-	-	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 16398 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.

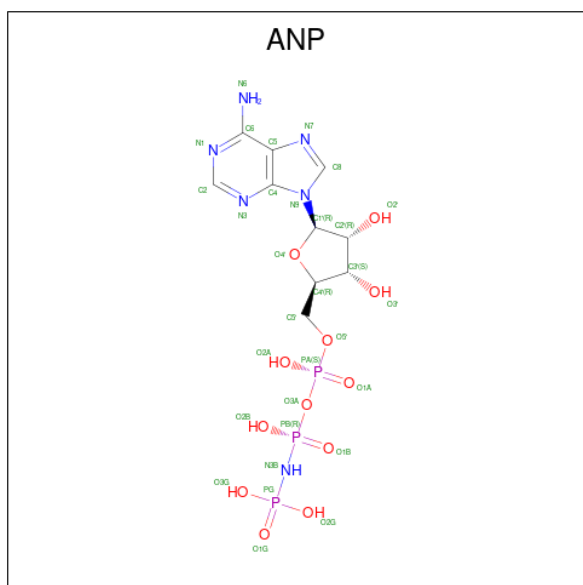
- Molecule 1 is a protein called REVERSE GYRASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	1020	Total	C	N	O	S	0	0	0
			8226	5246	1450	1505	25			
1	C	1005	Total	C	N	O	S	0	0	0
			8108	5169	1430	1483	26			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	719	LEU	PRO	engineered mutation	UNP O29238
B	1046	MET	LEU	engineered mutation	UNP O29238
C	719	LEU	PRO	engineered mutation	UNP O29238
C	1046	MET	LEU	engineered mutation	UNP O29238

- Molecule 2 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: $C_{10}H_{17}N_6O_{12}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
2	C	1	Total	C	N	O	P	0	0
			31	10	6	12	3		

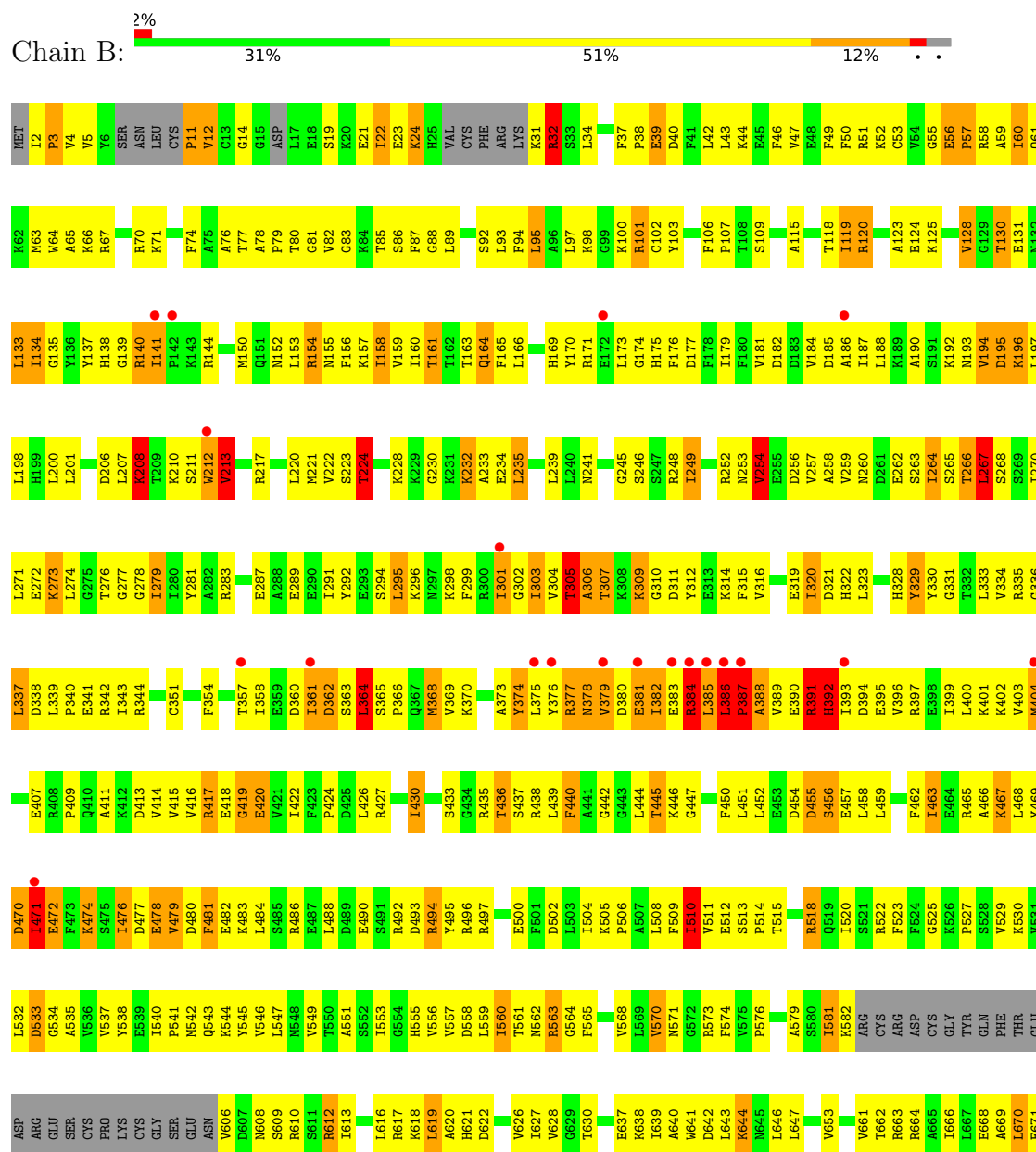
- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

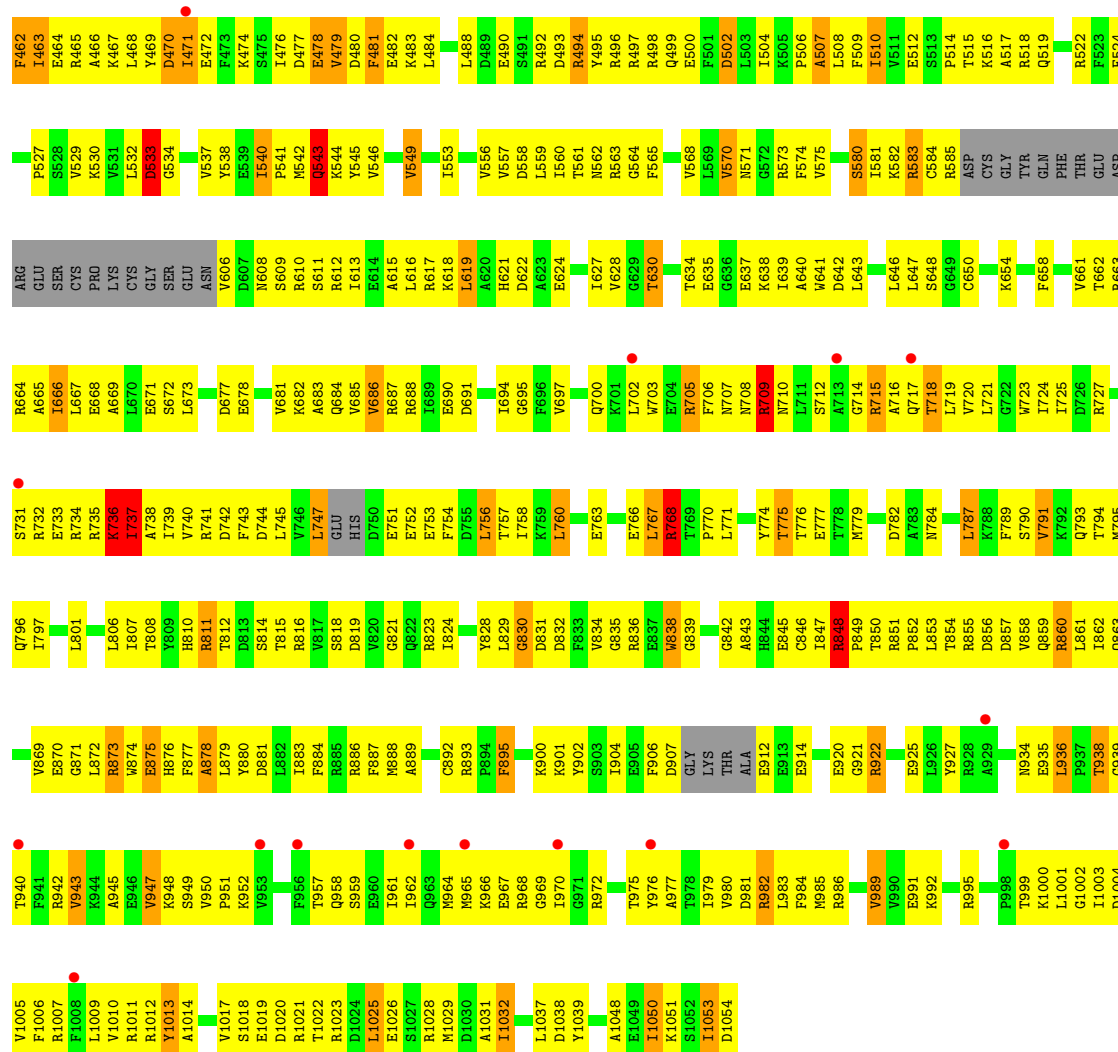
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

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: REVERSE GYRASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	132.41Å 68.69Å 134.00Å 90.00° 99.70° 90.00°	Depositor
Resolution (Å)	34.00 – 3.20 34.00 – 3.20	Depositor EDS
% Data completeness (in resolution range)	96.6 (34.00-3.20) 96.6 (34.00-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 3.18Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.256 , 0.332 0.247 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	69.1	Xtriage
Anisotropy	0.233	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 87.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.022 for l,-k,h	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	16398	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.78% 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: ANP, MG

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	B	0.60	0/8371	1.10	48/11265 (0.4%)
1	C	0.41	0/8246	0.92	35/11090 (0.3%)
All	All	0.51	0/16617	1.02	83/22355 (0.4%)

There are no bond length outliers.

All (83) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	391	ARG	N-CA-C	-17.87	91.57	113.01
1	B	388	ALA	N-CA-C	-17.20	92.68	113.15
1	B	386	LEU	CA-C-N	-14.30	101.96	119.84
1	B	386	LEU	C-N-CA	-14.30	101.96	119.84
1	B	246	SER	N-CA-C	10.08	123.84	110.53
1	B	478	GLU	N-CA-C	-9.64	100.08	113.20
1	C	305	THR	N-CA-C	8.89	129.73	110.80
1	C	478	GLU	N-CA-C	-8.66	101.43	113.20
1	B	878	ALA	N-CA-C	-8.10	102.45	111.28
1	C	384	ARG	N-CA-C	-8.02	103.62	113.41
1	B	169	HIS	N-CA-C	7.47	119.41	110.44
1	B	787	LEU	N-CA-C	-7.26	100.95	110.53
1	B	661	VAL	N-CA-C	7.15	113.86	106.21
1	B	384	ARG	N-CA-C	-6.96	104.30	114.39
1	B	986	ARG	N-CA-C	-6.86	104.89	112.72
1	C	224	THR	N-CA-C	-6.82	101.53	110.53
1	C	510	ILE	N-CA-C	6.80	118.34	107.73
1	B	305	THR	N-CA-C	6.73	125.13	110.80
1	B	768	ARG	N-CA-C	6.71	118.00	108.74
1	C	848	ARG	N-CA-C	6.69	114.15	108.13
1	B	3	PRO	N-CA-CB	6.65	110.23	103.25
1	C	304	VAL	N-CA-C	6.52	122.89	109.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1050	ILE	N-CA-C	6.49	116.63	110.53
1	B	718	THR	N-CA-C	-6.47	104.15	111.14
1	C	11	PRO	N-CA-CB	6.32	109.95	103.00
1	C	462	PHE	N-CA-C	-6.25	103.39	111.02
1	B	382	ILE	N-CA-C	-6.24	103.30	112.04
1	C	3	PRO	N-CA-CB	6.19	109.81	103.00
1	C	619	LEU	N-CA-C	-6.19	104.22	110.97
1	B	11	PRO	N-CA-CB	6.18	109.80	103.00
1	B	78	ALA	CA-C-N	6.16	126.13	119.78
1	B	78	ALA	C-N-CA	6.16	126.13	119.78
1	B	273	LYS	N-CA-C	-6.14	105.28	112.89
1	C	92	SER	N-CA-C	-6.13	104.52	111.14
1	B	733	GLU	N-CA-C	6.12	116.36	108.34
1	C	787	LEU	N-CA-C	-6.11	102.31	110.55
1	B	98	LYS	N-CA-C	-6.08	105.44	112.92
1	C	381	GLU	N-CA-C	-6.04	105.56	113.17
1	B	22	ILE	N-CA-C	6.03	118.14	108.85
1	C	718	THR	N-CA-C	-6.01	104.36	111.03
1	C	768	ARG	N-CA-C	5.96	118.46	109.41
1	B	211	SER	N-CA-C	-5.87	100.85	109.18
1	B	329	TYR	N-CA-C	5.86	117.47	111.14
1	B	769	THR	N-CA-C	5.86	117.08	109.64
1	C	368	MET	N-CA-C	-5.84	106.13	113.72
1	B	253	ASN	N-CA-C	-5.81	98.90	108.48
1	B	833	PHE	N-CA-C	5.76	118.07	109.25
1	B	619	LEU	N-CA-C	-5.71	104.74	110.97
1	B	581	ILE	N-CA-C	5.65	117.96	109.20
1	B	644	LYS	N-CA-C	-5.64	104.14	111.02
1	B	185	ASP	N-CA-C	-5.63	106.25	113.23
1	C	382	ILE	N-CA-C	-5.63	104.68	112.50
1	C	250	THR	N-CA-C	5.60	118.46	110.68
1	C	397	ARG	N-CA-C	5.58	119.57	112.87
1	C	412	LYS	N-CA-C	-5.58	105.07	113.89
1	C	155	ASN	N-CA-C	-5.56	106.50	113.72
1	C	583	ARG	N-CA-C	-5.49	106.44	113.02
1	B	224	THR	N-CA-C	-5.46	103.17	110.55
1	C	78	ALA	N-CA-C	5.46	113.04	108.13
1	B	510	ILE	N-CA-C	5.45	116.58	108.23
1	B	1012	ARG	N-CA-C	5.43	120.78	113.72
1	B	661	VAL	CB-CA-C	-5.41	107.16	113.22
1	C	144	ARG	N-CA-C	5.40	116.97	111.14
1	B	386	LEU	N-CA-C	-5.39	100.15	109.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	273	LYS	N-CA-C	-5.37	106.70	113.20
1	B	155	ASN	N-CA-C	-5.31	106.19	113.30
1	C	540	ILE	CA-C-N	5.28	126.44	119.84
1	C	540	ILE	C-N-CA	5.28	126.44	119.84
1	C	1048	ALA	N-CA-C	-5.26	105.23	110.97
1	C	775	THR	N-CA-C	-5.24	101.22	109.50
1	C	815	THR	N-CA-C	-5.21	106.94	113.72
1	C	878	ALA	N-CA-C	-5.20	105.04	111.33
1	C	85	THR	N-CA-C	-5.19	104.89	111.11
1	B	535	ALA	N-CA-C	5.16	116.91	108.76
1	C	98	LYS	N-CA-C	-5.16	106.30	112.59
1	B	254	VAL	N-CA-C	5.13	120.00	109.34
1	B	815	THR	N-CA-C	-5.09	105.98	113.61
1	C	843	ALA	N-CA-C	-5.09	105.98	113.61
1	B	78	ALA	N-CA-C	5.05	112.68	108.13
1	B	368	MET	N-CA-C	-5.04	106.05	113.61
1	B	513	SER	CA-C-N	5.01	124.31	118.85
1	B	513	SER	C-N-CA	5.01	124.31	118.85
1	B	1030	ASP	N-CA-C	-5.00	105.28	111.33

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	B	8226	0	8310	882	0
1	C	8108	0	8178	814	0
2	B	31	0	13	5	0
2	C	31	0	13	5	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
All	All	16398	0	16514	1693	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 51.

All (1693) 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:B:379:VAL:HG12	1:B:380:ASP:H	1.12	1.12
1:B:386:LEU:HD23	1:B:386:LEU:H	1.19	1.06
1:B:638:LYS:HB2	1:B:684:GLN:HG3	1.38	1.06
1:C:192:LYS:HB3	1:C:196:LYS:HE3	1.43	1.00
1:B:540:ILE:HG13	1:B:541:PRO:HD2	1.46	0.97
1:B:387:PRO:C	1:B:389:VAL:H	1.72	0.96
1:C:334:VAL:HG12	1:C:335:ARG:HG3	1.47	0.96
1:C:379:VAL:HG12	1:C:380:ASP:H	1.30	0.95
1:C:426:LEU:HD11	1:C:466:ALA:HA	1.49	0.94
1:C:301:ILE:O	1:C:320:ILE:HG13	1.67	0.94
1:C:305:THR:OG1	1:C:329:TYR:HB3	1.66	0.94
1:B:852:PRO:O	1:B:853:LEU:HD23	1.68	0.94
1:B:985:MET:HE2	1:B:986:ARG:HH21	1.33	0.93
1:B:760:LEU:HD23	1:B:760:LEU:H	1.34	0.93
1:C:244:ILE:HG22	1:C:245:GLY:N	1.84	0.93
1:B:771:LEU:HD12	1:B:927:TYR:HE2	1.34	0.92
1:B:358:ILE:H	1:B:358:ILE:HD12	1.31	0.92
1:C:556:VAL:HB	1:C:613:ILE:HD11	1.49	0.92
1:C:153:LEU:O	1:C:154:ARG:HB2	1.69	0.90
1:B:369:VAL:HG11	1:B:393:ILE:HG23	1.55	0.89
1:B:556:VAL:HB	1:B:613:ILE:HD11	1.52	0.88
1:C:386:LEU:HD23	1:C:386:LEU:H	1.37	0.88
1:B:904:ILE:HD11	1:B:911:ALA:HB3	1.56	0.87
1:B:784:ASN:HD22	1:B:959:SER:H	1.22	0.87
1:C:527:PRO:HB3	1:C:540:ILE:HB	1.56	0.87
1:B:869:VAL:HG21	1:B:872:LEU:HD12	1.56	0.87
1:B:556:VAL:HG23	1:B:557:VAL:HG13	1.54	0.86
1:B:784:ASN:ND2	1:B:959:SER:H	1.73	0.86
1:B:61:GLN:HE22	2:B:2055:ANP:HN61	0.91	0.86
1:B:322:HIS:O	1:B:323:LEU:HD12	1.75	0.86
1:C:892:CYS:SG	1:C:922:ARG:HB3	2.16	0.86
1:C:198:LEU:H	1:C:198:LEU:HD12	1.39	0.86
1:B:201:LEU:O	1:B:217:ARG:HB3	1.75	0.86
1:B:301:ILE:H	1:B:301:ILE:HD12	1.40	0.86
1:C:638:LYS:HB2	1:C:684:GLN:HG3	1.57	0.86
1:B:369:VAL:HG21	1:B:393:ILE:HG21	1.56	0.86
1:B:1001:LEU:O	1:B:1005:VAL:HG23	1.76	0.86
1:B:1028:ARG:O	1:B:1032:ILE:HG22	1.76	0.85
1:B:61:GLN:NE2	2:B:2055:ANP:HN61	1.74	0.85
1:B:384:ARG:HB3	1:B:384:ARG:HH11	1.40	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:395:GLU:O	1:B:399:ILE:HG12	1.77	0.85
1:C:39:GLU:CD	1:C:39:GLU:H	1.81	0.84
1:B:556:VAL:HB	1:B:613:ILE:CD1	2.06	0.84
1:B:508:LEU:HD21	1:B:510:ILE:HD11	1.56	0.84
1:B:745:LEU:HD13	1:B:747:LEU:HD11	1.57	0.83
1:C:1028:ARG:O	1:C:1032:ILE:HG22	1.78	0.83
1:C:305:THR:HG21	1:C:329:TYR:CD1	2.12	0.83
1:B:390:GLU:C	1:B:392:HIS:N	2.27	0.83
1:C:771:LEU:HD12	1:C:927:TYR:HE2	1.44	0.83
1:B:465:ARG:NH2	1:B:864:GLU:OE2	2.12	0.83
1:B:379:VAL:HG12	1:B:380:ASP:N	1.93	0.83
1:B:76:ALA:HB2	1:B:221:MET:HE2	1.61	0.82
1:C:610:ARG:HB3	1:C:610:ARG:HH11	1.43	0.82
1:C:834:VAL:HG12	1:C:835:GLY:H	1.43	0.82
1:B:103:TYR:HB2	1:B:176:PHE:CD2	2.15	0.82
1:B:198:LEU:H	1:B:198:LEU:HD12	1.44	0.81
1:B:381:GLU:O	1:B:384:ARG:HG2	1.79	0.81
1:B:32:ARG:HD3	1:B:32:ARG:H	1.44	0.81
1:B:58:ARG:H	1:B:61:GLN:HE21	1.29	0.81
1:C:609:SER:O	1:C:613:ILE:HD13	1.80	0.81
1:C:58:ARG:H	1:C:61:GLN:HE21	1.29	0.81
1:C:413:ASP:OD1	1:C:414:VAL:HG13	1.80	0.81
1:B:403:VAL:O	1:B:404:MET:HB3	1.77	0.80
1:C:784:ASN:ND2	1:C:959:SER:H	1.80	0.80
1:C:305:THR:HG22	1:C:306:ALA:H	1.45	0.80
1:C:134:ILE:HD13	1:C:135:GLY:N	1.96	0.80
1:C:628:VAL:HG11	1:C:640:ALA:HB2	1.63	0.80
1:B:1043:LEU:HD23	1:B:1046:MET:HE2	1.61	0.79
1:C:58:ARG:HB2	1:C:61:GLN:NE2	1.97	0.79
1:B:369:VAL:HG21	1:B:393:ILE:CG2	2.12	0.79
1:B:470:ASP:O	1:B:472:GLU:HB2	1.82	0.79
1:B:921:GLY:O	1:B:925:GLU:HG2	1.83	0.79
1:C:560:ILE:HG13	1:C:561:THR:H	1.48	0.79
1:B:555:HIS:HA	1:B:639:ILE:CD1	2.12	0.79
1:C:824:ILE:HG23	1:C:852:PRO:HG3	1.64	0.79
1:C:470:ASP:O	1:C:472:GLU:N	2.16	0.79
1:B:224:THR:HG22	1:B:245:GLY:O	1.82	0.78
1:B:256:ASP:HB3	1:B:474:LYS:HD3	1.65	0.78
1:C:610:ARG:HB3	1:C:610:ARG:NH1	1.98	0.78
1:B:492:ARG:HH11	1:B:492:ARG:HG3	1.49	0.78
1:C:1010:VAL:O	1:C:1014:ALA:HB2	1.83	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:470:ASP:O	1:B:472:GLU:N	2.17	0.78
1:B:719:LEU:HD11	1:B:1019:GLU:HA	1.66	0.78
1:C:921:GLY:O	1:C:925:GLU:HG2	1.84	0.78
1:C:357:THR:HB	1:C:360:ASP:HB2	1.66	0.78
1:C:58:ARG:HB2	1:C:61:GLN:HE21	1.49	0.78
1:B:95:LEU:HD13	1:B:100:LYS:HD3	1.65	0.77
1:C:301:ILE:HD12	1:C:301:ILE:H	1.49	0.77
1:B:163:THR:HG21	1:B:186:ALA:HB1	1.67	0.77
1:C:85:THR:HB	2:C:2055:ANP:O1A	1.84	0.77
1:C:797:ILE:HG23	1:C:876:HIS:ND1	2.00	0.77
1:C:103:TYR:HB2	1:C:176:PHE:CD2	2.19	0.77
1:B:56:GLU:OE1	1:B:57:PRO:HD2	1.83	0.77
1:B:709:ARG:HG2	1:B:710:ASN:H	1.49	0.77
1:B:768:ARG:HB3	1:B:768:ARG:NH1	2.00	0.77
1:B:608:ASN:HD22	1:B:610:ARG:HH11	1.33	0.76
1:B:22:ILE:O	1:B:576:PRO:HD2	1.86	0.76
1:C:992:LYS:HE2	1:C:992:LYS:HA	1.66	0.76
1:C:396:VAL:HG11	1:C:399:ILE:HB	1.66	0.76
1:C:712:SER:HB2	1:C:986:ARG:HH11	1.49	0.76
1:C:436:THR:HG23	1:C:447:GLY:HA3	1.68	0.76
1:C:848:ARG:HB2	1:C:887:PHE:CD2	2.21	0.76
1:C:712:SER:HB2	1:C:986:ARG:NH1	2.01	0.75
1:B:626:VAL:HB	1:B:653:VAL:HG22	1.69	0.75
1:B:279:ILE:HG22	1:B:281:TYR:CE1	2.22	0.75
1:B:687:ARG:HA	1:B:1029:MET:CE	2.17	0.75
1:C:414:VAL:HG12	1:C:424:PRO:HD3	1.68	0.75
1:B:58:ARG:HB2	1:B:61:GLN:HE21	1.50	0.75
1:C:985:MET:HE2	1:C:986:ARG:HH21	1.52	0.75
1:B:436:THR:HG23	1:B:447:GLY:HA3	1.69	0.75
1:B:387:PRO:C	1:B:389:VAL:N	2.37	0.75
1:B:391:ARG:O	1:B:391:ARG:HD3	1.86	0.75
1:C:720:VAL:HG13	1:C:961:ILE:HD11	1.68	0.75
1:B:763:GLU:HG3	1:B:900:LYS:HG2	1.69	0.74
1:C:50:PHE:CD2	1:C:57:PRO:HG3	2.22	0.74
1:B:687:ARG:HA	1:B:1029:MET:HE1	1.70	0.74
1:B:703:TRP:CD1	1:B:709:ARG:HA	2.21	0.74
1:B:32:ARG:H	1:B:32:ARG:CD	1.98	0.74
1:B:560:ILE:HD12	1:B:579:ALA:HB2	1.70	0.74
1:C:975:THR:HG22	1:C:979:ILE:HG13	1.67	0.74
1:C:787:LEU:HD13	1:C:789:PHE:CE2	2.23	0.74
1:C:1029:MET:O	1:C:1032:ILE:HG23	1.87	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:192:LYS:HB3	1:B:196:LYS:HE3	1.70	0.74
1:B:774:TYR:CZ	1:B:886:ARG:HG3	2.22	0.74
1:B:873:ARG:H	1:B:876:HIS:HD2	1.35	0.74
1:C:922:ARG:HH21	1:C:922:ARG:HG3	1.51	0.74
1:B:384:ARG:HB3	1:B:384:ARG:NH1	2.03	0.73
1:B:103:TYR:HE1	1:B:161:THR:HG23	1.53	0.73
1:C:63:MET:HE3	1:C:67:ARG:HH21	1.52	0.73
1:B:397:ARG:HA	1:B:400:LEU:HD12	1.68	0.73
1:C:303:ILE:HG12	1:C:304:VAL:H	1.52	0.73
1:B:476:ILE:HG22	1:B:478:GLU:OE2	1.89	0.73
1:C:790:SER:HA	1:C:958:GLN:HE22	1.53	0.73
1:C:438:ARG:O	1:C:444:LEU:HD12	1.87	0.73
1:B:735:ARG:HA	1:B:950:VAL:O	1.88	0.73
1:B:784:ASN:HD22	1:B:959:SER:N	1.86	0.73
1:B:262:GLU:HG2	1:B:454:ASP:CG	2.13	0.73
1:B:438:ARG:C	1:B:444:LEU:HD12	2.14	0.73
1:B:370:LYS:O	1:B:373:ALA:HB3	1.88	0.73
1:B:262:GLU:HG2	1:B:454:ASP:OD2	1.89	0.72
1:C:241:ASN:O	1:C:530:LYS:HD2	1.89	0.72
1:B:413:ASP:OD1	1:B:414:VAL:HG13	1.88	0.72
1:C:784:ASN:HD22	1:C:959:SER:H	1.37	0.72
1:B:529:VAL:HG12	1:B:530:LYS:N	2.05	0.72
1:B:608:ASN:HD22	1:B:610:ARG:NH1	1.87	0.72
1:B:904:ILE:HD12	1:B:904:ILE:O	1.89	0.72
1:C:135:GLY:HA3	1:C:156:PHE:CE1	2.24	0.72
1:C:875:GLU:O	1:C:878:ALA:HB3	1.90	0.72
1:B:70:ARG:HH21	1:B:618:LYS:HD2	1.53	0.72
1:B:271:LEU:HD23	1:B:481:PHE:HZ	1.53	0.72
1:B:690:GLU:OE2	1:B:715:ARG:NH1	2.23	0.72
1:C:608:ASN:HD22	1:C:610:ARG:NH1	1.88	0.72
1:B:1037:LEU:HD23	1:B:1038:ASP:N	2.04	0.72
1:C:181:VAL:HB	1:C:222:VAL:HG12	1.71	0.72
1:C:370:LYS:O	1:C:373:ALA:HB3	1.90	0.72
1:C:738:ALA:HA	1:C:947:VAL:HA	1.72	0.72
1:C:1007:ARG:HA	1:C:1010:VAL:HG12	1.70	0.72
1:B:376:TYR:O	1:B:377:ARG:HB3	1.90	0.72
1:B:212:TRP:O	1:B:213:VAL:HG13	1.90	0.71
1:B:895:PHE:H	1:B:895:PHE:HD2	1.38	0.71
1:B:1012:ARG:HG2	1:B:1012:ARG:HH11	1.54	0.71
1:B:58:ARG:H	1:B:61:GLN:NE2	1.87	0.71
1:B:736:LYS:O	1:B:737:ILE:HG22	1.89	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:540:ILE:HG13	1:B:541:PRO:CD	2.20	0.71
1:C:103:TYR:HB3	1:C:179:ILE:HG13	1.71	0.71
1:B:390:GLU:C	1:B:392:HIS:H	1.83	0.71
1:C:212:TRP:O	1:C:213:VAL:HG13	1.89	0.71
1:B:962:ILE:HD13	1:B:965:MET:HE3	1.72	0.71
1:C:763:GLU:HG3	1:C:900:LYS:HG2	1.72	0.71
1:B:731:SER:O	1:B:951:PRO:HB3	1.90	0.71
1:C:490:GLU:O	1:C:493:ASP:HB3	1.89	0.71
1:B:58:ARG:HH21	1:B:83:GLY:N	1.89	0.71
1:B:358:ILE:HD11	1:B:419:GLY:C	2.15	0.71
1:B:821:GLY:HA3	1:B:848:ARG:HH22	1.56	0.71
1:B:787:LEU:HD13	1:B:789:PHE:CE2	2.26	0.70
1:C:163:THR:HG21	1:C:186:ALA:HB1	1.72	0.70
1:C:176:PHE:HB2	1:C:179:ILE:HD11	1.73	0.70
1:C:480:ASP:O	1:C:482:GLU:N	2.24	0.70
1:C:617:ARG:HA	1:C:647:LEU:HD22	1.72	0.70
1:B:884:PHE:CE2	1:B:888:MET:HG3	2.26	0.70
1:B:1009:LEU:H	1:B:1009:LEU:HD22	1.56	0.70
1:C:492:ARG:HH11	1:C:492:ARG:HG3	1.55	0.70
1:C:736:LYS:C	1:C:737:ILE:HG23	2.17	0.70
1:B:768:ARG:HB3	1:B:768:ARG:HH11	1.56	0.70
1:C:735:ARG:HA	1:C:950:VAL:O	1.91	0.70
1:B:361:ILE:HG22	1:B:397:ARG:HB3	1.73	0.70
1:B:480:ASP:O	1:B:482:GLU:N	2.25	0.70
1:C:360:ASP:O	1:C:362:ASP:N	2.25	0.70
1:B:438:ARG:O	1:B:444:LEU:HD12	1.92	0.69
1:B:618:LYS:O	1:B:621:HIS:HB3	1.91	0.69
1:C:301:ILE:HG22	1:C:302:GLY:H	1.58	0.69
1:B:334:VAL:HG12	1:B:335:ARG:HG3	1.74	0.69
1:B:971:GLY:O	1:B:972:ARG:HD3	1.92	0.69
1:B:490:GLU:O	1:B:493:ASP:HB3	1.92	0.69
1:C:540:ILE:HG13	1:C:541:PRO:HD2	1.72	0.69
1:B:791:VAL:HG13	1:B:958:GLN:NE2	2.07	0.69
1:B:1037:LEU:HD23	1:B:1038:ASP:H	1.56	0.69
1:B:638:LYS:CB	1:B:684:GLN:HG3	2.21	0.69
1:C:244:ILE:HG22	1:C:245:GLY:H	1.55	0.69
1:C:608:ASN:HD22	1:C:610:ARG:HH11	1.37	0.69
1:C:747:LEU:HD12	1:C:747:LEU:N	2.07	0.69
1:C:1012:ARG:O	1:C:1013:TYR:HB2	1.92	0.69
1:B:309:LYS:H	1:B:309:LYS:HD2	1.57	0.69
1:B:426:LEU:HD11	1:B:466:ALA:HA	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:39:GLU:HG2	1:C:70:ARG:HG2	1.73	0.69
1:B:305:THR:CG2	1:B:306:ALA:N	2.55	0.69
1:B:364:LEU:HD23	1:B:368:MET:SD	2.33	0.69
1:B:479:VAL:HG23	1:B:481:PHE:H	1.56	0.69
1:B:470:ASP:O	1:B:471:ILE:C	2.35	0.68
1:C:834:VAL:HG12	1:C:835:GLY:N	2.07	0.68
1:C:510:ILE:CD1	1:C:616:LEU:HD22	2.23	0.68
1:C:642:ASP:C	1:C:646:LEU:HD12	2.18	0.68
1:C:1001:LEU:O	1:C:1005:VAL:HG23	1.93	0.68
1:B:379:VAL:CG1	1:B:380:ASP:H	1.97	0.68
1:C:58:ARG:H	1:C:61:GLN:NE2	1.91	0.68
1:C:95:LEU:HD13	1:C:100:LYS:HD3	1.74	0.68
1:B:617:ARG:HA	1:B:647:LEU:HD22	1.74	0.68
1:B:985:MET:HE2	1:B:986:ARG:NH2	2.07	0.68
1:B:747:LEU:HD12	1:B:747:LEU:N	2.08	0.68
1:C:476:ILE:HG22	1:C:478:GLU:OE2	1.93	0.68
1:B:212:TRP:CD1	1:B:212:TRP:H	2.11	0.68
1:B:419:GLY:O	1:B:420:GLU:HB2	1.93	0.68
1:C:337:LEU:HD21	1:C:343:ILE:HD11	1.76	0.68
1:C:259:VAL:HG12	1:C:261:ASP:H	1.59	0.68
1:C:754:PHE:O	1:C:943:VAL:HG23	1.94	0.68
1:B:736:LYS:HA	1:B:736:LYS:NZ	2.08	0.68
1:C:248:ARG:C	1:C:249:ILE:HG12	2.18	0.68
1:C:257:VAL:HB	1:C:450:PHE:CD1	2.28	0.68
1:B:610:ARG:NH1	1:B:610:ARG:HB3	2.08	0.67
1:C:296:LYS:HA	1:C:301:ILE:HG12	1.76	0.67
1:C:305:THR:O	1:C:307:THR:N	2.27	0.67
1:C:873:ARG:H	1:C:876:HIS:HD2	1.41	0.67
1:B:916:ILE:HG22	1:B:932:VAL:HA	1.76	0.67
1:C:305:THR:HG21	1:C:329:TYR:HD1	1.60	0.67
1:B:248:ARG:HG3	1:B:249:ILE:N	2.08	0.67
1:B:616:LEU:HA	1:B:619:LEU:HD12	1.77	0.67
1:B:2:ILE:N	1:B:11:PRO:N	2.43	0.67
1:B:426:LEU:O	1:B:430:ILE:HD12	1.94	0.67
1:B:784:ASN:ND2	1:B:959:SER:N	2.42	0.67
1:C:303:ILE:HG12	1:C:304:VAL:N	2.09	0.67
1:C:379:VAL:HG12	1:C:380:ASP:N	2.04	0.67
1:C:858:VAL:HG21	1:C:880:TYR:CE1	2.30	0.67
1:B:305:THR:HG23	1:B:329:TYR:CE1	2.30	0.67
1:B:363:SER:O	1:B:364:LEU:HB2	1.95	0.67
1:B:834:VAL:HG12	1:B:835:GLY:H	1.60	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:304:VAL:O	1:B:305:THR:HB	1.95	0.66
1:C:3:PRO:O	1:C:11:PRO:N	2.28	0.66
1:C:735:ARG:HB3	1:C:949:SER:HB3	1.78	0.66
1:B:265:SER:O	1:B:267:LEU:N	2.28	0.66
1:B:791:VAL:HG13	1:B:958:GLN:HE21	1.61	0.66
1:B:809:TYR:CE2	1:B:811:ARG:HB2	2.31	0.66
1:C:376:TYR:O	1:C:377:ARG:CB	2.42	0.66
1:C:608:ASN:ND2	1:C:610:ARG:HH11	1.93	0.66
1:C:510:ILE:HD11	1:C:616:LEU:HD22	1.75	0.66
1:B:194:VAL:O	1:B:195:ASP:C	2.36	0.66
1:B:800:GLU:OE2	1:B:876:HIS:HE1	1.78	0.66
1:B:305:THR:HG22	1:B:306:ALA:N	2.11	0.66
1:B:749:HIS:CD2	1:B:751:GLU:HB2	2.30	0.66
1:C:63:MET:HE3	1:C:67:ARG:NH2	2.11	0.66
1:C:694:ILE:HD13	1:C:715:ARG:NH2	2.10	0.66
1:C:904:ILE:HD12	1:C:904:ILE:O	1.96	0.66
1:B:855:ARG:O	1:B:859:GLN:HG3	1.96	0.66
1:B:872:LEU:HA	1:B:876:HIS:CD2	2.31	0.65
1:C:58:ARG:HH21	1:C:83:GLY:N	1.94	0.65
1:C:419:GLY:O	1:C:420:GLU:HB2	1.95	0.65
1:B:739:ILE:O	1:B:945:ALA:HA	1.96	0.65
1:B:85:THR:HB	2:B:2055:ANP:O1A	1.96	0.65
1:B:357:THR:HB	1:B:360:ASP:HB2	1.77	0.65
1:C:376:TYR:O	1:C:377:ARG:HB2	1.95	0.65
1:B:51:ARG:HA	1:B:55:GLY:O	1.96	0.65
1:C:700:GLN:HA	1:C:703:TRP:CE3	2.32	0.65
1:B:135:GLY:HA3	1:B:156:PHE:CD1	2.31	0.65
1:B:754:PHE:O	1:B:943:VAL:HG23	1.97	0.65
1:B:849:PRO:HD3	1:B:883:ILE:HG22	1.78	0.65
1:C:468:LEU:HD11	1:C:853:LEU:HD21	1.79	0.65
1:B:270:ILE:O	1:B:273:LYS:HB3	1.96	0.65
1:C:301:ILE:HD12	1:C:301:ILE:N	2.11	0.65
1:C:396:VAL:CG1	1:C:399:ILE:HB	2.27	0.65
1:C:756:LEU:O	1:C:940:THR:HG23	1.96	0.65
1:C:895:PHE:HB2	1:C:920:GLU:O	1.97	0.65
1:B:438:ARG:HG3	1:B:439:LEU:N	2.12	0.65
1:B:1010:VAL:O	1:B:1014:ALA:HB2	1.97	0.65
1:C:745:LEU:HD13	1:C:747:LEU:HD11	1.79	0.65
1:C:873:ARG:HB2	1:C:875:GLU:OE2	1.95	0.65
1:C:60:ILE:HD13	1:C:546:VAL:HG21	1.78	0.65
1:C:542:MET:HE1	1:C:544:LYS:HB2	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:301:ILE:HG22	1:B:302:GLY:H	1.62	0.64
1:B:771:LEU:HD12	1:B:927:TYR:CE2	2.24	0.64
1:B:905:GLU:HG2	1:B:910:THR:HB	1.79	0.64
1:B:259:VAL:HG22	1:B:477:ASP:OD1	1.97	0.64
1:B:439:LEU:CD1	1:B:444:LEU:HD13	2.27	0.64
1:B:439:LEU:HD13	1:B:444:LEU:HD13	1.79	0.64
1:B:553:ILE:HB	1:B:612:ARG:HE	1.60	0.64
1:C:705:ARG:O	1:C:706:PHE:HB3	1.96	0.64
1:C:1002:GLY:O	1:C:1005:VAL:HB	1.96	0.64
1:B:254:VAL:HG22	1:B:436:THR:CG2	2.28	0.64
1:B:512:GLU:OE1	1:B:630:THR:HB	1.97	0.64
1:C:1012:ARG:C	1:C:1013:TYR:HD1	2.06	0.64
1:B:465:ARG:HH22	1:B:864:GLU:CD	2.06	0.64
1:B:197:LEU:HA	1:B:200:LEU:HD12	1.80	0.64
1:C:254:VAL:HG22	1:C:436:THR:CG2	2.26	0.64
1:C:220:LEU:HD12	1:C:221:MET:H	1.62	0.64
1:C:756:LEU:O	1:C:940:THR:HA	1.98	0.64
1:C:109:SER:HB3	1:C:138:HIS:CD2	2.33	0.64
1:C:1037:LEU:HD23	1:C:1038:ASP:N	2.13	0.64
1:B:86:SER:N	2:B:2055:ANP:O1A	2.30	0.64
1:B:374:TYR:C	1:B:376:TYR:H	2.04	0.64
1:C:190:ALA:HB3	1:C:193:ASN:ND2	2.12	0.64
1:C:63:MET:O	1:C:63:MET:HG2	1.98	0.64
1:C:109:SER:HB3	1:C:138:HIS:HD2	1.62	0.64
1:C:224:THR:OG1	1:C:247:SER:HB3	1.98	0.64
1:B:529:VAL:CG1	1:B:530:LYS:N	2.61	0.64
1:B:706:PHE:O	1:B:706:PHE:CD1	2.51	0.64
1:B:738:ALA:HA	1:B:947:VAL:HA	1.79	0.64
1:B:50:PHE:CD2	1:B:57:PRO:HG3	2.33	0.63
1:C:264:ILE:HD13	1:C:264:ILE:C	2.23	0.63
1:C:194:VAL:O	1:C:198:LEU:HD12	1.97	0.63
1:C:718:THR:O	1:C:721:LEU:HB3	1.99	0.63
1:B:570:VAL:HG21	1:B:1047:TYR:OH	1.98	0.63
1:B:681:VAL:O	1:B:685:VAL:HG23	1.98	0.63
1:B:830:GLY:O	1:B:831:ASP:HB2	1.97	0.63
1:B:901:LYS:HG3	1:B:902:TYR:H	1.64	0.63
1:B:901:LYS:HG3	1:B:902:TYR:N	2.12	0.63
1:B:493:ASP:HA	1:B:496:ARG:NH2	2.14	0.63
1:B:883:ILE:H	1:B:883:ILE:HD12	1.62	0.63
1:B:709:ARG:O	1:B:710:ASN:HB2	1.98	0.63
1:B:1002:GLY:O	1:B:1005:VAL:HB	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:530:LYS:O	1:C:537:VAL:HG12	1.98	0.63
1:C:477:ASP:C	1:C:479:VAL:H	2.06	0.63
1:C:816:ARG:O	1:C:845:GLU:HG3	1.99	0.63
1:B:1042:ALA:O	1:B:1046:MET:HG3	1.98	0.63
1:C:267:LEU:O	1:C:270:ILE:HG22	1.99	0.63
1:C:470:ASP:O	1:C:471:ILE:C	2.41	0.63
1:C:560:ILE:HG13	1:C:561:THR:N	2.14	0.63
1:C:573:ARG:O	1:C:574:PHE:HB2	1.97	0.63
1:C:685:VAL:O	1:C:688:ARG:HB3	1.99	0.62
1:B:164:GLN:HE21	1:B:164:GLN:C	2.07	0.62
1:B:610:ARG:HH11	1:B:610:ARG:HB3	1.64	0.62
1:B:620:ALA:HB2	1:B:647:LEU:HD13	1.81	0.62
1:B:858:VAL:HG12	1:B:862:ILE:HD11	1.81	0.62
1:C:345:PHE:HA	1:C:436:THR:OG1	1.99	0.62
1:C:440:PHE:CD2	1:C:495:TYR:HA	2.34	0.62
1:C:735:ARG:O	1:C:949:SER:HA	1.99	0.62
1:B:55:GLY:O	1:B:57:PRO:HD3	2.00	0.62
1:B:103:TYR:HB3	1:B:179:ILE:HG12	1.81	0.62
1:B:207:LEU:N	1:B:207:LEU:HD23	2.14	0.62
1:B:472:GLU:OE2	1:B:472:GLU:HA	1.98	0.62
1:B:560:ILE:HD12	1:B:579:ALA:CB	2.29	0.62
1:C:895:PHE:H	1:C:895:PHE:HD2	1.48	0.62
1:B:426:LEU:HD12	1:B:465:ARG:HG3	1.80	0.62
1:B:1007:ARG:HA	1:B:1010:VAL:HG12	1.81	0.62
1:C:358:ILE:HD11	1:C:419:GLY:C	2.25	0.62
1:C:363:SER:O	1:C:364:LEU:HD12	2.00	0.62
1:C:398:GLU:HG3	1:C:399:ILE:HD12	1.81	0.62
1:C:855:ARG:HG3	1:C:877:PHE:HB3	1.81	0.62
1:B:267:LEU:O	1:B:270:ILE:HG22	1.99	0.62
1:C:883:ILE:HD12	1:C:883:ILE:H	1.63	0.62
1:B:384:ARG:O	1:B:387:PRO:HB3	1.99	0.62
1:C:37:PHE:HB2	1:C:621:HIS:CE1	2.34	0.62
1:C:849:PRO:HG2	1:C:884:PHE:HA	1.82	0.62
1:B:387:PRO:HB2	1:B:389:VAL:H	1.64	0.62
1:B:756:LEU:O	1:B:940:THR:HG23	2.00	0.62
1:C:154:ARG:HE	1:C:174:GLY:HA3	1.63	0.62
1:C:249:ILE:HG23	1:C:541:PRO:HG2	1.81	0.62
1:B:972:ARG:NH1	1:B:1026:GLU:HG3	2.14	0.62
1:C:1007:ARG:HA	1:C:1010:VAL:CG1	2.30	0.61
1:B:212:TRP:CZ2	1:B:235:LEU:HD11	2.35	0.61
1:B:721:LEU:O	1:B:725:ILE:HG12	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:740:VAL:HG12	1:B:741:ARG:N	2.14	0.61
1:B:904:ILE:HD12	1:B:911:ALA:H	1.65	0.61
1:B:975:THR:HG22	1:B:979:ILE:HG12	1.80	0.61
1:B:1009:LEU:HD22	1:B:1009:LEU:N	2.15	0.61
1:C:238:GLN:HG3	1:C:239:LEU:HD23	1.82	0.61
1:C:244:ILE:CG2	1:C:245:GLY:N	2.57	0.61
1:C:981:ASP:O	1:C:982:ARG:C	2.43	0.61
1:B:140:ARG:O	1:B:141:ILE:O	2.18	0.61
1:B:333:LEU:HD22	1:B:333:LEU:N	2.14	0.61
1:B:774:TYR:CE2	1:B:886:ARG:HG3	2.35	0.61
1:C:101:ARG:HH22	1:C:174:GLY:HA3	1.65	0.61
1:C:374:TYR:C	1:C:376:TYR:H	2.07	0.61
1:C:556:VAL:HB	1:C:613:ILE:CD1	2.29	0.61
1:C:743:PHE:C	1:C:745:LEU:H	2.08	0.61
1:B:190:ALA:HB3	1:B:193:ASN:ND2	2.16	0.61
1:B:314:LYS:HB3	1:B:320:ILE:HG12	1.82	0.61
1:B:957:THR:O	1:B:958:GLN:C	2.44	0.61
1:B:1010:VAL:HG13	1:B:1011:ARG:N	2.15	0.61
1:B:1049:GLU:O	1:B:1052:SER:HB3	1.99	0.61
1:C:212:TRP:CZ2	1:C:235:LEU:HD11	2.35	0.61
1:C:740:VAL:HG12	1:C:741:ARG:N	2.16	0.61
1:B:393:ILE:HG22	1:B:397:ARG:NE	2.14	0.61
1:B:612:ARG:HG3	1:B:612:ARG:NH1	2.14	0.61
1:B:690:GLU:HG3	1:B:694:ILE:HG13	1.83	0.61
1:C:852:PRO:O	1:C:853:LEU:HD23	2.01	0.61
1:B:1009:LEU:H	1:B:1009:LEU:CD2	2.12	0.61
1:C:731:SER:O	1:C:951:PRO:HB3	2.00	0.61
1:B:968:ARG:HD2	1:B:1019:GLU:OE1	2.00	0.61
1:C:266:THR:O	1:C:267:LEU:HB2	1.99	0.61
1:C:1010:VAL:HG13	1:C:1011:ARG:N	2.15	0.61
1:B:1025:LEU:O	1:B:1025:LEU:HD23	2.01	0.61
1:C:61:GLN:HE22	2:C:2055:ANP:HN61	1.47	0.61
1:C:849:PRO:HD3	1:C:883:ILE:HG22	1.83	0.61
1:C:705:ARG:O	1:C:706:PHE:CB	2.49	0.61
1:B:176:PHE:HB2	1:B:179:ILE:HD11	1.83	0.60
1:B:849:PRO:HG2	1:B:884:PHE:HA	1.82	0.60
1:B:70:ARG:NH2	1:B:618:LYS:HD2	2.15	0.60
1:B:849:PRO:HD2	1:B:887:PHE:HD2	1.66	0.60
1:B:771:LEU:HB3	1:B:886:ARG:NH1	2.16	0.60
1:C:494:ARG:O	1:C:497:ARG:N	2.34	0.60
1:C:687:ARG:HA	1:C:1029:MET:CE	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:63:MET:HG3	1:B:622:ASP:HB3	1.83	0.60
1:B:800:GLU:OE2	1:B:876:HIS:CE1	2.54	0.60
1:B:612:ARG:HG3	1:B:612:ARG:HH11	1.65	0.60
1:C:65:ALA:HA	1:C:91:MET:HE2	1.84	0.60
1:C:463:ILE:HD13	1:C:474:LYS:HG3	1.83	0.60
1:C:739:ILE:O	1:C:945:ALA:HA	2.01	0.60
1:B:709:ARG:HH11	1:B:709:ARG:HG3	1.66	0.60
1:B:735:ARG:O	1:B:949:SER:HA	2.02	0.60
1:B:301:ILE:HD12	1:B:301:ILE:N	2.14	0.60
1:B:560:ILE:HG13	1:B:561:THR:H	1.66	0.60
1:C:518:ARG:O	1:C:522:ARG:HD2	2.02	0.60
1:C:203:PHE:CD1	1:C:239:LEU:HD12	2.37	0.60
1:C:532:LEU:O	1:C:534:GLY:N	2.35	0.60
1:B:415:VAL:HG12	1:B:422:ILE:HB	1.83	0.59
1:C:63:MET:CE	1:C:67:ARG:HH21	2.15	0.59
1:B:66:LYS:O	1:B:70:ARG:HG3	2.01	0.59
1:B:463:ILE:HD13	1:B:474:LYS:HG3	1.82	0.59
1:B:532:LEU:HD12	1:B:619:LEU:HD21	1.83	0.59
1:C:357:THR:HG22	1:C:358:ILE:N	2.16	0.59
1:C:958:GLN:O	1:C:961:ILE:HG22	2.01	0.59
1:B:309:LYS:HD2	1:B:309:LYS:N	2.16	0.59
1:B:477:ASP:C	1:B:479:VAL:H	2.08	0.59
1:C:101:ARG:NH1	1:C:154:ARG:HA	2.17	0.59
1:B:260:ASN:ND2	1:B:260:ASN:O	2.35	0.59
1:B:339:LEU:HB2	1:B:343:ILE:HG12	1.85	0.59
1:B:471:ILE:HG22	1:B:471:ILE:O	2.01	0.59
1:C:56:GLU:OE1	1:C:57:PRO:HD2	2.03	0.59
1:C:609:SER:HA	1:C:612:ARG:NH1	2.17	0.59
1:B:107:PRO:HA	1:B:163:THR:HG22	1.84	0.59
1:B:257:VAL:HB	1:B:450:PHE:CD1	2.37	0.59
1:B:358:ILE:HD12	1:B:358:ILE:N	2.12	0.59
1:B:735:ARG:HB3	1:B:949:SER:CA	2.33	0.59
1:C:358:ILE:HG23	1:C:404:MET:HE1	1.85	0.59
1:B:387:PRO:CB	1:B:389:VAL:H	2.16	0.59
1:B:664:ARG:O	1:B:668:GLU:HG3	2.03	0.59
1:C:339:LEU:HB2	1:C:343:ILE:HG13	1.85	0.59
1:B:1051:LYS:C	1:B:1053:ILE:H	2.11	0.59
1:C:771:LEU:HD12	1:C:927:TYR:CE2	2.33	0.59
1:B:708:ASN:C	1:B:708:ASN:HD22	2.11	0.59
1:B:735:ARG:HB3	1:B:949:SER:HB3	1.85	0.59
1:C:459:LEU:O	1:C:463:ILE:HG12	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:848:ARG:HG3	1:C:887:PHE:CE2	2.38	0.59
1:B:376:TYR:O	1:B:377:ARG:CB	2.50	0.58
1:C:514:PRO:O	1:C:517:ALA:HB3	2.03	0.58
1:B:692:ARG:HG2	1:B:692:ARG:HH21	1.68	0.58
1:B:358:ILE:HD11	1:B:419:GLY:O	2.04	0.58
1:C:561:THR:OG1	1:C:585:ARG:HD2	2.03	0.58
1:B:440:PHE:HD1	1:B:440:PHE:H	1.52	0.58
1:B:555:HIS:O	1:B:581:ILE:HD12	2.02	0.58
1:C:760:LEU:HD21	1:C:938:THR:HB	1.85	0.58
1:B:682:LYS:HD3	1:B:1039:TYR:CD2	2.39	0.58
1:C:737:ILE:HD13	1:C:950:VAL:HG21	1.84	0.58
1:B:103:TYR:HB2	1:B:176:PHE:CE2	2.38	0.58
1:B:271:LEU:HD23	1:B:481:PHE:CZ	2.38	0.58
1:B:309:LYS:HG2	1:B:310:GLY:H	1.67	0.58
1:B:608:ASN:ND2	1:B:610:ARG:HB2	2.19	0.58
1:B:717:GLN:OE1	1:B:982:ARG:HD3	2.02	0.58
1:B:49:PHE:CD2	1:B:94:PHE:HD1	2.21	0.58
1:B:135:GLY:HA3	1:B:156:PHE:CE1	2.37	0.58
1:B:402:LYS:HG2	1:B:403:VAL:HG13	1.84	0.58
1:B:613:ILE:HG21	1:B:646:LEU:HD22	1.86	0.58
1:C:808:THR:HG23	1:C:848:ARG:HE	1.67	0.58
1:B:60:ILE:HD12	1:B:506:PRO:HG3	1.86	0.58
1:B:608:ASN:ND2	1:B:610:ARG:HH11	2.00	0.58
1:B:608:ASN:HD21	1:B:610:ARG:HB2	1.69	0.58
1:B:766:GLU:HB3	1:B:768:ARG:HE	1.69	0.58
1:B:824:ILE:HG23	1:B:852:PRO:HG3	1.85	0.58
1:B:858:VAL:O	1:B:862:ILE:HG13	2.04	0.58
1:C:509:PHE:O	1:C:549:VAL:HA	2.03	0.58
1:C:664:ARG:O	1:C:668:GLU:HG3	2.04	0.58
1:C:983:LEU:HB3	1:C:989:VAL:HG23	1.84	0.58
1:B:93:LEU:HA	1:B:158:ILE:HD11	1.85	0.57
1:B:268:SER:O	1:B:272:GLU:HG2	2.03	0.57
1:B:387:PRO:HG2	1:B:389:VAL:HA	1.86	0.57
1:B:883:ILE:HD12	1:B:883:ILE:N	2.19	0.57
1:B:901:LYS:HA	1:B:914:GLU:HA	1.85	0.57
1:B:972:ARG:HB3	1:B:973:PRO:HD2	1.85	0.57
1:C:466:ALA:HB1	1:C:472:GLU:HG3	1.85	0.57
1:C:714:GLY:O	1:C:716:ALA:N	2.37	0.57
1:B:981:ASP:O	1:B:982:ARG:C	2.46	0.57
1:C:279:ILE:O	1:C:346:ALA:HA	2.05	0.57
1:C:305:THR:CG2	1:C:329:TYR:CD1	2.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:791:VAL:H	1:C:958:GLN:NE2	2.02	0.57
1:B:386:LEU:HD23	1:B:386:LEU:N	2.03	0.57
1:B:804:ASN:O	1:B:806:LEU:HG	2.04	0.57
1:B:972:ARG:HB3	1:B:973:PRO:CD	2.34	0.57
1:C:618:LYS:O	1:C:621:HIS:HB3	2.05	0.57
1:B:365:SER:OG	1:B:368:MET:HB2	2.05	0.57
1:C:32:ARG:HD2	1:C:32:ARG:N	2.18	0.57
1:C:164:GLN:HE21	1:C:164:GLN:C	2.13	0.57
1:C:492:ARG:HG3	1:C:492:ARG:NH1	2.18	0.57
1:B:194:VAL:O	1:B:197:LEU:N	2.34	0.57
1:B:612:ARG:HH11	1:B:612:ARG:CG	2.17	0.57
1:B:775:THR:O	1:B:776:THR:C	2.47	0.57
1:C:79:PRO:HG2	1:C:504:ILE:HD13	1.86	0.57
1:C:426:LEU:O	1:C:430:ILE:HD12	2.04	0.57
1:B:970:ILE:HD11	1:B:1019:GLU:HG3	1.87	0.57
1:C:3:PRO:O	1:C:4:VAL:CB	2.53	0.57
1:B:23:GLU:O	1:B:24:LYS:CB	2.51	0.57
1:B:362:ASP:O	1:B:397:ARG:NH1	2.38	0.57
1:B:518:ARG:O	1:B:522:ARG:HD3	2.04	0.57
1:C:58:ARG:HG2	1:C:502:ASP:HB2	1.87	0.57
1:C:259:VAL:HB	1:C:452:LEU:CD2	2.34	0.57
1:C:418:GLU:HG3	1:C:418:GLU:O	2.05	0.57
1:C:532:LEU:HD12	1:C:619:LEU:HD21	1.87	0.57
1:C:758:ILE:HG22	1:C:938:THR:OG1	2.05	0.57
1:C:824:ILE:HG23	1:C:852:PRO:CG	2.34	0.57
1:B:279:ILE:HG22	1:B:281:TYR:HE1	1.67	0.57
1:B:380:ASP:HB3	1:B:383:GLU:OE2	2.05	0.57
1:C:93:LEU:HA	1:C:158:ILE:HD11	1.85	0.57
1:C:791:VAL:HG21	1:C:977:ALA:HA	1.85	0.57
1:C:968:ARG:HD2	1:C:1019:GLU:OE1	2.04	0.57
1:B:58:ARG:HB2	1:B:61:GLN:NE2	2.19	0.56
1:B:134:ILE:HD12	1:B:135:GLY:N	2.20	0.56
1:B:418:GLU:O	1:B:418:GLU:HG3	2.04	0.56
1:B:905:GLU:HA	1:B:910:THR:HA	1.87	0.56
1:C:638:LYS:HE2	1:C:642:ASP:OD2	2.05	0.56
1:B:510:ILE:CD1	1:B:616:LEU:HD22	2.36	0.56
1:B:706:PHE:O	1:B:706:PHE:HD1	1.88	0.56
1:C:687:ARG:HA	1:C:1029:MET:HE3	1.87	0.56
1:C:690:GLU:OE2	1:C:715:ARG:NH1	2.39	0.56
1:C:835:GLY:O	1:C:836:ARG:HB2	2.05	0.56
1:B:170:TYR:CZ	1:B:171:ARG:HG2	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:337:LEU:HD21	1:B:343:ILE:HD11	1.87	0.56
1:B:492:ARG:HG3	1:B:492:ARG:NH1	2.16	0.56
1:B:983:LEU:HB3	1:B:989:VAL:HG23	1.85	0.56
1:C:195:ASP:HA	1:C:198:LEU:HD13	1.87	0.56
1:C:207:LEU:N	1:C:207:LEU:HD23	2.21	0.56
1:C:703:TRP:CD1	1:C:709:ARG:HA	2.40	0.56
1:C:1012:ARG:O	1:C:1013:TYR:CB	2.52	0.56
1:B:476:ILE:HG13	1:B:476:ILE:O	2.05	0.56
1:B:124:GLU:C	1:B:125:LYS:HD2	2.30	0.56
1:B:315:PHE:HB2	1:B:323:LEU:HD11	1.87	0.56
1:B:385:LEU:O	1:B:387:PRO:HD3	2.05	0.56
1:C:74:PHE:CZ	1:C:221:MET:HE3	2.40	0.56
1:C:117:GLU:HB3	1:C:121:LYS:HZ3	1.70	0.56
1:C:249:ILE:HD12	1:C:541:PRO:CG	2.36	0.56
1:B:497:ARG:O	1:B:500:GLU:HG2	2.05	0.56
1:B:767:LEU:O	1:B:767:LEU:HD23	2.05	0.56
1:C:322:HIS:O	1:C:323:LEU:HD12	2.05	0.56
1:C:358:ILE:HG23	1:C:404:MET:SD	2.45	0.56
1:B:834:VAL:HG12	1:B:835:GLY:N	2.20	0.56
1:C:787:LEU:HD13	1:C:789:PHE:HE2	1.69	0.56
1:B:93:LEU:O	1:B:97:LEU:HG	2.05	0.56
1:B:391:ARG:O	1:B:392:HIS:HB2	2.05	0.56
1:C:860:ARG:HB3	1:C:860:ARG:HH11	1.71	0.56
1:B:197:LEU:O	1:B:201:LEU:HD12	2.05	0.56
1:B:305:THR:HG23	1:B:329:TYR:CZ	2.40	0.56
1:B:791:VAL:HG21	1:B:977:ALA:HA	1.86	0.56
1:C:476:ILE:H	1:C:478:GLU:HG2	1.71	0.56
1:C:1051:LYS:C	1:C:1053:ILE:H	2.14	0.56
1:B:67:ARG:HH11	1:B:67:ARG:HG2	1.71	0.56
1:B:377:ARG:O	1:B:378:ASN:C	2.49	0.56
1:B:433:SER:O	1:B:436:THR:HG22	2.06	0.56
1:B:47:VAL:HG13	1:B:57:PRO:HG2	1.88	0.55
1:B:344:ARG:HH12	1:B:488:LEU:HD12	1.69	0.55
1:B:736:LYS:HA	1:B:736:LYS:HZ3	1.70	0.55
1:B:936:LEU:HD12	1:B:936:LEU:O	2.06	0.55
1:C:74:PHE:HZ	1:C:221:MET:HE3	1.71	0.55
1:B:4:VAL:CB	1:B:32:ARG:HH21	2.19	0.55
1:B:179:ILE:O	1:B:220:LEU:HD12	2.05	0.55
1:B:305:THR:O	1:B:307:THR:N	2.40	0.55
1:B:743:PHE:C	1:B:745:LEU:H	2.13	0.55
1:C:426:LEU:HD11	1:C:466:ALA:CA	2.29	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:830:GLY:O	1:C:831:ASP:HB2	2.05	0.55
1:C:176:PHE:HB2	1:C:179:ILE:CD1	2.36	0.55
1:C:462:PHE:O	1:C:463:ILE:C	2.49	0.55
1:C:465:ARG:O	1:C:465:ARG:HD3	2.07	0.55
1:B:80:THR:HG21	1:B:438:ARG:NH1	2.21	0.55
1:B:559:LEU:HD11	1:B:689:ILE:HD13	1.88	0.55
1:C:63:MET:HG3	1:C:622:ASP:HB3	1.88	0.55
1:C:532:LEU:O	1:C:533:ASP:C	2.49	0.55
1:C:706:PHE:HD1	1:C:706:PHE:O	1.90	0.55
1:C:717:GLN:OE1	1:C:982:ARG:HD3	2.05	0.55
1:C:936:LEU:HD12	1:C:936:LEU:O	2.06	0.55
1:B:266:THR:O	1:B:267:LEU:HB2	2.05	0.55
1:B:555:HIS:HA	1:B:639:ILE:HD13	1.88	0.55
1:B:760:LEU:H	1:B:760:LEU:CD2	2.14	0.55
1:B:975:THR:HG22	1:B:979:ILE:CG1	2.36	0.55
1:B:79:PRO:HG2	1:B:504:ILE:HG21	1.88	0.55
1:B:692:ARG:HG2	1:B:692:ARG:NH2	2.21	0.55
1:B:961:ILE:HD13	1:B:961:ILE:O	2.06	0.55
1:C:274:LEU:HD22	1:C:450:PHE:HE2	1.70	0.55
1:C:508:LEU:HD21	1:C:510:ILE:HD11	1.88	0.55
1:C:757:THR:O	1:C:904:ILE:HA	2.06	0.55
1:C:784:ASN:HB3	1:C:959:SER:HB2	1.88	0.55
1:C:883:ILE:HD12	1:C:883:ILE:N	2.21	0.55
1:B:106:PHE:O	1:B:163:THR:N	2.39	0.55
1:B:608:ASN:ND2	1:B:610:ARG:CB	2.70	0.55
1:C:61:GLN:OE1	1:C:83:GLY:HA3	2.07	0.55
1:C:433:SER:O	1:C:436:THR:HG22	2.07	0.55
1:C:235:LEU:O	1:C:239:LEU:HB2	2.07	0.55
1:C:540:ILE:HG13	1:C:541:PRO:CD	2.35	0.55
1:C:616:LEU:HA	1:C:619:LEU:HD12	1.89	0.55
1:C:662:THR:O	1:C:663:ARG:C	2.50	0.55
1:C:774:TYR:CD2	1:C:886:ARG:HG2	2.42	0.55
1:C:848:ARG:CB	1:C:887:PHE:CD2	2.90	0.55
1:C:1012:ARG:C	1:C:1013:TYR:CD1	2.85	0.55
1:B:965:MET:CE	1:B:980:VAL:HG23	2.37	0.55
1:C:615:ALA:O	1:C:618:LYS:HB3	2.07	0.55
1:C:735:ARG:HB3	1:C:949:SER:CB	2.36	0.55
1:C:138:HIS:NE2	1:C:140:ARG:HB2	2.21	0.55
1:C:476:ILE:C	1:C:478:GLU:H	2.15	0.55
1:B:366:PRO:HA	1:B:393:ILE:HG12	1.87	0.54
1:B:376:TYR:OH	1:B:409:PRO:HD3	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:714:GLY:O	1:B:715:ARG:HB3	2.07	0.54
1:C:194:VAL:HG12	1:C:198:LEU:HD11	1.88	0.54
1:C:616:LEU:HD23	1:C:619:LEU:HD12	1.89	0.54
1:B:134:ILE:HD12	1:B:135:GLY:H	1.72	0.54
1:B:171:ARG:HG3	1:B:171:ARG:HH11	1.73	0.54
1:B:248:ARG:CG	1:B:249:ILE:N	2.70	0.54
1:B:558:ASP:CG	1:B:688:ARG:HH21	2.15	0.54
1:B:570:VAL:O	1:B:571:ASN:HB2	2.08	0.54
1:C:47:VAL:HG12	1:C:51:ARG:NH1	2.22	0.54
1:C:198:LEU:O	1:C:201:LEU:N	2.37	0.54
1:B:153:LEU:O	1:B:154:ARG:HB2	2.07	0.54
1:B:416:VAL:HG13	1:B:417:ARG:N	2.22	0.54
1:B:801:LEU:HB3	1:B:807:ILE:CD1	2.38	0.54
1:C:3:PRO:HA	1:C:16:ASP:CB	2.38	0.54
1:C:775:THR:O	1:C:776:THR:C	2.50	0.54
1:B:61:GLN:NE2	2:B:2055:ANP:N6	2.40	0.54
1:B:64:TRP:O	1:B:67:ARG:HB2	2.07	0.54
1:B:638:LYS:HE2	1:B:642:ASP:OD2	2.07	0.54
1:B:749:HIS:HD2	1:B:751:GLU:HB2	1.72	0.54
1:B:849:PRO:HD2	1:B:887:PHE:CD2	2.42	0.54
1:B:972:ARG:CZ	1:B:1026:GLU:HG3	2.38	0.54
1:C:687:ARG:NH2	1:C:972:ARG:HB2	2.23	0.54
1:B:735:ARG:NH1	1:B:951:PRO:HG3	2.23	0.54
1:B:828:TYR:HD2	1:B:829:LEU:HD23	1.72	0.54
1:C:194:VAL:O	1:C:195:ASP:C	2.50	0.54
1:B:344:ARG:NH1	1:B:488:LEU:HD12	2.23	0.54
1:B:377:ARG:HG3	1:B:382:ILE:HG13	1.88	0.54
1:C:170:TYR:CZ	1:C:171:ARG:HG2	2.42	0.54
1:C:542:MET:O	1:C:544:LYS:N	2.40	0.54
1:B:455:ASP:O	1:B:456:SER:C	2.51	0.54
1:B:687:ARG:HA	1:B:1029:MET:HE3	1.89	0.54
1:B:1012:ARG:HH11	1:B:1012:ARG:CG	2.20	0.54
1:C:43:LEU:HD23	1:C:43:LEU:C	2.33	0.54
1:C:66:LYS:HD2	1:C:622:ASP:HA	1.88	0.54
1:B:194:VAL:HG12	1:B:195:ASP:N	2.22	0.54
1:B:532:LEU:O	1:B:533:ASP:C	2.49	0.54
1:B:809:TYR:CZ	1:B:811:ARG:HB2	2.43	0.54
1:C:330:TYR:O	1:C:333:LEU:HD23	2.08	0.54
1:C:351:CYS:HB3	1:C:462:PHE:CD1	2.43	0.54
1:C:556:VAL:HG23	1:C:557:VAL:HG13	1.89	0.54
1:C:970:ILE:HD11	1:C:1019:GLU:HG3	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:682:LYS:HD3	1:B:1039:TYR:HD2	1.73	0.54
1:B:700:GLN:HA	1:B:703:TRP:CE3	2.43	0.54
1:B:935:GLU:HG2	1:B:936:LEU:N	2.23	0.54
1:C:381:GLU:O	1:C:384:ARG:HG2	2.08	0.54
1:C:468:LEU:HD22	1:C:468:LEU:H	1.73	0.54
1:B:137:TYR:CD1	1:B:137:TYR:C	2.85	0.53
1:C:80:THR:HG21	1:C:438:ARG:NH1	2.23	0.53
1:C:344:ARG:O	1:C:345:PHE:HB3	2.08	0.53
1:C:991:GLU:HG3	1:C:995:ARG:O	2.08	0.53
1:B:37:PHE:O	1:B:40:ASP:HB2	2.09	0.53
1:B:358:ILE:H	1:B:358:ILE:CD1	2.08	0.53
1:B:469:TYR:O	1:B:470:ASP:C	2.51	0.53
1:B:1022:THR:O	1:B:1026:GLU:HB2	2.08	0.53
1:C:824:ILE:CG2	1:C:852:PRO:HG3	2.36	0.53
1:C:860:ARG:HH11	1:C:860:ARG:CG	2.21	0.53
1:B:58:ARG:NH2	1:B:81:GLY:O	2.41	0.53
1:B:102:CYS:C	1:B:176:PHE:HD2	2.16	0.53
1:B:153:LEU:O	1:B:154:ARG:CB	2.56	0.53
1:B:542:MET:HE1	1:B:670:LEU:HD12	1.89	0.53
1:B:854:THR:O	1:B:858:VAL:HG23	2.09	0.53
1:C:255:GLU:OE1	1:C:484:LEU:HD11	2.09	0.53
1:C:736:LYS:C	1:C:737:ILE:CG2	2.81	0.53
1:C:348:PHE:HZ	1:C:432:GLY:HA3	1.74	0.53
1:C:479:VAL:HG23	1:C:481:PHE:H	1.74	0.53
1:B:101:ARG:HA	1:B:157:LYS:O	2.07	0.53
1:B:256:ASP:CB	1:B:474:LYS:HD3	2.39	0.53
1:C:88:GLY:O	1:C:92:SER:HB2	2.09	0.53
1:B:361:ILE:CG2	1:B:397:ARG:HB3	2.38	0.53
1:C:683:ALA:O	1:C:687:ARG:HB2	2.07	0.53
1:C:887:PHE:C	1:C:889:ALA:H	2.16	0.53
1:C:1009:LEU:H	1:C:1009:LEU:HD22	1.72	0.53
1:C:251:VAL:O	1:C:252:ARG:HB3	2.07	0.53
1:C:859:GLN:O	1:C:863:GLN:HG2	2.09	0.53
1:C:784:ASN:CG	1:C:958:GLN:HB2	2.34	0.53
1:B:399:ILE:O	1:B:403:VAL:HG22	2.08	0.53
1:C:468:LEU:HD22	1:C:468:LEU:N	2.24	0.53
1:C:609:SER:HA	1:C:612:ARG:HH11	1.73	0.53
1:C:715:ARG:HG2	1:C:716:ALA:N	2.24	0.53
1:C:725:ILE:C	1:C:727:ARG:H	2.17	0.53
1:C:736:LYS:N	1:C:736:LYS:HD2	2.23	0.53
1:B:100:LYS:HG2	1:B:177:ASP:CG	2.34	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:720:VAL:HG13	1:B:961:ILE:HD11	1.92	0.52
1:B:1000:LYS:O	1:B:1001:LEU:C	2.52	0.52
1:B:1029:MET:O	1:B:1032:ILE:HG23	2.09	0.52
1:C:34:LEU:HD13	1:C:648:SER:HB3	1.91	0.52
1:C:935:GLU:HG2	1:C:936:LEU:N	2.22	0.52
1:B:207:LEU:O	1:B:210:LYS:HG3	2.09	0.52
1:B:555:HIS:HA	1:B:639:ILE:HD11	1.88	0.52
1:C:463:ILE:O	1:C:466:ALA:HB3	2.09	0.52
1:B:895:PHE:N	1:B:895:PHE:CD2	2.77	0.52
1:C:192:LYS:O	1:C:193:ASN:C	2.51	0.52
1:C:869:VAL:HG21	1:C:872:LEU:HD12	1.91	0.52
1:C:981:ASP:O	1:C:984:PHE:N	2.42	0.52
1:B:74:PHE:CZ	1:B:221:MET:HG3	2.44	0.52
1:B:763:GLU:CD	1:B:900:LYS:HE2	2.35	0.52
1:B:981:ASP:O	1:B:984:PHE:N	2.36	0.52
1:C:414:VAL:HG12	1:C:424:PRO:CD	2.39	0.52
1:C:480:ASP:O	1:C:481:PHE:C	2.51	0.52
1:B:314:LYS:CB	1:B:320:ILE:HG12	2.39	0.52
1:B:397:ARG:O	1:B:401:LYS:HG3	2.10	0.52
1:B:493:ASP:HA	1:B:496:ARG:CZ	2.39	0.52
1:C:107:PRO:HG3	1:C:183:ASP:HB3	1.91	0.52
1:C:314:LYS:HB3	1:C:320:ILE:HD13	1.92	0.52
1:C:818:SER:O	1:C:821:GLY:N	2.41	0.52
1:B:305:THR:HG23	1:B:329:TYR:CD1	2.44	0.52
1:B:309:LYS:HG2	1:B:310:GLY:N	2.25	0.52
1:B:466:ALA:HB1	1:B:472:GLU:HG3	1.90	0.52
1:C:364:LEU:HD22	1:C:368:MET:SD	2.50	0.52
1:C:425:ASP:OD2	1:C:428:THR:HB	2.09	0.52
1:C:446:LYS:HD3	1:C:447:GLY:H	1.73	0.52
1:C:628:VAL:HG21	1:C:640:ALA:HA	1.91	0.52
1:B:207:LEU:HD23	1:B:207:LEU:H	1.75	0.52
1:B:462:PHE:O	1:B:463:ILE:C	2.51	0.52
1:B:463:ILE:CD1	1:B:474:LYS:HG3	2.40	0.52
1:B:708:ASN:C	1:B:708:ASN:ND2	2.68	0.52
1:B:709:ARG:HG2	1:B:710:ASN:N	2.22	0.52
1:B:1028:ARG:O	1:B:1031:ALA:HB3	2.09	0.52
1:C:265:SER:O	1:C:267:LEU:N	2.42	0.52
1:C:276:THR:HA	1:C:322:HIS:CE1	2.45	0.52
1:C:296:LYS:HA	1:C:301:ILE:CG1	2.38	0.52
1:C:957:THR:O	1:C:958:GLN:C	2.53	0.52
1:B:212:TRP:CH2	1:B:235:LEU:HD21	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:697:VAL:HG21	1:B:1050:ILE:HD12	1.92	0.52
1:B:1010:VAL:CG1	1:B:1011:ARG:N	2.73	0.52
1:B:1029:MET:HA	1:B:1032:ILE:CG2	2.40	0.52
1:C:123:ALA:C	1:C:125:LYS:H	2.18	0.52
1:C:712:SER:CB	1:C:986:ARG:HH11	2.22	0.52
1:B:291:ILE:O	1:B:294:SER:N	2.43	0.52
1:B:292:TYR:CD2	1:B:303:ILE:HB	2.44	0.52
1:B:760:LEU:HD22	1:B:938:THR:OG1	2.10	0.52
1:B:897:VAL:HG13	1:B:897:VAL:O	2.10	0.52
1:C:207:LEU:HG	1:C:208:LYS:N	2.23	0.52
1:C:231:LYS:O	1:C:234:GLU:HB2	2.09	0.52
1:C:460:SER:O	1:C:463:ILE:HB	2.10	0.52
1:C:808:THR:CG2	1:C:848:ARG:NE	2.73	0.52
1:B:241:ASN:O	1:B:530:LYS:HD2	2.10	0.52
1:B:384:ARG:HH22	1:C:497:ARG:HD3	1.75	0.52
1:C:60:ILE:HD12	1:C:506:PRO:HG3	1.91	0.52
1:C:160:ILE:O	1:C:161:THR:HG22	2.10	0.52
1:C:464:GLU:O	1:C:467:LYS:HB2	2.10	0.52
1:B:182:ASP:O	1:B:223:SER:HB3	2.09	0.51
1:B:271:LEU:C	1:B:273:LYS:H	2.19	0.51
1:B:523:PHE:CE2	1:B:663:ARG:N	2.78	0.51
1:B:529:VAL:CG1	1:B:530:LYS:H	2.22	0.51
1:B:832:ASP:HA	1:B:893:ARG:HG3	1.91	0.51
1:C:873:ARG:N	1:C:876:HIS:HD2	2.09	0.51
1:B:653:VAL:O	1:B:653:VAL:HG12	2.09	0.51
1:B:662:THR:O	1:B:663:ARG:C	2.53	0.51
1:C:113:ILE:HD11	1:C:138:HIS:CE1	2.44	0.51
1:C:117:GLU:HB3	1:C:121:LYS:NZ	2.25	0.51
1:C:135:GLY:HA3	1:C:156:PHE:CD1	2.45	0.51
1:C:498:ARG:NH1	2:C:2055:ANP:O1G	2.43	0.51
1:B:476:ILE:O	1:B:477:ASP:HB2	2.11	0.51
1:C:66:LYS:O	1:C:70:ARG:HG3	2.10	0.51
1:C:438:ARG:CG	1:C:439:LEU:N	2.74	0.51
1:C:709:ARG:O	1:C:710:ASN:HB2	2.10	0.51
1:B:553:ILE:O	1:B:553:ILE:HG22	2.11	0.51
1:B:832:ASP:O	1:B:893:ARG:HG3	2.09	0.51
1:C:58:ARG:HH21	1:C:83:GLY:H	1.58	0.51
1:B:14:GLY:HA2	1:B:21:GLU:O	2.10	0.51
1:B:65:ALA:O	1:B:66:LYS:C	2.53	0.51
1:B:76:ALA:CB	1:B:221:MET:HE2	2.36	0.51
1:B:573:ARG:O	1:B:574:PHE:HB2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:910:THR:HG23	1:B:910:THR:O	2.10	0.51
1:C:137:TYR:CD1	1:C:137:TYR:C	2.88	0.51
1:C:872:LEU:HA	1:C:876:HIS:CD2	2.46	0.51
1:B:370:LYS:O	1:B:373:ALA:CB	2.55	0.51
1:B:581:ILE:HG22	1:B:582:LYS:N	2.26	0.51
1:C:22:ILE:O	1:C:575:VAL:HG13	2.11	0.51
1:C:637:GLU:OE1	1:C:637:GLU:HA	2.11	0.51
1:C:736:LYS:O	1:C:737:ILE:HG23	2.09	0.51
1:C:795:MET:HE1	1:C:811:ARG:HD2	1.91	0.51
1:B:708:ASN:ND2	1:B:709:ARG:O	2.43	0.51
1:B:907:ASP:O	1:B:909:LYS:HG2	2.10	0.51
1:C:816:ARG:HA	1:C:838:TRP:O	2.11	0.51
1:B:43:LEU:O	1:B:46:PHE:N	2.44	0.51
1:C:171:ARG:HH11	1:C:171:ARG:HG3	1.76	0.51
1:B:419:GLY:O	1:B:420:GLU:CB	2.58	0.51
1:B:715:ARG:HG2	1:B:716:ALA:N	2.26	0.51
1:B:757:THR:O	1:B:904:ILE:HA	2.11	0.51
1:B:784:ASN:HB3	1:B:959:SER:HB2	1.92	0.51
1:B:906:PHE:O	1:B:907:ASP:C	2.54	0.51
1:C:940:THR:O	1:C:940:THR:HG22	2.10	0.51
1:C:512:GLU:OE1	1:C:630:THR:HB	2.10	0.51
1:C:1053:ILE:HG13	1:C:1054:ASP:N	2.26	0.51
1:B:387:PRO:HB2	1:B:389:VAL:N	2.27	0.50
1:B:682:LYS:HB2	1:B:1032:ILE:HD11	1.93	0.50
1:C:719:LEU:HD21	1:C:1017:VAL:O	2.11	0.50
1:C:821:GLY:HA3	1:C:848:ARG:HH22	1.76	0.50
1:B:58:ARG:HH21	1:B:83:GLY:H	1.57	0.50
1:B:682:LYS:CB	1:B:1032:ILE:HD11	2.42	0.50
1:B:791:VAL:HG23	1:B:977:ALA:HB1	1.93	0.50
1:B:807:ILE:N	1:B:850:THR:HG23	2.26	0.50
1:C:352:PRO:O	1:C:425:ASP:HB3	2.11	0.50
1:C:466:ALA:HB1	1:C:472:GLU:CG	2.41	0.50
1:C:627:ILE:HD13	1:C:673:LEU:HD21	1.93	0.50
1:C:824:ILE:HD12	1:C:852:PRO:HD3	1.93	0.50
1:C:1017:VAL:O	1:C:1017:VAL:HG12	2.11	0.50
1:B:754:PHE:CE2	1:B:906:PHE:HB2	2.47	0.50
1:C:55:GLY:O	1:C:57:PRO:HD3	2.11	0.50
1:C:134:ILE:HD13	1:C:134:ILE:C	2.36	0.50
1:C:248:ARG:HG3	1:C:249:ILE:N	2.26	0.50
1:C:254:VAL:HG22	1:C:436:THR:HG23	1.92	0.50
1:C:258:ALA:HA	1:C:451:LEU:O	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:438:ARG:C	1:C:444:LEU:HD12	2.35	0.50
1:C:874:TRP:CG	1:C:875:GLU:N	2.79	0.50
1:B:170:TYR:CG	1:B:171:ARG:N	2.79	0.50
1:B:322:HIS:C	1:B:323:LEU:HD12	2.35	0.50
1:B:907:ASP:C	1:B:909:LYS:H	2.20	0.50
1:B:1052:SER:O	1:B:1053:ILE:HG23	2.11	0.50
1:C:278:GLY:O	1:C:279:ILE:HD13	2.11	0.50
1:C:533:ASP:OD1	1:C:533:ASP:N	2.45	0.50
1:C:735:ARG:HB3	1:C:949:SER:CA	2.41	0.50
1:B:184:VAL:O	1:B:188:LEU:HG	2.12	0.50
1:C:52:LYS:NZ	1:C:126:ALA:HB3	2.27	0.50
1:C:105:ILE:HD12	1:C:181:VAL:HG22	1.93	0.50
1:C:879:LEU:O	1:C:880:TYR:C	2.53	0.50
1:B:95:LEU:HD13	1:B:100:LYS:CD	2.39	0.50
1:B:330:TYR:O	1:B:333:LEU:HD23	2.12	0.50
1:C:87:PHE:O	1:C:88:GLY:C	2.54	0.50
1:C:103:TYR:HE1	1:C:161:THR:HG23	1.76	0.50
1:C:490:GLU:O	1:C:493:ASP:N	2.42	0.50
1:B:101:ARG:HH22	1:B:174:GLY:HA3	1.76	0.50
1:C:570:VAL:O	1:C:571:ASN:HB2	2.12	0.50
1:C:734:ARG:HB2	1:C:952:LYS:HA	1.93	0.50
1:C:947:VAL:O	1:C:948:LYS:HG3	2.11	0.50
1:B:344:ARG:NH1	1:B:488:LEU:HB2	2.27	0.50
1:C:261:ASP:CG	1:C:262:GLU:H	2.20	0.50
1:C:329:TYR:O	1:C:332:THR:N	2.44	0.50
1:C:745:LEU:HD22	1:C:747:LEU:HD11	1.94	0.50
1:B:138:HIS:NE2	1:B:140:ARG:HB2	2.27	0.50
1:B:735:ARG:HB3	1:B:949:SER:CB	2.41	0.50
1:B:1050:ILE:O	1:B:1053:ILE:HG12	2.12	0.50
1:C:64:TRP:O	1:C:65:ALA:C	2.55	0.50
1:C:720:VAL:O	1:C:724:ILE:HG13	2.12	0.50
1:B:874:TRP:CG	1:B:875:GLU:H	2.29	0.49
1:C:56:GLU:OE1	1:C:56:GLU:HA	2.12	0.49
1:C:58:ARG:N	1:C:61:GLN:HE21	2.04	0.49
1:C:791:VAL:H	1:C:958:GLN:HE22	1.58	0.49
1:C:806:LEU:C	1:C:850:THR:HG23	2.36	0.49
1:B:128:VAL:HG12	1:B:130:THR:H	1.76	0.49
1:B:884:PHE:C	1:B:884:PHE:CD2	2.90	0.49
1:C:106:PHE:CE2	1:C:115:ALA:HB2	2.48	0.49
1:C:747:LEU:HD12	1:C:747:LEU:H	1.77	0.49
1:C:760:LEU:CD2	1:C:938:THR:OG1	2.59	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:181:VAL:HB	1:B:222:VAL:HG12	1.94	0.49
1:B:254:VAL:HG22	1:B:436:THR:HG22	1.93	0.49
1:B:383:GLU:OE1	1:C:497:ARG:NH2	2.45	0.49
1:C:712:SER:HB3	1:C:986:ARG:HG3	1.93	0.49
1:B:342:ARG:O	1:B:343:ILE:C	2.56	0.49
1:B:391:ARG:HD3	1:B:391:ARG:C	2.34	0.49
1:B:465:ARG:HA	1:B:468:LEU:HD23	1.94	0.49
1:C:254:VAL:HA	1:C:447:GLY:C	2.37	0.49
1:C:524:PHE:HB2	1:C:540:ILE:HG12	1.94	0.49
1:C:870:GLU:O	1:C:872:LEU:HG	2.12	0.49
1:C:874:TRP:CG	1:C:875:GLU:H	2.30	0.49
1:B:374:TYR:C	1:B:376:TYR:N	2.70	0.49
1:B:397:ARG:O	1:B:400:LEU:HB2	2.13	0.49
1:B:518:ARG:HD3	1:B:538:TYR:OH	2.13	0.49
1:B:874:TRP:CG	1:B:875:GLU:N	2.80	0.49
1:C:98:LYS:O	1:C:99:GLY:C	2.54	0.49
1:C:760:LEU:HD23	1:C:938:THR:OG1	2.11	0.49
1:C:808:THR:HG23	1:C:848:ARG:NE	2.27	0.49
1:C:61:GLN:NE2	2:C:2055:ANP:HN61	2.10	0.49
1:C:103:TYR:HB2	1:C:176:PHE:CE2	2.48	0.49
1:C:249:ILE:HD12	1:C:541:PRO:HG3	1.94	0.49
1:C:330:TYR:HA	1:C:333:LEU:HD23	1.94	0.49
1:B:212:TRP:C	1:B:213:VAL:HG22	2.38	0.49
1:B:298:LYS:O	1:B:299:PHE:CD1	2.65	0.49
1:B:304:VAL:O	1:B:305:THR:CB	2.60	0.49
1:B:480:ASP:O	1:B:481:PHE:C	2.55	0.49
1:B:725:ILE:C	1:B:727:ARG:H	2.21	0.49
1:B:802:PHE:CD1	1:B:809:TYR:HA	2.47	0.49
1:B:828:TYR:CB	1:B:888:MET:HE3	2.42	0.49
1:C:86:SER:O	1:C:87:PHE:C	2.55	0.49
1:C:124:GLU:C	1:C:125:LYS:HD2	2.38	0.49
1:C:259:VAL:O	1:C:452:LEU:HA	2.13	0.49
1:C:968:ARG:HG2	1:C:968:ARG:HH11	1.77	0.49
1:B:195:ASP:HA	1:B:198:LEU:HD13	1.95	0.49
1:B:278:GLY:H	1:B:322:HIS:CD2	2.31	0.49
1:B:320:ILE:HD13	1:B:320:ILE:N	2.27	0.49
1:B:363:SER:O	1:B:364:LEU:CD1	2.60	0.49
1:C:198:LEU:HB3	1:C:203:PHE:HB2	1.93	0.49
1:C:494:ARG:O	1:C:495:TYR:C	2.56	0.49
1:C:681:VAL:O	1:C:684:GLN:HB3	2.12	0.49
1:C:767:LEU:HD23	1:C:767:LEU:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:808:THR:CG2	1:C:848:ARG:HE	2.26	0.49
1:C:35:CYS:CB	1:C:650:CYS:N	2.76	0.49
1:C:103:TYR:HD1	1:C:159:VAL:O	1.96	0.49
1:C:337:LEU:HB3	1:C:435:ARG:HE	1.78	0.49
1:C:463:ILE:CD1	1:C:474:LYS:HG3	2.42	0.49
1:C:737:ILE:HA	1:C:747:LEU:O	2.13	0.49
1:C:834:VAL:HG21	1:C:893:ARG:HA	1.95	0.49
1:B:217:ARG:O	1:B:217:ARG:HD2	2.13	0.49
1:B:382:ILE:O	1:B:385:LEU:HB3	2.11	0.49
1:B:415:VAL:CG1	1:B:422:ILE:HB	2.43	0.49
1:B:736:LYS:HA	1:B:736:LYS:HZ2	1.78	0.49
1:B:760:LEU:HD23	1:B:760:LEU:N	2.16	0.49
1:B:774:TYR:CE1	1:B:886:ARG:HG3	2.48	0.49
1:B:821:GLY:HA3	1:B:848:ARG:NH2	2.27	0.49
1:B:31:LYS:N	1:B:32:ARG:NE	2.61	0.48
1:B:58:ARG:O	1:B:59:ALA:C	2.56	0.48
1:B:393:ILE:HD12	1:B:393:ILE:N	2.27	0.48
1:B:440:PHE:CD2	1:B:495:TYR:HA	2.47	0.48
1:B:685:VAL:O	1:B:686:VAL:C	2.56	0.48
1:B:966:LYS:O	1:B:967:GLU:C	2.55	0.48
1:B:1039:TYR:HA	1:B:1042:ALA:HB3	1.95	0.48
1:C:477:ASP:C	1:C:479:VAL:N	2.71	0.48
1:C:583:ARG:HB3	1:C:608:ASN:CG	2.37	0.48
1:C:721:LEU:O	1:C:724:ILE:HB	2.12	0.48
1:C:856:ASP:O	1:C:857:ASP:C	2.56	0.48
1:C:887:PHE:C	1:C:889:ALA:N	2.66	0.48
1:C:991:GLU:HG3	1:C:995:ARG:C	2.38	0.48
1:B:712:SER:HB2	1:B:986:ARG:HH11	1.79	0.48
1:C:79:PRO:HG2	1:C:504:ILE:HG21	1.95	0.48
1:C:197:LEU:O	1:C:200:LEU:N	2.44	0.48
1:C:351:CYS:SG	1:C:451:LEU:HD21	2.53	0.48
1:C:518:ARG:HG2	1:C:538:TYR:CE2	2.48	0.48
1:C:745:LEU:HD22	1:C:747:LEU:CD1	2.43	0.48
1:C:756:LEU:HD13	1:C:757:THR:H	1.78	0.48
1:C:1007:ARG:CA	1:C:1010:VAL:HG12	2.40	0.48
1:B:750:ASP:C	1:B:751:GLU:OE1	2.57	0.48
1:B:1007:ARG:HA	1:B:1010:VAL:CG1	2.42	0.48
1:C:208:LYS:C	1:C:210:LYS:H	2.21	0.48
1:C:613:ILE:N	1:C:613:ILE:HD12	2.28	0.48
1:B:119:ILE:O	1:B:120:ARG:C	2.55	0.48
1:B:376:TYR:HE2	1:B:407:GLU:HB3	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:387:PRO:CG	1:B:389:VAL:HA	2.43	0.48
1:B:38:PRO:C	1:B:40:ASP:H	2.20	0.48
1:B:131:GLU:O	1:B:133:LEU:N	2.46	0.48
1:B:320:ILE:CG2	1:B:322:HIS:O	2.62	0.48
1:B:361:ILE:HD11	1:B:404:MET:HE1	1.96	0.48
1:C:379:VAL:CG1	1:C:380:ASP:H	2.07	0.48
1:B:527:PRO:HB3	1:B:540:ILE:HB	1.95	0.48
1:B:562:ASN:O	1:B:563:ARG:C	2.57	0.48
1:B:735:ARG:HB3	1:B:949:SER:HA	1.94	0.48
1:B:857:ASP:O	1:B:861:LEU:HG	2.13	0.48
1:B:905:GLU:CG	1:B:910:THR:HB	2.42	0.48
1:C:207:LEU:HD23	1:C:207:LEU:H	1.77	0.48
1:C:365:SER:OG	1:C:368:MET:HB2	2.13	0.48
1:B:291:ILE:O	1:B:292:TYR:C	2.56	0.48
1:B:621:HIS:HD2	1:B:622:ASP:OD2	1.97	0.48
1:C:122:TYR:O	1:C:125:LYS:HD3	2.14	0.48
1:C:476:ILE:C	1:C:478:GLU:N	2.70	0.48
1:B:87:PHE:O	1:B:88:GLY:C	2.56	0.48
1:B:669:ALA:C	1:B:671:GLU:H	2.20	0.48
1:C:274:LEU:HD23	1:C:345:PHE:CD2	2.49	0.48
1:C:358:ILE:HG23	1:C:404:MET:CE	2.44	0.48
1:C:666:ILE:O	1:C:667:LEU:C	2.56	0.48
1:B:456:SER:O	1:B:457:GLU:C	2.57	0.48
1:B:697:VAL:HG21	1:B:1050:ILE:CD1	2.44	0.48
1:B:754:PHE:HE2	1:B:906:PHE:HB2	1.77	0.48
1:B:828:TYR:CD2	1:B:829:LEU:HD23	2.49	0.48
1:C:194:VAL:O	1:C:197:LEU:HB2	2.14	0.48
1:C:274:LEU:HD23	1:C:345:PHE:CE2	2.49	0.48
1:C:446:LYS:HD3	1:C:447:GLY:N	2.28	0.48
1:C:1000:LYS:O	1:C:1001:LEU:C	2.55	0.48
1:B:386:LEU:HD12	1:C:500:GLU:OE1	2.14	0.48
1:C:278:GLY:HA3	1:C:345:PHE:CZ	2.49	0.48
1:C:661:VAL:O	1:C:661:VAL:HG12	2.13	0.48
1:C:768:ARG:HH11	1:C:768:ARG:HG3	1.79	0.48
1:B:80:THR:HG21	1:B:438:ARG:HH11	1.78	0.47
1:B:426:LEU:O	1:B:427:ARG:C	2.57	0.47
1:B:975:THR:O	1:B:979:ILE:HG12	2.13	0.47
1:C:249:ILE:CG2	1:C:541:PRO:HG2	2.43	0.47
1:C:484:LEU:O	1:C:488:LEU:HG	2.14	0.47
1:B:374:TYR:O	1:B:375:LEU:HB2	2.14	0.47
1:B:477:ASP:C	1:B:479:VAL:N	2.73	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:609:SER:O	1:B:613:ILE:HG12	2.14	0.47
1:B:816:ARG:HA	1:B:838:TRP:O	2.14	0.47
1:B:856:ASP:O	1:B:857:ASP:C	2.57	0.47
1:C:490:GLU:O	1:C:493:ASP:CB	2.60	0.47
1:C:754:PHE:HE2	1:C:906:PHE:HB2	1.78	0.47
1:B:272:GLU:HB2	1:B:299:PHE:CZ	2.50	0.47
1:B:685:VAL:O	1:B:688:ARG:HB3	2.14	0.47
1:B:740:VAL:CG1	1:B:741:ARG:N	2.78	0.47
1:C:610:ARG:HH11	1:C:610:ARG:CB	2.21	0.47
1:C:791:VAL:CG2	1:C:977:ALA:HA	2.43	0.47
1:C:887:PHE:C	1:C:887:PHE:CD1	2.93	0.47
1:B:154:ARG:CZ	1:B:174:GLY:HA2	2.45	0.47
1:B:333:LEU:N	1:B:333:LEU:CD2	2.76	0.47
1:B:758:ILE:HG22	1:B:938:THR:OG1	2.14	0.47
1:C:172:GLU:HG2	1:C:172:GLU:O	2.14	0.47
1:C:455:ASP:O	1:C:456:SER:C	2.57	0.47
1:C:476:ILE:O	1:C:476:ILE:HG23	2.14	0.47
1:C:581:ILE:HG22	1:C:582:LYS:N	2.30	0.47
1:C:834:VAL:CG1	1:C:835:GLY:H	2.21	0.47
1:C:858:VAL:HG21	1:C:880:TYR:HE1	1.79	0.47
1:B:95:LEU:HB2	1:B:102:CYS:SG	2.54	0.47
1:B:468:LEU:HD22	1:B:468:LEU:N	2.29	0.47
1:B:512:GLU:CD	1:B:630:THR:HB	2.39	0.47
1:B:705:ARG:HD3	1:B:1008:PHE:CE1	2.49	0.47
1:B:720:VAL:O	1:B:721:LEU:C	2.58	0.47
1:B:795:MET:HE1	1:B:811:ARG:HD2	1.95	0.47
1:C:106:PHE:O	1:C:163:THR:N	2.44	0.47
1:C:301:ILE:N	1:C:301:ILE:CD1	2.76	0.47
1:C:716:ALA:O	1:C:720:VAL:HG23	2.14	0.47
1:C:828:TYR:CB	1:C:888:MET:HE3	2.45	0.47
1:B:123:ALA:C	1:B:125:LYS:H	2.22	0.47
1:C:140:ARG:O	1:C:141:ILE:O	2.32	0.47
1:B:279:ILE:HD11	1:B:315:PHE:CE1	2.48	0.47
1:B:305:THR:O	1:B:306:ALA:C	2.57	0.47
1:B:394:ASP:HA	1:B:397:ARG:HG3	1.97	0.47
1:B:559:LEU:HD12	1:B:688:ARG:HG3	1.96	0.47
1:B:719:LEU:HD13	1:B:970:ILE:CD1	2.45	0.47
1:B:720:VAL:C	1:B:722:GLY:N	2.73	0.47
1:B:749:HIS:CG	1:B:906:PHE:HE2	2.32	0.47
1:B:972:ARG:NH1	1:B:1026:GLU:CG	2.78	0.47
1:C:71:LYS:HG2	1:C:219:CYS:HB3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:105:ILE:CG2	1:C:163:THR:HA	2.44	0.47
1:C:160:ILE:C	1:C:161:THR:CG2	2.87	0.47
1:C:170:TYR:CE2	1:C:171:ARG:HG2	2.50	0.47
1:C:201:LEU:O	1:C:217:ARG:HB2	2.14	0.47
1:C:260:ASN:O	1:C:260:ASN:ND2	2.48	0.47
1:C:283:ARG:NH2	1:C:453:GLU:OE2	2.47	0.47
1:C:544:LYS:O	1:C:545:TYR:CG	2.68	0.47
1:C:557:VAL:C	1:C:581:ILE:HD11	2.40	0.47
1:C:754:PHE:CE2	1:C:906:PHE:HB2	2.50	0.47
1:C:818:SER:O	1:C:819:ASP:C	2.58	0.47
1:C:912:GLU:OE1	1:C:912:GLU:N	2.47	0.47
1:B:369:VAL:CG1	1:B:393:ILE:HG23	2.38	0.47
1:B:476:ILE:C	1:B:478:GLU:H	2.23	0.47
1:C:212:TRP:CD1	1:C:212:TRP:H	2.33	0.47
1:C:259:VAL:HB	1:C:452:LEU:HD22	1.96	0.47
1:C:363:SER:O	1:C:364:LEU:HB2	2.15	0.47
1:C:413:ASP:OD1	1:C:413:ASP:C	2.57	0.47
1:B:60:ILE:HG21	1:B:82:VAL:HG11	1.96	0.47
1:B:360:ASP:O	1:B:361:ILE:C	2.57	0.47
1:B:510:ILE:HD11	1:B:616:LEU:HD22	1.97	0.47
1:B:807:ILE:CA	1:B:850:THR:HG23	2.45	0.47
1:C:95:LEU:HB2	1:C:102:CYS:SG	2.54	0.47
1:C:131:GLU:O	1:C:133:LEU:N	2.48	0.47
1:C:236:PHE:O	1:C:240:LEU:N	2.45	0.47
1:C:830:GLY:C	1:C:832:ASP:H	2.23	0.47
1:C:884:PHE:CE1	1:C:888:MET:HG3	2.49	0.47
1:B:175:HIS:C	1:B:176:PHE:CD1	2.93	0.47
1:B:192:LYS:O	1:B:193:ASN:C	2.58	0.47
1:B:271:LEU:CD2	1:B:481:PHE:HZ	2.24	0.47
1:B:370:LYS:O	1:B:373:ALA:N	2.48	0.47
1:B:440:PHE:N	1:B:440:PHE:CD1	2.83	0.47
1:B:444:LEU:O	1:B:445:THR:C	2.57	0.47
1:B:708:ASN:C	1:B:709:ARG:O	2.55	0.47
1:B:276:THR:HA	1:B:322:HIS:CE1	2.50	0.46
1:B:470:ASP:O	1:B:472:GLU:CB	2.58	0.46
1:B:1037:LEU:CD2	1:B:1038:ASP:N	2.76	0.46
1:C:65:ALA:O	1:C:66:LYS:C	2.58	0.46
1:C:73:SER:HA	1:C:220:LEU:O	2.14	0.46
1:C:249:ILE:HD12	1:C:541:PRO:HG2	1.97	0.46
1:C:469:TYR:O	1:C:470:ASP:C	2.58	0.46
1:C:779:MET:HE2	1:C:879:LEU:HD11	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:ILE:HD13	1:B:546:VAL:HG21	1.97	0.46
1:B:64:TRP:O	1:B:65:ALA:C	2.57	0.46
1:B:745:LEU:HD22	1:B:747:LEU:CD1	2.45	0.46
1:B:801:LEU:HD12	1:B:801:LEU:HA	1.63	0.46
1:C:179:ILE:O	1:C:220:LEU:HD12	2.14	0.46
1:C:439:LEU:CD1	1:C:444:LEU:HD13	2.44	0.46
1:C:544:LYS:HB3	1:C:545:TYR:CD1	2.50	0.46
1:C:873:ARG:H	1:C:876:HIS:CD2	2.28	0.46
1:B:103:TYR:CE1	1:B:161:THR:HG23	2.43	0.46
1:B:292:TYR:HD2	1:B:303:ILE:HB	1.80	0.46
1:B:361:ILE:O	1:B:362:ASP:C	2.58	0.46
1:B:492:ARG:NH1	1:B:492:ARG:CG	2.77	0.46
1:B:565:PHE:N	1:B:568:VAL:O	2.48	0.46
1:B:736:LYS:O	1:B:948:LYS:O	2.34	0.46
1:B:830:GLY:C	1:B:832:ASP:H	2.23	0.46
1:B:957:THR:O	1:B:959:SER:N	2.49	0.46
1:B:957:THR:HG22	1:B:995:ARG:HG2	1.97	0.46
1:C:154:ARG:NE	1:C:174:GLY:CA	2.78	0.46
1:C:608:ASN:C	1:C:610:ARG:N	2.72	0.46
1:C:1010:VAL:CG1	1:C:1011:ARG:N	2.77	0.46
1:B:58:ARG:N	1:B:61:GLN:HE21	2.06	0.46
1:B:301:ILE:N	1:B:301:ILE:CD1	2.76	0.46
1:B:385:LEU:O	1:B:387:PRO:CD	2.63	0.46
1:B:477:ASP:HA	1:B:479:VAL:HG13	1.97	0.46
1:B:720:VAL:O	1:B:723:TRP:N	2.48	0.46
1:C:314:LYS:HA	1:C:319:GLU:OE2	2.16	0.46
1:C:561:THR:CB	1:C:585:ARG:HH11	2.29	0.46
1:C:723:TRP:CZ3	1:C:964:MET:HG2	2.50	0.46
1:B:194:VAL:O	1:B:198:LEU:HD12	2.15	0.46
1:B:387:PRO:HB2	1:B:389:VAL:CA	2.45	0.46
1:B:628:VAL:HG21	1:B:640:ALA:HA	1.97	0.46
1:B:873:ARG:N	1:B:876:HIS:HD2	2.09	0.46
1:B:1038:ASP:C	1:B:1038:ASP:OD2	2.59	0.46
1:C:212:TRP:C	1:C:213:VAL:HG22	2.41	0.46
1:C:682:LYS:HD3	1:C:1039:TYR:HD2	1.80	0.46
1:C:743:PHE:C	1:C:745:LEU:N	2.74	0.46
1:C:793:GLN:O	1:C:797:ILE:HG12	2.15	0.46
1:B:309:LYS:CG	1:B:310:GLY:H	2.27	0.46
1:B:544:LYS:HB3	1:B:545:TYR:CD1	2.51	0.46
1:B:983:LEU:HB3	1:B:989:VAL:CG2	2.45	0.46
1:C:212:TRP:CH2	1:C:235:LEU:HD21	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:756:LEU:HD22	1:C:906:PHE:HB3	1.97	0.46
1:C:975:THR:O	1:C:976:TYR:C	2.57	0.46
1:B:301:ILE:O	1:B:320:ILE:HG23	2.16	0.46
1:B:330:TYR:CD1	1:B:337:LEU:HD12	2.51	0.46
1:B:366:PRO:C	1:B:368:MET:H	2.23	0.46
1:C:52:LYS:HE3	1:C:125:LYS:O	2.16	0.46
1:C:67:ARG:HD3	1:C:74:PHE:CD2	2.51	0.46
1:C:68:ILE:HB	1:C:91:MET:HE3	1.98	0.46
1:C:259:VAL:HB	1:C:452:LEU:HD23	1.96	0.46
1:C:374:TYR:C	1:C:376:TYR:N	2.73	0.46
1:C:377:ARG:O	1:C:378:ASN:C	2.57	0.46
1:C:413:ASP:OD2	1:C:465:ARG:NH2	2.49	0.46
1:C:1005:VAL:O	1:C:1006:PHE:C	2.58	0.46
1:C:1009:LEU:HD22	1:C:1009:LEU:N	2.31	0.46
1:B:11:PRO:O	1:B:12:VAL:CB	2.64	0.46
1:B:381:GLU:N	1:B:381:GLU:CD	2.74	0.46
1:B:720:VAL:O	1:B:722:GLY:N	2.49	0.46
1:B:775:THR:HG22	1:B:814:SER:O	2.15	0.46
1:B:834:VAL:CG2	1:B:893:ARG:HA	2.46	0.46
1:C:720:VAL:O	1:C:721:LEU:C	2.59	0.46
1:C:981:ASP:O	1:C:983:LEU:N	2.49	0.46
1:B:254:VAL:HA	1:B:447:GLY:C	2.40	0.46
1:B:556:VAL:HG23	1:B:557:VAL:N	2.29	0.46
1:B:733:GLU:HB2	1:B:952:LYS:HB2	1.97	0.46
1:C:268:SER:O	1:C:272:GLU:HG2	2.16	0.46
1:C:510:ILE:HD13	1:C:616:LEU:HD22	1.98	0.46
1:C:558:ASP:OD2	1:C:559:LEU:N	2.47	0.46
1:C:635:GLU:HG2	1:C:684:GLN:HE21	1.81	0.46
1:C:763:GLU:CD	1:C:900:LYS:HE2	2.41	0.46
1:C:1007:ARG:C	1:C:1010:VAL:HG12	2.41	0.46
1:C:1029:MET:HA	1:C:1032:ILE:CG2	2.46	0.46
1:B:166:LEU:HD11	1:B:200:LEU:HD11	1.97	0.46
1:C:413:ASP:CG	1:C:860:ARG:NH2	2.74	0.46
1:C:747:LEU:N	1:C:747:LEU:CD1	2.79	0.46
1:C:760:LEU:HD21	1:C:938:THR:CB	2.46	0.46
1:C:975:THR:HG22	1:C:979:ILE:CG1	2.42	0.46
1:B:187:ILE:HG23	1:B:193:ASN:HB3	1.98	0.45
1:B:267:LEU:O	1:B:270:ILE:N	2.48	0.45
1:B:386:LEU:N	1:B:386:LEU:CD2	2.73	0.45
1:B:832:ASP:HA	1:B:893:ARG:CD	2.46	0.45
1:C:380:ASP:HB3	1:C:383:GLU:OE2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:363:SER:O	1:B:364:LEU:CB	2.62	0.45
1:B:714:GLY:O	1:B:716:ALA:N	2.49	0.45
1:B:816:ARG:O	1:B:845:GLU:HG3	2.17	0.45
1:C:758:ILE:HB	1:C:939:GLY:H	1.82	0.45
1:C:832:ASP:HA	1:C:893:ARG:CD	2.47	0.45
1:B:264:ILE:O	1:B:264:ILE:HG12	2.16	0.45
1:B:331:GLY:O	1:B:336:GLY:HA2	2.17	0.45
1:B:466:ALA:HB1	1:B:472:GLU:CG	2.46	0.45
1:B:754:PHE:CZ	1:B:907:ASP:HB2	2.51	0.45
1:B:879:LEU:O	1:B:880:TYR:C	2.57	0.45
1:C:253:ASN:OD1	1:C:446:LYS:HG2	2.17	0.45
1:C:257:VAL:HB	1:C:450:PHE:HD1	1.77	0.45
1:C:320:ILE:CG2	1:C:321:ASP:N	2.79	0.45
1:C:682:LYS:HD3	1:C:1039:TYR:CD2	2.52	0.45
1:C:725:ILE:HD13	1:C:1003:ILE:HG12	1.98	0.45
1:C:740:VAL:HG12	1:C:741:ARG:H	1.81	0.45
1:C:858:VAL:O	1:C:862:ILE:HG13	2.16	0.45
1:B:74:PHE:CZ	1:B:221:MET:HE3	2.51	0.45
1:B:697:VAL:O	1:B:701:LYS:HG2	2.17	0.45
1:B:740:VAL:HG12	1:B:741:ARG:H	1.80	0.45
1:B:766:GLU:HA	1:B:766:GLU:OE2	2.16	0.45
1:C:68:ILE:C	1:C:70:ARG:H	2.24	0.45
1:C:508:LEU:HD13	1:C:619:LEU:HB3	1.99	0.45
1:C:1009:LEU:H	1:C:1009:LEU:CD2	2.29	0.45
1:B:52:LYS:HG2	1:B:125:LYS:HB3	1.99	0.45
1:B:134:ILE:HD11	1:B:160:ILE:HD12	1.97	0.45
1:B:315:PHE:CG	1:B:323:LEU:HD13	2.51	0.45
1:B:725:ILE:O	1:B:727:ARG:N	2.43	0.45
1:B:986:ARG:HD3	1:B:986:ARG:HA	1.66	0.45
1:C:194:VAL:O	1:C:197:LEU:N	2.41	0.45
1:C:279:ILE:HD12	1:C:323:LEU:HB2	1.97	0.45
1:C:580:SER:HB3	1:C:584:CYS:SG	2.57	0.45
1:C:627:ILE:HA	1:C:654:LYS:O	2.15	0.45
1:B:257:VAL:HG21	1:B:450:PHE:CE1	2.51	0.45
1:B:279:ILE:HG12	1:B:343:ILE:CD1	2.47	0.45
1:B:393:ILE:O	1:B:394:ASP:C	2.60	0.45
1:B:440:PHE:HD1	1:B:440:PHE:N	2.15	0.45
1:B:853:LEU:HD13	1:B:861:LEU:HD12	1.98	0.45
1:C:39:GLU:CD	1:C:39:GLU:N	2.58	0.45
1:C:43:LEU:HD23	1:C:47:VAL:HG23	1.99	0.45
1:C:341:GLU:OE2	1:C:496:ARG:HG2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:292:TYR:O	1:B:295:LEU:O	2.35	0.45
1:B:426:LEU:HD13	1:B:469:TYR:CD1	2.52	0.45
1:C:277:GLY:N	1:C:321:ASP:O	2.46	0.45
1:C:333:LEU:N	1:C:333:LEU:HD22	2.32	0.45
1:C:416:VAL:HG13	1:C:417:ARG:N	2.31	0.45
1:C:719:LEU:HD11	1:C:1019:GLU:HA	1.97	0.45
1:B:100:LYS:O	1:B:177:ASP:OD1	2.35	0.45
1:B:712:SER:HB3	1:B:986:ARG:HG3	1.98	0.45
1:C:304:VAL:HB	1:C:305:THR:H	1.57	0.45
1:B:37:PHE:HB3	1:B:39:GLU:OE2	2.16	0.45
1:B:277:GLY:N	1:B:321:ASP:O	2.40	0.45
1:B:511:VAL:O	1:B:551:ALA:HA	2.16	0.45
1:B:743:PHE:C	1:B:745:LEU:N	2.75	0.45
1:B:791:VAL:CG2	1:B:977:ALA:HA	2.47	0.45
1:C:363:SER:O	1:C:364:LEU:CB	2.64	0.45
1:C:807:ILE:N	1:C:850:THR:HG23	2.31	0.45
1:B:248:ARG:CG	1:B:249:ILE:H	2.29	0.45
1:B:263:SER:OG	1:B:267:LEU:HB2	2.17	0.45
1:B:426:LEU:C	1:B:430:ILE:HD12	2.42	0.45
1:B:620:ALA:CB	1:B:647:LEU:HD13	2.46	0.45
1:B:643:LEU:O	1:B:644:LYS:C	2.57	0.45
1:C:384:ARG:CZ	1:C:396:VAL:HB	2.47	0.45
1:C:584:CYS:O	1:C:585:ARG:HB2	2.16	0.45
1:C:790:SER:HA	1:C:958:GLN:NE2	2.28	0.45
1:C:986:ARG:HD3	1:C:986:ARG:HA	1.63	0.45
1:B:2:ILE:CA	1:B:11:PRO:N	2.80	0.44
1:B:89:LEU:HD12	1:B:118:THR:HG21	2.00	0.44
1:B:170:TYR:CE2	1:B:171:ARG:HG2	2.53	0.44
1:B:194:VAL:CG1	1:B:198:LEU:HD11	2.48	0.44
1:B:437:SER:HA	1:B:445:THR:OG1	2.17	0.44
1:B:438:ARG:HG3	1:B:439:LEU:H	1.81	0.44
1:B:825:ALA:HA	1:B:888:MET:HE1	1.99	0.44
1:B:860:ARG:O	1:B:863:GLN:N	2.50	0.44
1:C:887:PHE:O	1:C:889:ALA:N	2.50	0.44
1:B:303:ILE:HD13	1:B:304:VAL:N	2.32	0.44
1:B:693:TRP:O	1:B:697:VAL:HG23	2.18	0.44
1:B:699:SER:O	1:B:702:LEU:HB2	2.17	0.44
1:B:703:TRP:CH2	1:B:709:ARG:NH1	2.85	0.44
1:B:760:LEU:CD2	1:B:938:THR:OG1	2.65	0.44
1:C:640:ALA:HA	1:C:643:LEU:HD12	1.98	0.44
1:C:966:LYS:O	1:C:969:GLY:N	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:PHE:O	1:B:49:PHE:HB3	2.17	0.44
1:B:100:LYS:HG2	1:B:177:ASP:CB	2.47	0.44
1:B:377:ARG:HG3	1:B:382:ILE:CG1	2.46	0.44
1:B:381:GLU:N	1:B:381:GLU:OE1	2.51	0.44
1:B:689:ILE:H	1:B:689:ILE:HG12	1.64	0.44
1:B:1050:ILE:HG22	1:B:1051:LYS:N	2.33	0.44
1:C:562:ASN:O	1:C:563:ARG:C	2.61	0.44
1:C:611:SER:O	1:C:615:ALA:N	2.50	0.44
1:C:739:ILE:HG22	1:C:740:VAL:O	2.17	0.44
1:C:1017:VAL:O	1:C:1017:VAL:CG1	2.66	0.44
1:C:54:VAL:HG12	1:C:55:GLY:N	2.32	0.44
1:C:252:ARG:NH2	1:C:256:ASP:OD1	2.50	0.44
1:C:553:ILE:HB	1:C:612:ARG:NE	2.32	0.44
1:C:965:MET:HE1	1:C:980:VAL:HG23	2.00	0.44
1:B:700:GLN:O	1:B:703:TRP:N	2.40	0.44
1:B:791:VAL:H	1:B:958:GLN:HE22	1.65	0.44
1:C:43:LEU:O	1:C:46:PHE:N	2.50	0.44
1:C:529:VAL:HG12	1:C:530:LYS:N	2.33	0.44
1:C:854:THR:O	1:C:858:VAL:HG23	2.17	0.44
1:B:144:ARG:HG2	1:B:144:ARG:HH11	1.82	0.44
1:B:206:ASP:O	1:B:207:LEU:C	2.61	0.44
1:B:737:ILE:HD11	1:B:746:VAL:CG1	2.47	0.44
1:B:767:LEU:HA	1:B:895:PHE:O	2.18	0.44
1:B:873:ARG:H	1:B:876:HIS:CD2	2.25	0.44
1:B:1005:VAL:O	1:B:1009:LEU:HD23	2.17	0.44
1:B:1012:ARG:C	1:B:1013:TYR:CD1	2.96	0.44
1:C:138:HIS:O	1:C:140:ARG:N	2.50	0.44
1:C:254:VAL:HA	1:C:447:GLY:CA	2.48	0.44
1:C:617:ARG:HG3	1:C:647:LEU:HD23	1.99	0.44
1:C:774:TYR:CG	1:C:886:ARG:HG2	2.53	0.44
1:C:1020:ASP:CG	1:C:1023:ARG:HH22	2.26	0.44
1:B:386:LEU:H	1:B:386:LEU:CD2	1.97	0.44
1:B:901:LYS:CG	1:B:902:TYR:N	2.79	0.44
1:C:95:LEU:HB3	1:C:100:LYS:HB3	1.99	0.44
1:C:197:LEU:O	1:C:198:LEU:C	2.59	0.44
1:C:252:ARG:CA	1:C:444:LEU:HD21	2.47	0.44
1:C:439:LEU:HD13	1:C:444:LEU:HD13	2.00	0.44
1:C:565:PHE:N	1:C:568:VAL:O	2.51	0.44
1:C:883:ILE:H	1:C:883:ILE:CD1	2.31	0.44
1:B:12:VAL:HA	1:B:19:SER:O	2.17	0.44
1:B:470:ASP:HB3	1:B:471:ILE:H	1.61	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:669:ALA:C	1:B:671:GLU:N	2.76	0.44
1:B:860:ARG:O	1:B:861:LEU:C	2.61	0.44
1:B:1029:MET:HA	1:B:1032:ILE:HG23	1.99	0.44
1:B:320:ILE:HG21	1:B:322:HIS:O	2.17	0.44
1:C:11:PRO:HA	1:C:16:ASP:CB	2.47	0.44
1:C:166:LEU:HD11	1:C:200:LEU:HD11	2.00	0.44
1:C:436:THR:HG23	1:C:436:THR:O	2.18	0.44
1:C:685:VAL:O	1:C:686:VAL:C	2.60	0.44
1:C:702:LEU:HD21	1:C:1009:LEU:HD21	1.99	0.44
1:B:31:LYS:N	1:B:32:ARG:CZ	2.81	0.43
1:B:546:VAL:O	1:B:546:VAL:HG12	2.17	0.43
1:B:736:LYS:NZ	1:B:736:LYS:CA	2.79	0.43
1:C:58:ARG:O	1:C:61:GLN:N	2.51	0.43
1:C:121:LYS:O	1:C:125:LYS:NZ	2.48	0.43
1:C:154:ARG:NE	1:C:174:GLY:HA3	2.33	0.43
1:C:768:ARG:HG3	1:C:768:ARG:NH1	2.31	0.43
1:C:1028:ARG:O	1:C:1031:ALA:HB3	2.17	0.43
1:B:276:THR:HG21	1:B:321:ASP:CG	2.43	0.43
1:B:366:PRO:CA	1:B:393:ILE:HG12	2.48	0.43
1:B:382:ILE:O	1:B:386:LEU:CD2	2.66	0.43
1:B:468:LEU:HD22	1:B:468:LEU:H	1.83	0.43
1:B:509:PHE:O	1:B:549:VAL:HA	2.17	0.43
1:B:832:ASP:CA	1:B:893:ARG:HG3	2.47	0.43
1:C:43:LEU:O	1:C:44:LYS:C	2.59	0.43
1:C:436:THR:HG23	1:C:447:GLY:CA	2.44	0.43
1:C:477:ASP:HA	1:C:479:VAL:HG13	1.99	0.43
1:C:609:SER:CB	1:C:612:ARG:NH1	2.81	0.43
1:C:740:VAL:CG1	1:C:741:ARG:N	2.80	0.43
1:C:754:PHE:C	1:C:943:VAL:HG23	2.43	0.43
1:B:233:ALA:C	1:B:235:LEU:N	2.73	0.43
1:B:314:LYS:HA	1:B:319:GLU:OE2	2.19	0.43
1:B:393:ILE:N	1:B:393:ILE:CD1	2.82	0.43
1:B:763:GLU:OE1	1:B:900:LYS:HE2	2.18	0.43
1:B:907:ASP:O	1:B:909:LYS:N	2.51	0.43
1:C:212:TRP:HB2	1:C:213:VAL:H	1.59	0.43
1:C:337:LEU:CD2	1:C:343:ILE:HD11	2.46	0.43
1:C:570:VAL:O	1:C:573:ARG:HB2	2.17	0.43
1:C:734:ARG:C	1:C:736:LYS:H	2.25	0.43
1:B:207:LEU:HG	1:B:208:LYS:N	2.33	0.43
1:B:518:ARG:HG2	1:B:522:ARG:CZ	2.48	0.43
1:B:638:LYS:HB2	1:B:684:GLN:CG	2.28	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:702:LEU:HD21	1:B:1009:LEU:HD21	1.99	0.43
1:B:756:LEU:O	1:B:940:THR:HA	2.18	0.43
1:B:832:ASP:HA	1:B:893:ARG:CG	2.48	0.43
1:B:947:VAL:O	1:B:948:LYS:HG3	2.19	0.43
1:C:125:LYS:O	1:C:126:ALA:HB3	2.19	0.43
1:C:276:THR:HG21	1:C:321:ASP:OD1	2.18	0.43
1:C:291:ILE:O	1:C:292:TYR:C	2.61	0.43
1:B:797:ILE:O	1:B:798:ALA:C	2.61	0.43
1:C:118:THR:HG22	1:C:122:TYR:CZ	2.54	0.43
1:C:274:LEU:HG	1:C:274:LEU:O	2.19	0.43
1:C:357:THR:CG2	1:C:358:ILE:N	2.81	0.43
1:C:372:LEU:O	1:C:373:ALA:C	2.62	0.43
1:C:706:PHE:O	1:C:708:ASN:N	2.52	0.43
1:C:824:ILE:HG23	1:C:852:PRO:CD	2.48	0.43
1:C:1012:ARG:O	1:C:1013:TYR:CD1	2.72	0.43
1:B:440:PHE:C	1:B:442:GLY:N	2.74	0.43
1:C:13:CYS:O	1:C:20:LYS:HA	2.18	0.43
1:C:320:ILE:CG2	1:C:322:HIS:O	2.66	0.43
1:C:735:ARG:NH1	1:C:951:PRO:HG3	2.34	0.43
1:C:736:LYS:HB3	1:C:737:ILE:H	1.48	0.43
1:B:315:PHE:HA	1:B:320:ILE:HB	2.00	0.43
1:B:522:ARG:HA	1:B:525:GLY:O	2.19	0.43
1:B:628:VAL:HG13	1:B:630:THR:HG23	1.99	0.43
1:B:797:ILE:HG23	1:B:876:HIS:ND1	2.33	0.43
1:B:883:ILE:H	1:B:883:ILE:CD1	2.29	0.43
1:C:94:PHE:O	1:C:98:LYS:HD3	2.19	0.43
1:C:109:SER:O	1:C:112:VAL:HB	2.17	0.43
1:C:236:PHE:HD1	1:C:244:ILE:HD11	1.83	0.43
1:C:259:VAL:N	1:C:451:LEU:O	2.46	0.43
1:C:427:ARG:HH11	1:C:427:ARG:HG3	1.83	0.43
1:C:517:ALA:O	1:C:518:ARG:C	2.61	0.43
1:C:613:ILE:CD1	1:C:613:ILE:H	2.31	0.43
1:C:719:LEU:HD13	1:C:970:ILE:HD13	2.00	0.43
1:C:991:GLU:O	1:C:992:LYS:HE2	2.18	0.43
1:B:38:PRO:C	1:B:40:ASP:N	2.77	0.43
1:B:101:ARG:NH1	1:B:154:ARG:HA	2.33	0.43
1:B:361:ILE:O	1:B:362:ASP:O	2.36	0.43
1:B:369:VAL:HG21	1:B:393:ILE:HG23	1.95	0.43
1:B:436:THR:HG23	1:B:447:GLY:CA	2.46	0.43
1:B:560:ILE:HG13	1:B:561:THR:N	2.32	0.43
1:B:873:ARG:O	1:B:874:TRP:C	2.59	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:999:THR:O	1:B:1000:LYS:C	2.59	0.43
1:C:35:CYS:HB2	1:C:648:SER:O	2.18	0.43
1:C:89:LEU:O	1:C:92:SER:HB3	2.18	0.43
1:C:677:ASP:O	1:C:678:GLU:C	2.62	0.43
1:C:706:PHE:O	1:C:706:PHE:CD1	2.70	0.43
1:B:43:LEU:O	1:B:44:LYS:C	2.62	0.43
1:B:60:ILE:CG2	1:B:82:VAL:HG11	2.49	0.43
1:B:103:TYR:HD1	1:B:159:VAL:O	2.02	0.43
1:B:304:VAL:HB	1:B:305:THR:H	1.63	0.43
1:B:377:ARG:O	1:B:379:VAL:O	2.37	0.43
1:B:694:ILE:H	1:B:694:ILE:HG12	1.60	0.43
1:B:743:PHE:N	1:B:743:PHE:CD1	2.86	0.43
1:B:964:MET:O	1:B:965:MET:C	2.62	0.43
1:C:63:MET:CE	1:C:67:ARG:NH2	2.79	0.43
1:C:236:PHE:CD1	1:C:244:ILE:HD11	2.53	0.43
1:C:263:SER:C	1:C:265:SER:N	2.76	0.43
1:C:358:ILE:HD11	1:C:419:GLY:O	2.17	0.43
1:C:470:ASP:HB3	1:C:471:ILE:H	1.54	0.43
1:C:666:ILE:O	1:C:669:ALA:N	2.51	0.43
1:C:758:ILE:CG2	1:C:938:THR:HA	2.49	0.43
1:C:793:GLN:O	1:C:794:THR:C	2.62	0.43
1:C:1003:ILE:O	1:C:1004:ASP:C	2.62	0.43
1:B:267:LEU:O	1:B:268:SER:C	2.62	0.43
1:B:274:LEU:HD22	1:B:450:PHE:HE2	1.83	0.43
1:B:357:THR:HG22	1:B:358:ILE:N	2.34	0.43
1:B:374:TYR:O	1:B:376:TYR:N	2.52	0.43
1:B:440:PHE:CE2	1:B:442:GLY:HA3	2.54	0.43
1:B:505:LYS:O	1:B:545:TYR:HB3	2.19	0.43
1:B:1040:LEU:C	1:B:1042:ALA:H	2.27	0.43
1:B:1050:ILE:C	1:B:1052:SER:N	2.77	0.43
1:C:22:ILE:CB	1:C:641:TRP:HH2	2.32	0.43
1:C:154:ARG:NH2	1:C:174:GLY:HA2	2.34	0.43
1:C:260:ASN:O	1:C:454:ASP:HA	2.18	0.43
1:C:296:LYS:O	1:C:301:ILE:HD11	2.18	0.43
1:C:305:THR:HG22	1:C:306:ALA:N	2.23	0.43
1:C:665:ALA:HA	1:C:668:GLU:OE2	2.19	0.43
1:C:766:GLU:HB3	1:C:768:ARG:HE	1.83	0.43
1:C:972:ARG:NH1	1:C:1026:GLU:HG3	2.34	0.43
1:B:530:LYS:HB3	1:B:537:VAL:CG1	2.49	0.42
1:B:936:LEU:O	1:B:936:LEU:CD1	2.67	0.42
1:C:260:ASN:HA	1:C:453:GLU:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:753:GLU:HG3	1:B:754:PHE:N	2.34	0.42
1:C:122:TYR:O	1:C:123:ALA:C	2.62	0.42
1:C:220:LEU:HD12	1:C:221:MET:N	2.33	0.42
1:C:537:VAL:HG13	1:C:537:VAL:O	2.18	0.42
1:C:942:ARG:H	1:C:942:ARG:HG2	1.64	0.42
1:B:283:ARG:HG3	1:B:287:GLU:OE1	2.18	0.42
1:B:296:LYS:HA	1:B:301:ILE:HG12	2.02	0.42
1:B:544:LYS:O	1:B:545:TYR:CG	2.72	0.42
1:B:774:TYR:OH	1:B:782:ASP:CB	2.67	0.42
1:B:935:GLU:HG2	1:B:936:LEU:H	1.84	0.42
1:B:1029:MET:O	1:B:1030:ASP:C	2.61	0.42
1:C:697:VAL:O	1:C:700:GLN:HB3	2.20	0.42
1:C:743:PHE:O	1:C:745:LEU:N	2.51	0.42
1:C:829:LEU:O	1:C:830:GLY:O	2.37	0.42
1:C:860:ARG:HH11	1:C:860:ARG:CB	2.30	0.42
1:C:860:ARG:O	1:C:861:LEU:C	2.61	0.42
1:B:39:GLU:H	1:B:39:GLU:CD	2.26	0.42
1:B:176:PHE:HB2	1:B:179:ILE:CD1	2.49	0.42
1:B:834:VAL:HG21	1:B:893:ARG:HA	2.01	0.42
1:C:63:MET:HG3	1:C:622:ASP:O	2.19	0.42
1:C:170:TYR:CG	1:C:171:ARG:N	2.87	0.42
1:C:315:PHE:HB2	1:C:323:LEU:HD11	2.00	0.42
1:C:476:ILE:O	1:C:476:ILE:HG13	2.18	0.42
1:C:583:ARG:O	1:C:583:ARG:HG2	2.19	0.42
1:C:608:ASN:C	1:C:610:ARG:H	2.25	0.42
1:C:756:LEU:CD1	1:C:904:ILE:HB	2.49	0.42
1:B:60:ILE:CD1	1:B:546:VAL:HG11	2.50	0.42
1:B:396:VAL:O	1:B:400:LEU:HG	2.20	0.42
1:B:829:LEU:HB2	1:B:833:PHE:HB2	2.00	0.42
1:B:853:LEU:O	1:B:854:THR:O	2.38	0.42
1:B:879:LEU:C	1:B:881:ASP:N	2.77	0.42
1:C:101:ARG:HD3	1:C:176:PHE:CD1	2.54	0.42
1:C:152:ASN:O	1:C:153:LEU:C	2.62	0.42
1:C:154:ARG:HE	1:C:174:GLY:CA	2.29	0.42
1:C:507:ALA:O	1:C:508:LEU:C	2.61	0.42
1:C:658:PHE:CD1	1:C:658:PHE:N	2.87	0.42
1:B:690:GLU:HG3	1:B:694:ILE:CG1	2.49	0.42
1:B:764:ARG:HD3	1:B:764:ARG:H	1.83	0.42
1:B:838:TRP:O	1:B:839:GLY:O	2.38	0.42
1:B:912:GLU:O	1:B:912:GLU:HG2	2.19	0.42
1:B:912:GLU:OE1	1:B:912:GLU:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:989:VAL:HG12	1:B:990:VAL:N	2.34	0.42
1:C:43:LEU:O	1:C:46:PHE:HB3	2.19	0.42
1:C:58:ARG:CB	1:C:61:GLN:HE21	2.27	0.42
1:C:148:ASN:N	1:C:148:ASN:ND2	2.67	0.42
1:C:638:LYS:CB	1:C:684:GLN:HG3	2.39	0.42
1:C:812:THR:CG2	1:C:847:ILE:HG13	2.49	0.42
1:C:857:ASP:O	1:C:861:LEU:HG	2.20	0.42
1:C:860:ARG:CG	1:C:860:ARG:NH1	2.82	0.42
1:C:1012:ARG:NH1	1:C:1012:ARG:HB3	2.34	0.42
1:B:58:ARG:CB	1:B:61:GLN:HE21	2.26	0.42
1:B:508:LEU:CD2	1:B:510:ILE:HD11	2.39	0.42
1:B:991:GLU:HG3	1:B:995:ARG:O	2.19	0.42
1:C:43:LEU:HD23	1:C:47:VAL:CG2	2.50	0.42
1:C:370:LYS:O	1:C:373:ALA:CB	2.63	0.42
1:C:530:LYS:HB3	1:C:537:VAL:CG1	2.50	0.42
1:C:557:VAL:O	1:C:581:ILE:HD11	2.20	0.42
1:C:570:VAL:HB	1:C:571:ASN:H	1.62	0.42
1:C:687:ARG:HA	1:C:1029:MET:HE1	2.01	0.42
1:C:763:GLU:OE1	1:C:900:LYS:HE2	2.20	0.42
1:C:999:THR:O	1:C:1000:LYS:C	2.62	0.42
1:B:58:ARG:NH2	1:B:82:VAL:HA	2.35	0.42
1:B:153:LEU:O	1:B:154:ARG:HG3	2.19	0.42
1:B:260:ASN:O	1:B:260:ASN:CG	2.63	0.42
1:B:426:LEU:HD11	1:B:466:ALA:CA	2.45	0.42
1:B:483:LYS:O	1:B:484:LEU:C	2.61	0.42
1:B:756:LEU:HD11	1:B:904:ILE:HB	2.02	0.42
1:B:756:LEU:HD13	1:B:757:THR:H	1.85	0.42
1:B:772:PRO:O	1:B:886:ARG:NH1	2.46	0.42
1:B:872:LEU:HD22	1:B:876:HIS:CG	2.55	0.42
1:B:1051:LYS:C	1:B:1053:ILE:N	2.77	0.42
1:C:187:ILE:HG23	1:C:193:ASN:HB3	2.02	0.42
1:C:283:ARG:HG3	1:C:287:GLU:OE1	2.20	0.42
1:C:305:THR:OG1	1:C:329:TYR:CB	2.52	0.42
1:C:318:GLY:HA2	1:C:342:ARG:HH11	1.85	0.42
1:C:351:CYS:HB3	1:C:462:PHE:CE1	2.55	0.42
1:B:341:GLU:OE2	1:B:496:ARG:HG2	2.19	0.42
1:B:494:ARG:O	1:B:497:ARG:N	2.53	0.42
1:B:533:ASP:OD1	1:B:534:GLY:N	2.44	0.42
1:B:641:TRP:O	1:B:642:ASP:C	2.62	0.42
1:B:709:ARG:HH11	1:B:709:ARG:CG	2.33	0.42
1:B:756:LEU:HD22	1:B:906:PHE:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:86:SER:N	2:C:2055:ANP:O1A	2.52	0.42
1:C:138:HIS:H	1:C:141:ILE:CD1	2.32	0.42
1:C:357:THR:HG22	1:C:358:ILE:H	1.85	0.42
1:C:374:TYR:O	1:C:376:TYR:N	2.53	0.42
1:C:480:ASP:OD1	1:C:483:LYS:HG3	2.19	0.42
1:C:879:LEU:O	1:C:881:ASP:N	2.52	0.42
1:B:58:ARG:HH21	1:B:82:VAL:HA	1.85	0.42
1:B:103:TYR:O	1:B:179:ILE:HA	2.19	0.42
1:B:258:ALA:O	1:B:476:ILE:HD12	2.19	0.42
1:B:906:PHE:N	1:B:906:PHE:CD1	2.88	0.42
1:B:1012:ARG:CG	1:B:1012:ARG:NH1	2.82	0.42
1:C:37:PHE:CD2	1:C:621:HIS:ND1	2.88	0.42
1:C:49:PHE:O	1:C:50:PHE:C	2.61	0.42
1:C:58:ARG:HE	1:C:82:VAL:HA	1.85	0.42
1:C:122:TYR:C	1:C:124:GLU:N	2.78	0.42
1:C:252:ARG:N	1:C:444:LEU:HD21	2.35	0.42
1:C:515:THR:O	1:C:516:LYS:C	2.61	0.42
1:C:580:SER:OG	1:C:610:ARG:HB2	2.19	0.42
1:C:617:ARG:HG3	1:C:647:LEU:CD2	2.50	0.42
1:C:642:ASP:O	1:C:646:LEU:HD12	2.20	0.42
1:C:686:VAL:O	1:C:687:ARG:C	2.63	0.42
1:C:784:ASN:ND2	1:C:958:GLN:HB2	2.35	0.42
1:C:874:TRP:CD2	1:C:875:GLU:N	2.88	0.42
1:B:63:MET:HE3	1:B:67:ARG:NH2	2.35	0.41
1:B:160:ILE:C	1:B:161:THR:CG2	2.93	0.41
1:B:314:LYS:HG2	1:B:319:GLU:OE1	2.20	0.41
1:B:333:LEU:CD2	1:B:333:LEU:H	2.32	0.41
1:B:483:LYS:O	1:B:486:ARG:N	2.53	0.41
1:C:46:PHE:HA	1:C:94:PHE:CE1	2.54	0.41
1:C:198:LEU:HD12	1:C:198:LEU:N	2.21	0.41
1:C:272:GLU:OE1	1:C:322:HIS:CD2	2.73	0.41
1:C:426:LEU:O	1:C:427:ARG:C	2.63	0.41
1:C:480:ASP:O	1:C:483:LYS:N	2.51	0.41
1:C:540:ILE:HG13	1:C:541:PRO:N	2.35	0.41
1:C:608:ASN:ND2	1:C:610:ARG:CB	2.83	0.41
1:C:613:ILE:CD1	1:C:613:ILE:N	2.82	0.41
1:C:879:LEU:C	1:C:881:ASP:N	2.77	0.41
1:B:400:LEU:O	1:B:403:VAL:O	2.38	0.41
1:B:417:ARG:HB3	1:B:418:GLU:H	1.62	0.41
1:B:469:TYR:O	1:B:470:ASP:O	2.38	0.41
1:B:544:LYS:HE3	1:B:671:GLU:OE1	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:734:ARG:HB2	1:B:952:LYS:HA	2.02	0.41
1:B:934:ASN:HD22	1:B:935:GLU:H	1.68	0.41
1:C:380:ASP:C	1:C:381:GLU:OE1	2.63	0.41
1:C:440:PHE:CE2	1:C:495:TYR:HA	2.55	0.41
1:C:634:THR:HB	1:C:684:GLN:HB2	2.02	0.41
1:C:733:GLU:O	1:C:952:LYS:N	2.53	0.41
1:C:784:ASN:ND2	1:C:959:SER:N	2.59	0.41
1:B:393:ILE:HG22	1:B:397:ARG:CG	2.50	0.41
1:B:608:ASN:ND2	1:B:610:ARG:HB3	2.35	0.41
1:B:750:ASP:OD1	1:B:751:GLU:OE1	2.38	0.41
1:C:119:ILE:O	1:C:120:ARG:C	2.63	0.41
1:C:138:HIS:CE1	1:C:140:ARG:HB3	2.56	0.41
1:C:253:ASN:O	1:C:254:VAL:O	2.37	0.41
1:C:814:SER:OG	1:C:816:ARG:HB2	2.21	0.41
1:B:39:GLU:HG3	1:B:70:ARG:HG2	2.03	0.41
1:B:181:VAL:CG2	1:B:222:VAL:HG12	2.49	0.41
1:B:747:LEU:N	1:B:747:LEU:CD1	2.79	0.41
1:B:825:ALA:HA	1:B:888:MET:SD	2.60	0.41
1:B:881:ASP:OD1	1:B:882:LEU:HD12	2.19	0.41
1:C:472:GLU:OE2	1:C:472:GLU:HA	2.20	0.41
1:C:703:TRP:CH2	1:C:709:ARG:NH1	2.87	0.41
1:C:851:ARG:O	1:C:853:LEU:N	2.53	0.41
1:C:984:PHE:CD1	1:C:989:VAL:HB	2.55	0.41
1:B:309:LYS:C	1:B:311:ASP:H	2.28	0.41
1:B:358:ILE:HG23	1:B:404:MET:SD	2.60	0.41
1:B:361:ILE:HD13	1:B:400:LEU:HB2	2.03	0.41
1:B:674:ARG:HH11	1:B:674:ARG:HG3	1.85	0.41
1:B:696:PHE:CD1	1:B:696:PHE:C	2.99	0.41
1:B:784:ASN:ND2	1:B:958:GLN:HB2	2.35	0.41
1:B:1025:LEU:HD23	1:B:1025:LEU:C	2.45	0.41
1:C:285:GLY:O	1:C:288:ALA:HB3	2.19	0.41
1:C:320:ILE:HG22	1:C:321:ASP:N	2.36	0.41
1:C:381:GLU:CD	1:C:381:GLU:N	2.78	0.41
1:C:958:GLN:O	1:C:962:ILE:HG12	2.21	0.41
1:C:965:MET:HB3	1:C:970:ILE:O	2.19	0.41
1:B:271:LEU:C	1:B:273:LYS:N	2.79	0.41
1:B:329:TYR:O	1:B:330:TYR:C	2.64	0.41
1:B:426:LEU:CD1	1:B:466:ALA:HA	2.48	0.41
1:B:719:LEU:O	1:B:720:VAL:C	2.64	0.41
1:B:756:LEU:CD1	1:B:904:ILE:HB	2.51	0.41
1:C:66:LYS:HB3	1:C:70:ARG:NH1	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:266:THR:O	1:C:267:LEU:CB	2.68	0.41
1:C:336:GLY:O	1:C:337:LEU:HB3	2.21	0.41
1:C:757:THR:O	1:C:904:ILE:HG22	2.20	0.41
1:B:38:PRO:O	1:B:40:ASP:N	2.54	0.41
1:B:320:ILE:HG22	1:B:321:ASP:N	2.34	0.41
1:B:351:CYS:HA	1:B:451:LEU:HD11	2.03	0.41
1:B:514:PRO:O	1:B:515:THR:C	2.63	0.41
1:B:558:ASP:OD2	1:B:559:LEU:N	2.49	0.41
1:B:687:ARG:NH2	1:B:1026:GLU:OE1	2.49	0.41
1:B:774:TYR:HH	1:B:782:ASP:CG	2.28	0.41
1:B:975:THR:O	1:B:976:TYR:C	2.60	0.41
1:B:1002:GLY:O	1:B:1005:VAL:N	2.53	0.41
1:C:187:ILE:C	1:C:189:LYS:N	2.78	0.41
1:C:426:LEU:HD13	1:C:469:TYR:CD1	2.55	0.41
1:C:458:LEU:O	1:C:459:LEU:C	2.62	0.41
1:C:616:LEU:O	1:C:617:ARG:C	2.63	0.41
1:C:790:SER:CA	1:C:958:GLN:HE22	2.28	0.41
1:B:106:PHE:CD1	1:B:182:ASP:HB2	2.56	0.41
1:B:109:SER:HB3	1:B:138:HIS:CD2	2.55	0.41
1:B:230:GLY:C	1:B:232:LYS:H	2.29	0.41
1:B:388:ALA:O	1:B:389:VAL:HB	2.21	0.41
1:B:628:VAL:HG11	1:B:640:ALA:HB2	2.02	0.41
1:B:712:SER:CB	1:B:986:ARG:HH11	2.34	0.41
1:B:822:GLN:HE21	1:B:822:GLN:HB3	1.65	0.41
1:B:828:TYR:HB3	1:B:888:MET:HE3	2.02	0.41
1:B:1007:ARG:CA	1:B:1010:VAL:HG12	2.50	0.41
1:C:46:PHE:CZ	1:C:91:MET:HB2	2.56	0.41
1:C:76:ALA:N	1:C:222:VAL:O	2.52	0.41
1:C:112:VAL:HG21	1:C:162:THR:HG23	2.00	0.41
1:C:184:VAL:CG1	1:C:225:ALA:HA	2.50	0.41
1:C:382:ILE:O	1:C:386:LEU:CD2	2.68	0.41
1:C:419:GLY:O	1:C:420:GLU:CB	2.65	0.41
1:C:691:ASP:O	1:C:695:GLY:HA3	2.20	0.41
1:C:725:ILE:C	1:C:727:ARG:N	2.78	0.41
1:C:810:HIS:O	1:C:847:ILE:HD11	2.21	0.41
1:C:958:GLN:C	1:C:961:ILE:HG22	2.45	0.41
1:B:140:ARG:HB3	1:B:141:ILE:H	1.63	0.41
1:B:171:ARG:HG3	1:B:171:ARG:NH1	2.36	0.41
1:B:289:GLU:C	1:B:291:ILE:N	2.78	0.41
1:B:330:TYR:HA	1:B:333:LEU:HD23	2.03	0.41
1:B:466:ALA:O	1:B:467:LYS:C	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:497:ARG:O	1:B:500:GLU:CG	2.68	0.41
1:B:545:TYR:CE1	1:B:670:LEU:HD13	2.56	0.41
1:B:556:VAL:HB	1:B:613:ILE:HD12	1.96	0.41
1:B:637:GLU:HG2	1:B:680:LEU:HB2	2.03	0.41
1:B:698:LEU:HD22	1:B:1009:LEU:HD12	2.03	0.41
1:B:700:GLN:C	1:B:702:LEU:N	2.79	0.41
1:B:745:LEU:CD1	1:B:747:LEU:HD11	2.41	0.41
1:B:849:PRO:CG	1:B:884:PHE:HA	2.47	0.41
1:B:870:GLU:O	1:B:872:LEU:HG	2.21	0.41
1:B:904:ILE:O	1:B:911:ALA:N	2.54	0.41
1:B:1007:ARG:C	1:B:1010:VAL:HG12	2.46	0.41
1:C:58:ARG:O	1:C:59:ALA:C	2.64	0.41
1:C:171:ARG:HG3	1:C:171:ARG:NH1	2.36	0.41
1:C:360:ASP:O	1:C:361:ILE:C	2.63	0.41
1:C:425:ASP:OD2	1:C:428:THR:CB	2.69	0.41
1:C:466:ALA:O	1:C:467:LYS:C	2.63	0.41
1:C:497:ARG:O	1:C:500:GLU:HB2	2.21	0.41
1:C:628:VAL:HG13	1:C:630:THR:HG23	2.02	0.41
1:C:681:VAL:O	1:C:685:VAL:HG23	2.20	0.41
1:C:791:VAL:HG11	1:C:980:VAL:HG11	2.01	0.41
1:C:877:PHE:O	1:C:878:ALA:C	2.62	0.41
1:C:1022:THR:O	1:C:1026:GLU:HB2	2.20	0.41
1:B:279:ILE:HD11	1:B:315:PHE:CZ	2.55	0.41
1:B:337:LEU:HB3	1:B:435:ARG:HE	1.85	0.41
1:B:365:SER:HA	1:B:366:PRO:HD3	1.82	0.41
1:B:581:ILE:HG22	1:B:582:LYS:H	1.86	0.41
1:B:639:ILE:O	1:B:640:ALA:C	2.62	0.41
1:B:791:VAL:HG23	1:B:977:ALA:CB	2.52	0.41
1:B:962:ILE:HD13	1:B:962:ILE:HA	1.87	0.41
1:C:167:SER:OG	1:C:193:ASN:OD1	2.37	0.41
1:C:208:LYS:C	1:C:210:LYS:N	2.78	0.41
1:C:440:PHE:HD1	1:C:440:PHE:H	1.68	0.41
1:C:498:ARG:C	1:C:500:GLU:N	2.79	0.41
1:C:624:GLU:HA	1:C:650:CYS:O	2.20	0.41
1:C:669:ALA:C	1:C:671:GLU:H	2.29	0.41
1:C:736:LYS:O	1:C:948:LYS:O	2.40	0.41
1:C:808:THR:HG21	1:C:848:ARG:NE	2.36	0.41
1:C:901:LYS:HA	1:C:914:GLU:HA	2.02	0.41
1:C:906:PHE:O	1:C:907:ASP:C	2.63	0.41
1:B:124:GLU:O	1:B:125:LYS:HD2	2.21	0.40
1:B:279:ILE:CG2	1:B:281:TYR:CE1	2.98	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:314:LYS:HG2	1:B:319:GLU:CD	2.46	0.40
1:B:323:LEU:HB3	1:B:330:TYR:OH	2.21	0.40
1:B:336:GLY:O	1:B:337:LEU:HB3	2.21	0.40
1:B:403:VAL:O	1:B:404:MET:CB	2.58	0.40
1:B:413:ASP:O	1:B:424:PRO:HG2	2.21	0.40
1:B:754:PHE:CE2	1:B:907:ASP:N	2.90	0.40
1:C:261:ASP:CG	1:C:262:GLU:N	2.79	0.40
1:C:374:TYR:O	1:C:375:LEU:HB2	2.22	0.40
1:C:743:PHE:CD1	1:C:743:PHE:N	2.88	0.40
1:C:895:PHE:CD2	1:C:895:PHE:N	2.82	0.40
1:C:1018:SER:O	1:C:1021:ARG:N	2.54	0.40
1:B:338:ASP:OD1	1:B:340:PRO:HD3	2.21	0.40
1:B:377:ARG:O	1:B:377:ARG:HG2	2.21	0.40
1:B:637:GLU:HG2	1:B:680:LEU:HD12	2.04	0.40
1:B:758:ILE:HB	1:B:938:THR:HA	2.03	0.40
1:B:942:ARG:H	1:B:942:ARG:HG2	1.66	0.40
1:B:991:GLU:O	1:B:992:LYS:HE2	2.21	0.40
1:C:35:CYS:HB2	1:C:650:CYS:N	2.35	0.40
1:C:289:GLU:O	1:C:293:GLU:OE2	2.40	0.40
1:C:737:ILE:HG23	1:C:950:VAL:HG22	2.03	0.40
1:C:966:LYS:O	1:C:967:GLU:C	2.64	0.40
1:B:152:ASN:O	1:B:152:ASN:ND2	2.55	0.40
1:B:266:THR:O	1:B:267:LEU:CB	2.70	0.40
1:B:276:THR:HA	1:B:322:HIS:NE2	2.37	0.40
1:B:433:SER:O	1:B:436:THR:CG2	2.68	0.40
1:B:542:MET:HE3	1:B:547:LEU:HD11	2.02	0.40
1:B:878:ALA:O	1:B:881:ASP:HB3	2.20	0.40
1:C:278:GLY:HA2	1:C:345:PHE:O	2.21	0.40
1:C:901:LYS:HG3	1:C:902:TYR:H	1.86	0.40
1:C:1025:LEU:HD23	1:C:1025:LEU:O	2.21	0.40
1:B:106:PHE:CE2	1:B:115:ALA:HB2	2.57	0.40
1:B:195:ASP:O	1:B:196:LYS:C	2.64	0.40
1:B:387:PRO:CA	1:B:389:VAL:H	2.32	0.40
1:B:438:ARG:O	1:B:444:LEU:HA	2.22	0.40
1:B:1053:ILE:HG13	1:B:1054:ASP:N	2.35	0.40
1:C:243:ASP:HB3	1:C:529:VAL:HB	2.02	0.40
1:C:267:LEU:HD12	1:C:452:LEU:HD13	2.04	0.40
1:C:296:LYS:HG3	1:C:301:ILE:HD13	2.04	0.40
1:C:337:LEU:CD2	1:C:337:LEU:C	2.94	0.40
1:C:366:PRO:C	1:C:368:MET:H	2.28	0.40
1:C:542:MET:HB2	1:C:543:GLN:HE21	1.87	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1007:ARG:O	1:C:1011:ARG:HG3	2.22	0.40
1:C:1050:ILE:O	1:C:1053:ILE:HG12	2.22	0.40
1:B:305:THR:CG2	1:B:329:TYR:CZ	3.05	0.40
1:B:312:TYR:CE2	1:B:316:VAL:CG2	3.04	0.40
1:B:458:LEU:O	1:B:459:LEU:C	2.62	0.40
1:B:496:ARG:HG3	1:B:496:ARG:HH11	1.86	0.40
1:B:504:ILE:HG22	1:B:546:VAL:HG23	2.02	0.40
1:B:734:ARG:C	1:B:736:LYS:H	2.29	0.40
1:C:162:THR:C	1:C:164:GLN:H	2.29	0.40
1:C:170:TYR:HB3	1:C:196:LYS:HB3	2.03	0.40
1:C:430:ILE:O	1:C:431:GLN:C	2.65	0.40
1:C:444:LEU:O	1:C:445:THR:C	2.64	0.40
1:C:642:ASP:HB3	1:C:646:LEU:HD11	2.03	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	B	1010/1054 (96%)	752 (74%)	177 (18%)	81 (8%)	1	4
1	C	987/1054 (94%)	696 (70%)	214 (22%)	77 (8%)	1	5
All	All	1997/2108 (95%)	1448 (72%)	391 (20%)	158 (8%)	1	5

All (158) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	3	PRO
1	B	5	VAL
1	B	12	VAL
1	B	24	LYS
1	B	32	ARG

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Mol	Chain	Res	Type
1	B	128	VAL
1	B	141	ILE
1	B	154	ARG
1	B	213	VAL
1	B	266	THR
1	B	267	LEU
1	B	306	ALA
1	B	307	THR
1	B	361	ILE
1	B	362	ASP
1	B	364	LEU
1	B	387	PRO
1	B	420	GLU
1	B	445	THR
1	B	471	ILE
1	B	476	ILE
1	B	481	PHE
1	B	947	VAL
1	C	4	VAL
1	C	154	ARG
1	C	213	VAL
1	C	244	ILE
1	C	254	VAL
1	C	266	THR
1	C	267	LEU
1	C	305	THR
1	C	306	ALA
1	C	307	THR
1	C	361	ILE
1	C	362	ASP
1	C	364	LEU
1	C	379	VAL
1	C	420	GLU
1	C	471	ILE
1	C	481	PHE
1	C	533	ASP
1	C	543	GLN
1	C	580	SER
1	C	630	THR
1	C	707	ASN
1	C	737	ILE
1	C	947	VAL

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Mol	Chain	Res	Type
1	B	130	THR
1	B	139	GLY
1	B	194	VAL
1	B	228	LYS
1	B	254	VAL
1	B	377	ARG
1	B	392	HIS
1	B	404	MET
1	B	419	GLY
1	B	470	ASP
1	B	502	ASP
1	B	533	ASP
1	B	709	ARG
1	B	736	LYS
1	B	737	ILE
1	B	830	GLY
1	B	838	TRP
1	B	854	THR
1	B	871	GLY
1	B	908	GLY
1	C	128	VAL
1	C	133	LEU
1	C	139	GLY
1	C	208	LYS
1	C	262	GLU
1	C	374	TYR
1	C	377	ARG
1	C	378	ASN
1	C	564	GLY
1	C	715	ARG
1	C	732	ARG
1	C	752	GLU
1	C	830	GLY
1	C	838	TRP
1	C	871	GLY
1	C	1013	TYR
1	B	173	LEU
1	B	378	ASN
1	B	455	ASP
1	B	456	SER
1	B	563	ARG
1	B	707	ASN

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Mol	Chain	Res	Type
1	B	752	GLU
1	B	839	GLY
1	B	982	ARG
1	B	1052	SER
1	B	1053	ILE
1	C	76	ALA
1	C	130	THR
1	C	502	ASP
1	C	709	ARG
1	C	770	PRO
1	C	777	GLU
1	C	846	CYS
1	C	982	ARG
1	B	39	GLU
1	B	101	ARG
1	B	165	PHE
1	B	195	ASP
1	B	196	LYS
1	B	208	LYS
1	B	305	THR
1	B	374	TYR
1	B	411	ALA
1	B	417	ARG
1	B	721	LEU
1	B	726	ASP
1	B	805	GLY
1	C	101	ARG
1	C	141	ILE
1	C	249	ILE
1	C	283	ARG
1	C	417	ARG
1	C	456	SER
1	C	507	ALA
1	C	744	ASP
1	B	133	LEU
1	B	249	ILE
1	B	379	VAL
1	B	564	GLY
1	B	715	ARG
1	B	786	ILE
1	B	860	ARG
1	C	5	VAL

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Mol	Chain	Res	Type
1	C	411	ALA
1	C	519	GLN
1	C	570	VAL
1	C	736	LYS
1	C	1053	ILE
1	B	706	PHE
1	B	735	ARG
1	C	12	VAL
1	C	153	LEU
1	C	173	LEU
1	C	194	VAL
1	C	252	ARG
1	C	337	LEU
1	C	470	ASP
1	C	842	GLY
1	C	54	VAL
1	C	666	ILE
1	B	842	GLY
1	C	83	GLY
1	B	570	VAL
1	C	839	GLY
1	B	463	ILE
1	B	835	GLY
1	C	686	VAL
1	B	685	VAL
1	C	142	PRO
1	C	463	ILE

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	876/925 (95%)	772 (88%)	104 (12%)	5	23
1	C	863/925 (93%)	772 (90%)	91 (10%)	6	28
All	All	1739/1850 (94%)	1544 (89%)	195 (11%)	6	25

All (195) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	32	ARG
1	B	34	LEU
1	B	42	LEU
1	B	53	CYS
1	B	56	GLU
1	B	57	PRO
1	B	60	ILE
1	B	71	LYS
1	B	77	THR
1	B	92	SER
1	B	95	LEU
1	B	119	ILE
1	B	120	ARG
1	B	134	ILE
1	B	140	ARG
1	B	150	MET
1	B	158	ILE
1	B	161	THR
1	B	164	GLN
1	B	208	LYS
1	B	212	TRP
1	B	213	VAL
1	B	224	THR
1	B	232	LYS
1	B	234	GLU
1	B	235	LEU
1	B	239	LEU
1	B	252	ARG
1	B	264	ILE
1	B	267	LEU
1	B	279	ILE
1	B	295	LEU
1	B	301	ILE
1	B	303	ILE
1	B	309	LYS
1	B	320	ILE
1	B	328	HIS
1	B	337	LEU
1	B	354	PHE
1	B	364	LEU
1	B	381	GLU
1	B	384	ARG

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Mol	Chain	Res	Type
1	B	385	LEU
1	B	386	LEU
1	B	387	PRO
1	B	391	ARG
1	B	392	HIS
1	B	430	ILE
1	B	436	THR
1	B	440	PHE
1	B	446	LYS
1	B	452	LEU
1	B	467	LYS
1	B	471	ILE
1	B	472	GLU
1	B	474	LYS
1	B	479	VAL
1	B	494	ARG
1	B	510	ILE
1	B	518	ARG
1	B	520	ILE
1	B	543	GLN
1	B	560	ILE
1	B	606	VAL
1	B	612	ARG
1	B	627	ILE
1	B	666	ILE
1	B	670	LEU
1	B	672	SER
1	B	694	ILE
1	B	708	ASN
1	B	724	ILE
1	B	725	ILE
1	B	736	LYS
1	B	742	ASP
1	B	745	LEU
1	B	747	LEU
1	B	751	GLU
1	B	756	LEU
1	B	760	LEU
1	B	764	ARG
1	B	767	LEU
1	B	768	ARG
1	B	777	GLU

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Mol	Chain	Res	Type
1	B	791	VAL
1	B	796	GLN
1	B	801	LEU
1	B	811	ARG
1	B	848	ARG
1	B	895	PHE
1	B	916	ILE
1	B	931	TRP
1	B	938	THR
1	B	943	VAL
1	B	961	ILE
1	B	965	MET
1	B	970	ILE
1	B	980	VAL
1	B	1009	LEU
1	B	1012	ARG
1	B	1017	VAL
1	B	1027	SER
1	B	1032	ILE
1	B	1050	ILE
1	C	32	ARG
1	C	34	LEU
1	C	39	GLU
1	C	53	CYS
1	C	60	ILE
1	C	61	GLN
1	C	71	LYS
1	C	77	THR
1	C	120	ARG
1	C	134	ILE
1	C	140	ARG
1	C	150	MET
1	C	155	ASN
1	C	158	ILE
1	C	161	THR
1	C	164	GLN
1	C	198	LEU
1	C	204	HIS
1	C	212	TRP
1	C	213	VAL
1	C	215	GLU
1	C	217	ARG

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Mol	Chain	Res	Type
1	C	221	MET
1	C	224	THR
1	C	232	LYS
1	C	234	GLU
1	C	235	LEU
1	C	249	ILE
1	C	252	ARG
1	C	264	ILE
1	C	267	LEU
1	C	295	LEU
1	C	301	ILE
1	C	305	THR
1	C	309	LYS
1	C	320	ILE
1	C	328	HIS
1	C	337	LEU
1	C	354	PHE
1	C	364	LEU
1	C	377	ARG
1	C	381	GLU
1	C	384	ARG
1	C	385	LEU
1	C	386	LEU
1	C	413	ASP
1	C	430	ILE
1	C	436	THR
1	C	440	PHE
1	C	446	LYS
1	C	479	VAL
1	C	494	ARG
1	C	499	GLN
1	C	533	ASP
1	C	543	GLN
1	C	549	VAL
1	C	606	VAL
1	C	639	ILE
1	C	672	SER
1	C	705	ARG
1	C	709	ARG
1	C	736	LYS
1	C	737	ILE
1	C	742	ASP

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Mol	Chain	Res	Type
1	C	747	LEU
1	C	751	GLU
1	C	753	GLU
1	C	756	LEU
1	C	760	LEU
1	C	767	LEU
1	C	768	ARG
1	C	782	ASP
1	C	791	VAL
1	C	796	GLN
1	C	801	LEU
1	C	811	ARG
1	C	823	ARG
1	C	848	ARG
1	C	860	ARG
1	C	873	ARG
1	C	875	GLU
1	C	895	PHE
1	C	922	ARG
1	C	934	ASN
1	C	936	LEU
1	C	938	THR
1	C	943	VAL
1	C	989	VAL
1	C	1025	LEU
1	C	1032	ILE
1	C	1050	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (55) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	61	GLN
1	B	114	GLN
1	B	138	HIS
1	B	148	ASN
1	B	151	GLN
1	B	152	ASN
1	B	164	GLN
1	B	193	ASN
1	B	253	ASN
1	B	297	ASN
1	B	322	HIS

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Mol	Chain	Res	Type
1	B	431	GLN
1	B	519	GLN
1	B	566	HIS
1	B	608	ASN
1	B	621	HIS
1	B	708	ASN
1	B	710	ASN
1	B	784	ASN
1	B	793	GLN
1	B	796	GLN
1	B	799	GLN
1	B	804	ASN
1	B	822	GLN
1	B	876	HIS
1	B	934	ASN
1	B	958	GLN
1	B	963	GLN
1	C	61	GLN
1	C	138	HIS
1	C	148	ASN
1	C	151	GLN
1	C	152	ASN
1	C	164	GLN
1	C	169	HIS
1	C	175	HIS
1	C	253	ASN
1	C	431	GLN
1	C	499	GLN
1	C	519	GLN
1	C	543	GLN
1	C	608	ASN
1	C	621	HIS
1	C	684	GLN
1	C	708	ASN
1	C	710	ASN
1	C	784	ASN
1	C	793	GLN
1	C	796	GLN
1	C	804	ASN
1	C	822	GLN
1	C	876	HIS
1	C	934	ASN

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Mol	Chain	Res	Type
1	C	958	GLN
1	C	963	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	ANP	B	2055	-	33,33,33	1.86	7 (21%)	45,52,52	2.69	16 (35%)
2	ANP	C	2055	-	33,33,33	2.32	13 (39%)	45,52,52	2.64	15 (33%)

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	ANP	B	2055	-	1/1/7/8	4/18/38/38	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ANP	C	2055	-	-	6/18/38/38	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2055	ANP	PB-O3A	7.95	1.69	1.59
2	B	2055	ANP	PB-O3A	4.39	1.64	1.59
2	B	2055	ANP	PB-O2B	-4.03	1.46	1.56
2	B	2055	ANP	PG-O1G	3.91	1.52	1.46
2	C	2055	ANP	PB-O2B	-3.82	1.46	1.56
2	C	2055	ANP	PG-O1G	3.80	1.51	1.46
2	C	2055	ANP	PB-O1B	3.35	1.51	1.46
2	C	2055	ANP	PA-O3A	2.91	1.62	1.59
2	B	2055	ANP	C5-N7	-2.79	1.34	1.39
2	C	2055	ANP	C5-N7	-2.49	1.34	1.39
2	B	2055	ANP	PG-N3B	2.42	1.69	1.63
2	C	2055	ANP	PG-N3B	2.36	1.69	1.63
2	B	2055	ANP	C2'-C1'	-2.29	1.46	1.53
2	C	2055	ANP	C2-N1	2.22	1.37	1.33
2	B	2055	ANP	PB-O1B	2.19	1.49	1.46
2	C	2055	ANP	PB-N3B	-2.18	1.57	1.63
2	C	2055	ANP	C2'-C1'	-2.09	1.46	1.53
2	C	2055	ANP	C1'-N9	2.06	1.52	1.46
2	C	2055	ANP	C4-N3	2.05	1.38	1.34
2	C	2055	ANP	PG-O3G	-2.03	1.51	1.56

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2055	ANP	C5-C4-N3	-7.30	116.66	126.72
2	C	2055	ANP	C5-C4-N3	-7.25	116.74	126.72
2	B	2055	ANP	N3-C4-N9	6.28	137.86	127.17
2	C	2055	ANP	N3-C4-N9	6.22	137.74	127.17
2	C	2055	ANP	O1B-PB-N3B	-5.55	103.59	111.77
2	B	2055	ANP	O1B-PB-N3B	-5.46	103.73	111.77
2	B	2055	ANP	O1G-PG-N3B	-5.43	103.78	111.77
2	C	2055	ANP	O1G-PG-N3B	-5.36	103.87	111.77
2	C	2055	ANP	O4'-C1'-N9	5.30	118.27	108.09
2	B	2055	ANP	O4'-C1'-N9	5.21	118.09	108.09
2	B	2055	ANP	C4'-O4'-C1'	4.80	120.07	109.47
2	C	2055	ANP	C4'-O4'-C1'	4.55	119.52	109.47
2	B	2055	ANP	C2-N3-C4	3.94	121.46	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2055	ANP	C2-N3-C4	3.92	121.40	111.83
2	B	2055	ANP	O3A-PB-N3B	3.88	117.35	106.59
2	C	2055	ANP	O3A-PB-N3B	3.86	117.30	106.59
2	B	2055	ANP	C3'-C2'-C1'	3.82	108.69	101.46
2	B	2055	ANP	C6-C5-C4	3.31	121.70	117.18
2	C	2055	ANP	C6-C5-C4	3.18	121.52	117.18
2	C	2055	ANP	O4'-C1'-C2'	-2.93	100.34	106.62
2	B	2055	ANP	O4'-C1'-C2'	-2.77	100.68	106.62
2	C	2055	ANP	C3'-C2'-C1'	2.56	106.31	101.46
2	B	2055	ANP	N3-C2-N1	-2.56	124.71	128.58
2	C	2055	ANP	N3-C2-N1	-2.51	124.79	128.58
2	C	2055	ANP	O4'-C4'-C3'	-2.29	100.60	105.15
2	C	2055	ANP	O2G-PG-O3G	2.17	113.43	107.59
2	C	2055	ANP	O3'-C3'-C4'	-2.16	104.88	111.08
2	B	2055	ANP	O3'-C3'-C4'	-2.16	104.88	111.08
2	B	2055	ANP	O2G-PG-O3G	2.12	113.29	107.59
2	B	2055	ANP	O3'-C3'-C2'	-2.09	105.12	111.82
2	B	2055	ANP	C6-C5-N7	-2.02	128.20	132.09

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	2055	ANP	C2'

All (10) torsion outliers are listed below:

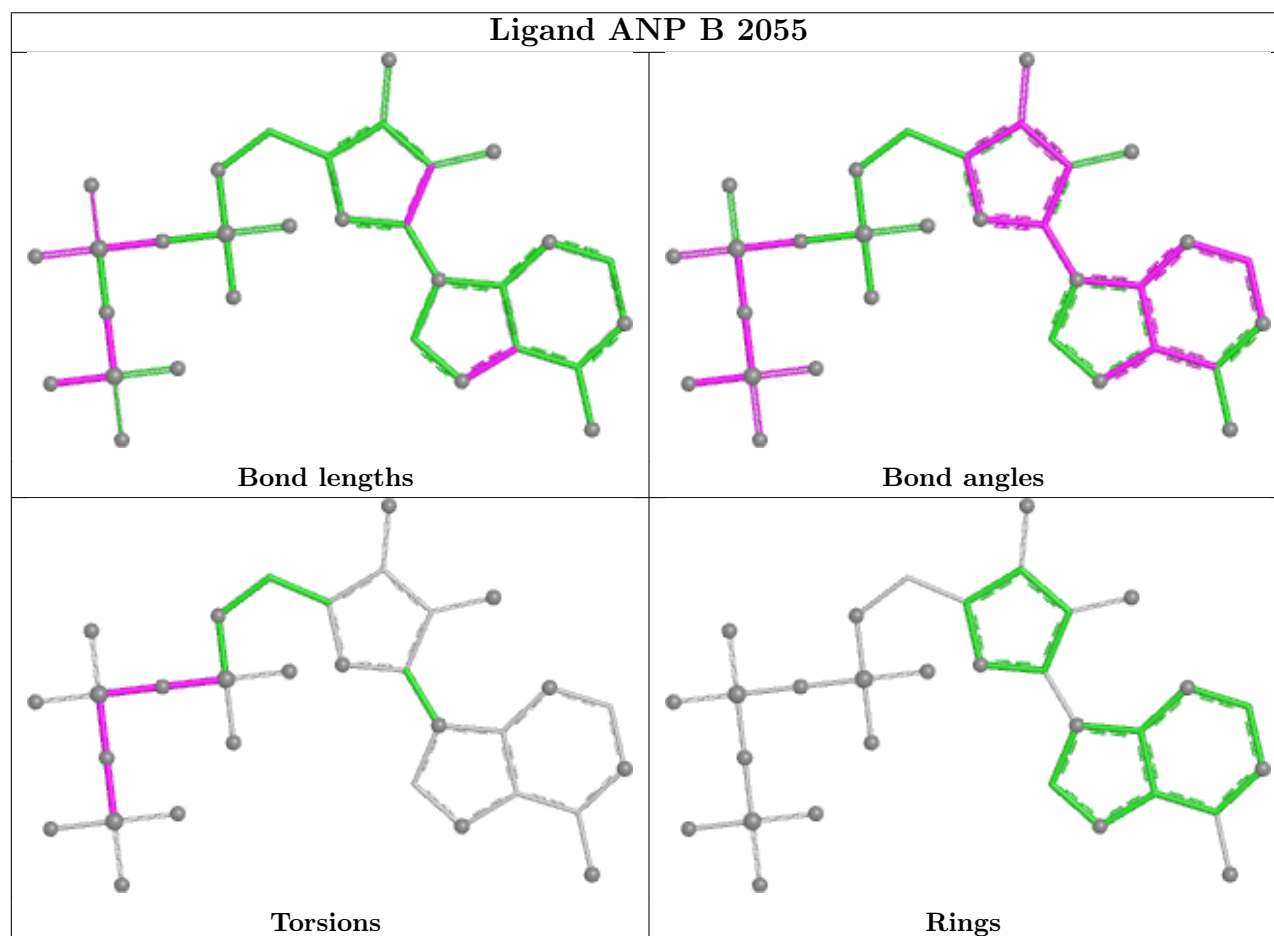
Mol	Chain	Res	Type	Atoms
2	B	2055	ANP	PB-N3B-PG-O1G
2	C	2055	ANP	PB-N3B-PG-O1G
2	C	2055	ANP	PA-O3A-PB-O2B
2	C	2055	ANP	C3'-C4'-C5'-O5'
2	C	2055	ANP	PB-O3A-PA-O2A
2	B	2055	ANP	PG-N3B-PB-O1B
2	B	2055	ANP	PA-O3A-PB-O1B
2	C	2055	ANP	PA-O3A-PB-O1B
2	B	2055	ANP	PB-O3A-PA-O1A
2	C	2055	ANP	PB-O3A-PA-O1A

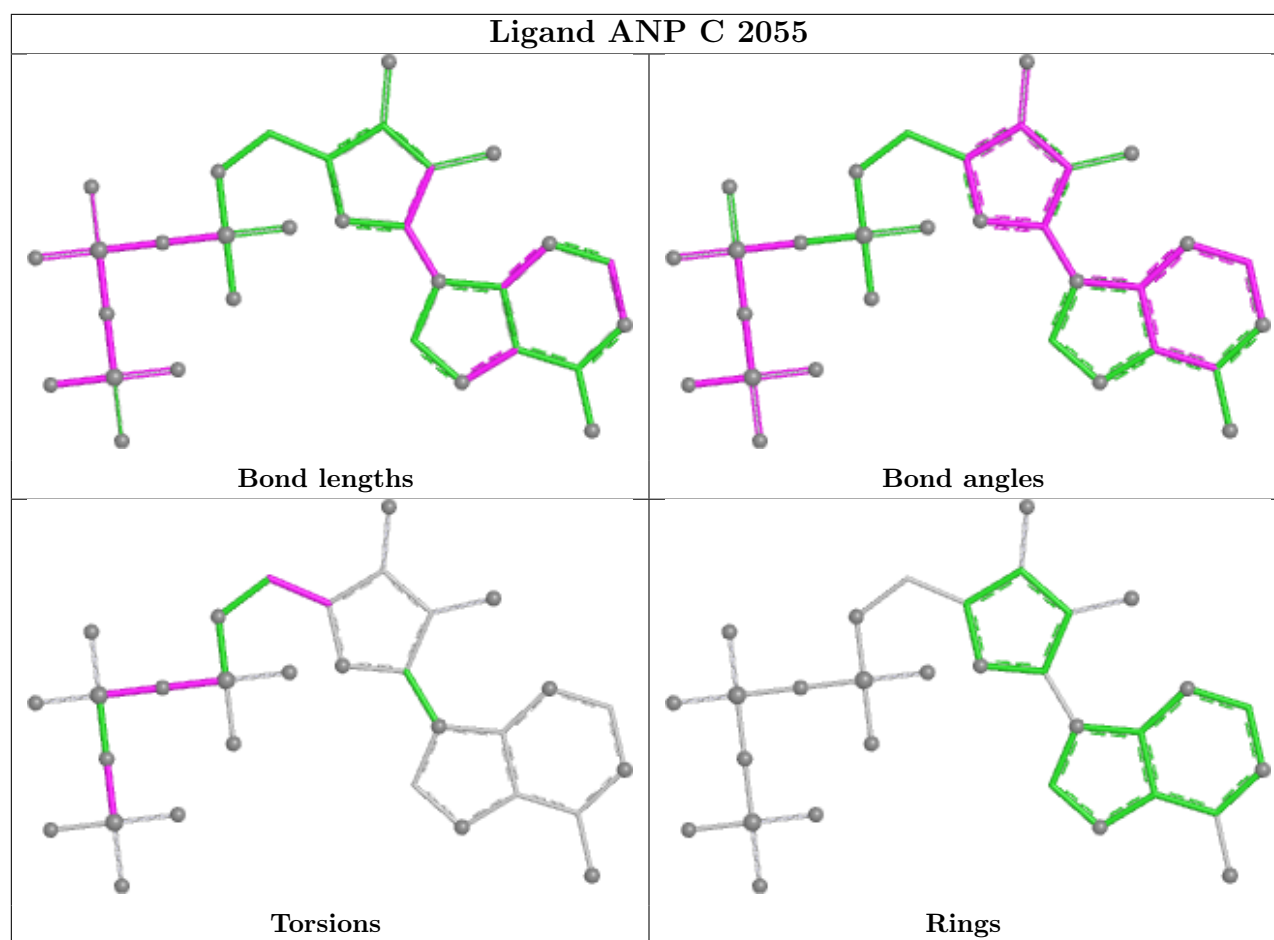
There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2055	ANP	5	0
2	C	2055	ANP	5	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	C	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	15:GLY	C	16:ASP	N	13.42

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	B	1020/1054 (96%)	-0.16	26 (2%) 58 39	1, 38, 137, 192	0
1	C	1005/1054 (95%)	0.39	28 (2%) 55 35	24, 92, 160, 195	0
All	All	2025/2108 (96%)	0.11	54 (2%) 56 36	1, 68, 153, 195	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	706	PHE	4.3
1	B	937	PRO	4.2
1	B	212	TRP	4.0
1	B	384	ARG	3.9
1	B	376	TYR	3.9
1	C	965	MET	3.6
1	C	301	ILE	3.4
1	B	1013	TYR	3.3
1	C	953	VAL	3.3
1	C	998	PRO	3.2
1	C	471	ILE	3.2
1	B	141	ILE	3.0
1	B	357	THR	2.8
1	B	393	ILE	2.8
1	C	213	VAL	2.7
1	C	929	ALA	2.7
1	B	760	LEU	2.7
1	C	1008	PHE	2.7
1	B	375	LEU	2.7
1	B	387	PRO	2.6
1	B	301	ILE	2.5
1	C	970	ILE	2.5
1	C	381	GLU	2.4
1	C	136	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	142	PRO	2.4
1	B	379	VAL	2.4
1	C	360	ASP	2.3
1	B	381	GLU	2.3
1	B	383	GLU	2.3
1	B	386	LEU	2.3
1	B	471	ILE	2.3
1	C	380	ASP	2.3
1	C	962	ILE	2.2
1	C	717	GLN	2.2
1	B	172	GLU	2.2
1	C	212	TRP	2.2
1	B	186	ALA	2.2
1	C	375	LEU	2.2
1	C	940	THR	2.2
1	C	713	ALA	2.2
1	C	976	TYR	2.2
1	B	385	LEU	2.1
1	C	3	PRO	2.1
1	B	361	ILE	2.1
1	C	385	LEU	2.1
1	B	743	PHE	2.1
1	C	7	SER	2.1
1	B	404	MET	2.1
1	C	352	PRO	2.1
1	C	415	VAL	2.1
1	C	731	SER	2.0
1	B	759	LYS	2.0
1	C	956	PHE	2.0
1	C	702	LEU	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 ⓘ

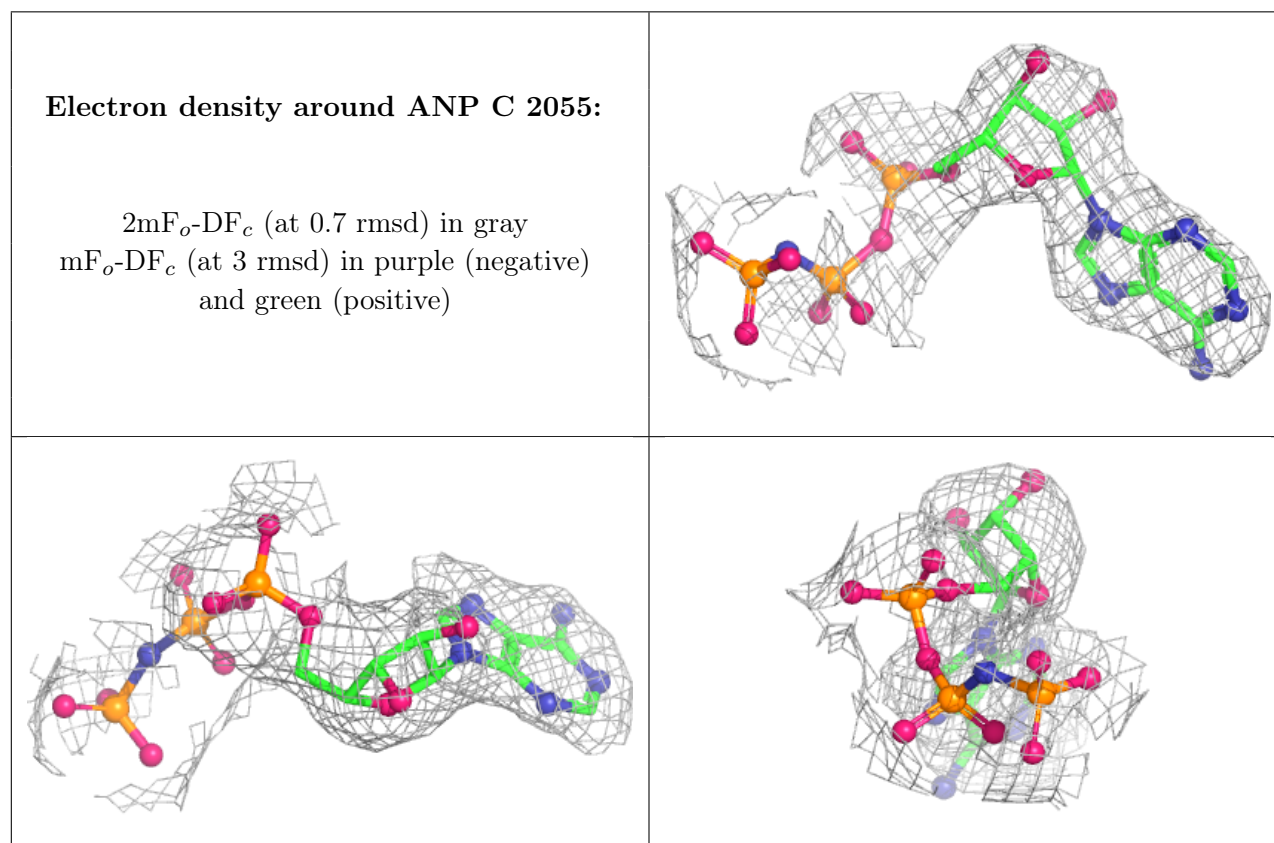
There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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.

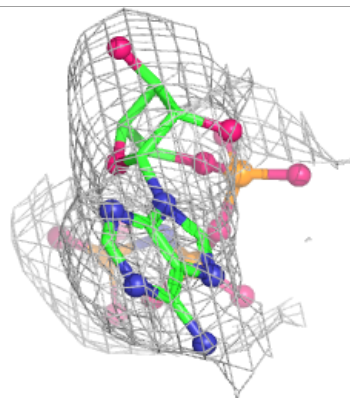
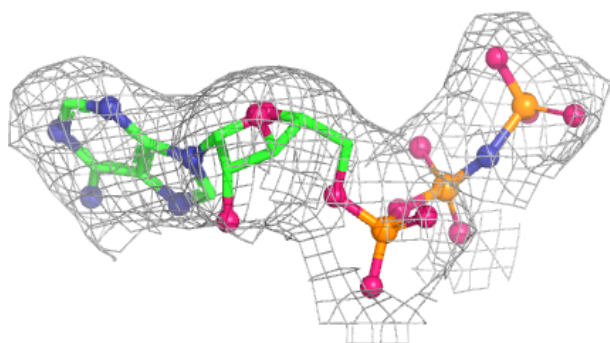
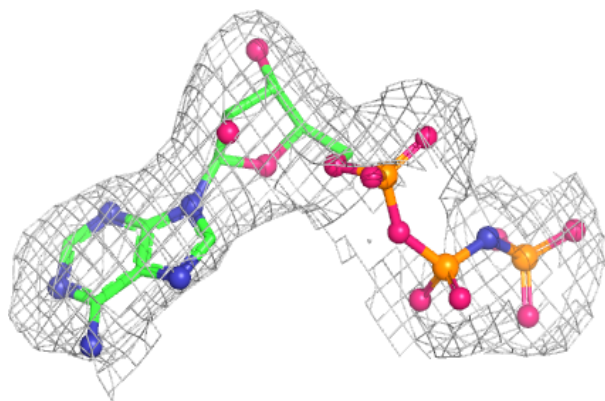
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MG	C	2056	1/1	0.47	0.20	70,70,70,70	0
3	MG	B	2056	1/1	0.80	0.08	40,40,40,40	0
2	ANP	C	2055	31/31	0.88	0.10	38,57,75,77	0
2	ANP	B	2055	31/31	0.94	0.08	24,34,69,77	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around ANP B 2055:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.