



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:09 AM UTC

PDB ID : 1DO0 / pdb_00001do0
Title : ORTHORHOMBIC CRYSTAL FORM OF HEAT SHOCK LOCUS U (HSLU)
FROM ESCHERICHIA COLI
Authors : Bochtler, M.; Hartmann, C.; Song, H.K.; Bourenkov, G.P.; Bartunik, H.D.
Deposited on : 1999-12-18
Resolution : 3.00 Å(reported)

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

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

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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

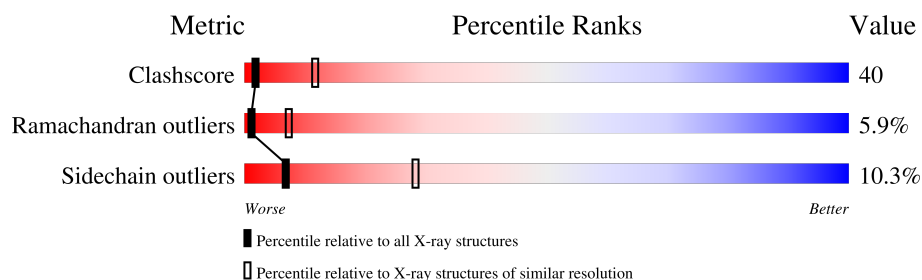
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	442	46% 36% 10% 8%
1	B	442	40% 42% 9% 8%
1	C	442	40% 43% 9% 8%
1	D	442	46% 37% 9% 8%
1	E	442	39% 43% 9% 8%
1	F	442	39% 45% 7% 8%

2 Entry composition [i](#)

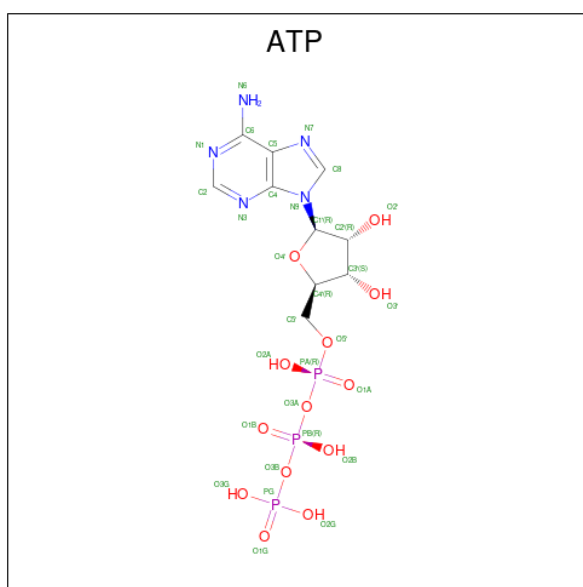
There are 4 unique types of molecules in this entry. The entry contains 19366 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROTEIN (HEAT SHOCK LOCUS U).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	406	Total	C	N	O	S	628	0	0
			3205	2001	570	624	10			
1	B	406	Total	C	N	O	S	143	0	0
			3205	2001	570	624	10			
1	C	406	Total	C	N	O	S	677	0	0
			3205	2001	570	624	10			
1	D	406	Total	C	N	O	S	579	0	0
			3205	2001	570	624	10			
1	E	406	Total	C	N	O	S	135	0	0
			3205	2001	570	624	10			
1	F	406	Total	C	N	O	S	600	0	0
			3205	2001	570	624	10			

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	D	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	E	1	Total	C	N	O	P	0	0
			31	10	5	13	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	E	1	Total	Mg	0	0
			1	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



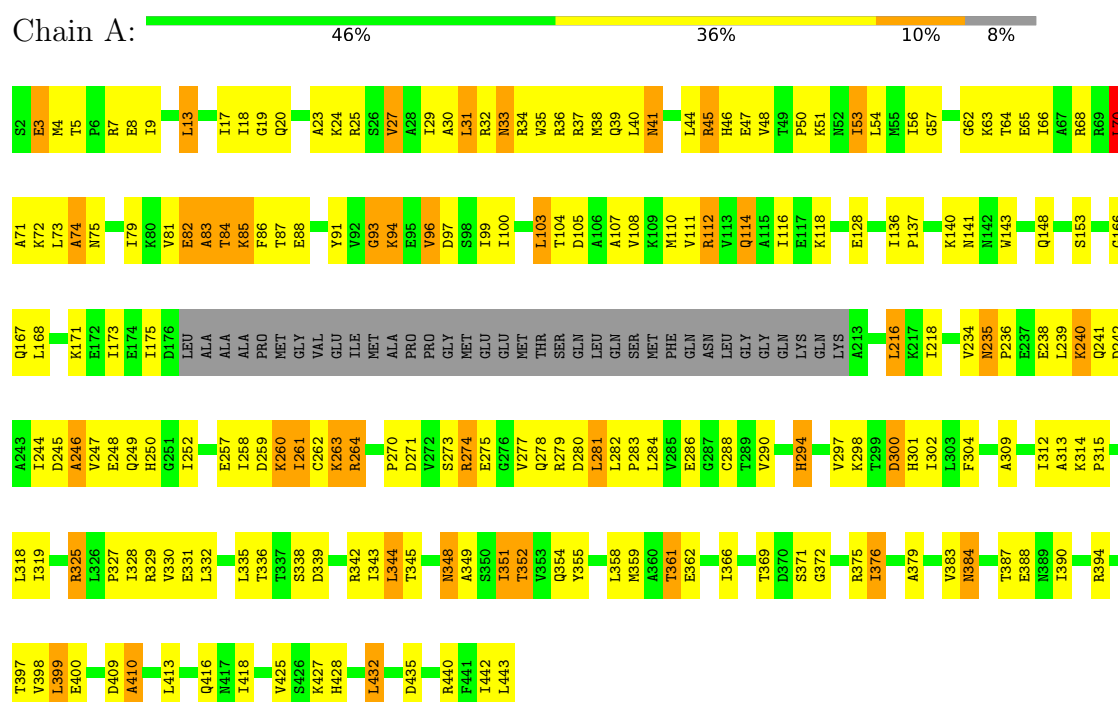
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		

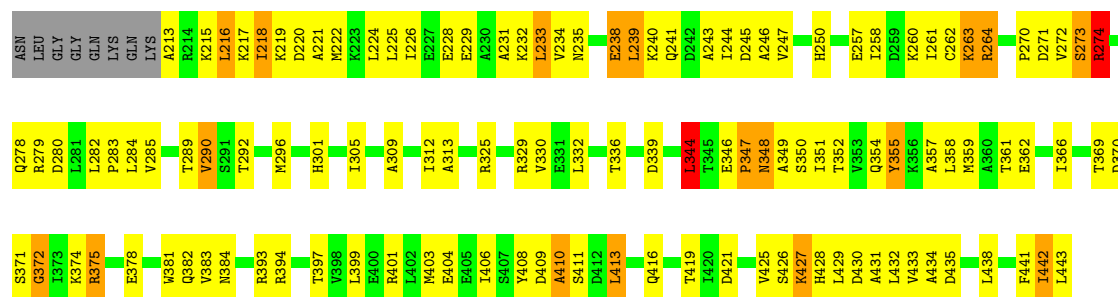
3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

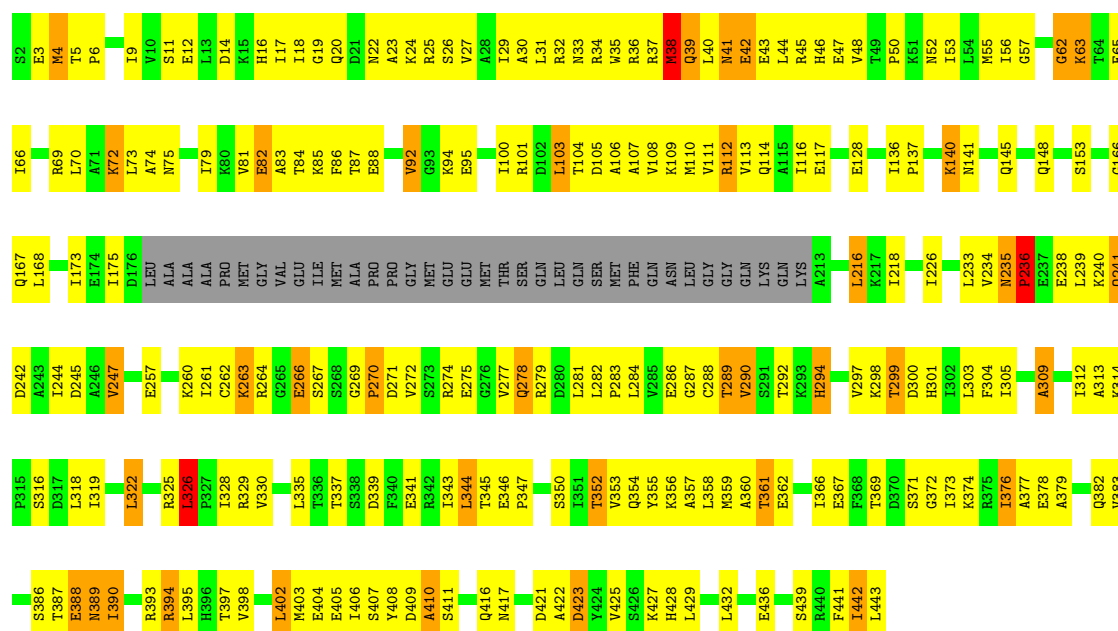
• Molecule 1: PROTEIN (HEAT SHOCK LOCUS U)





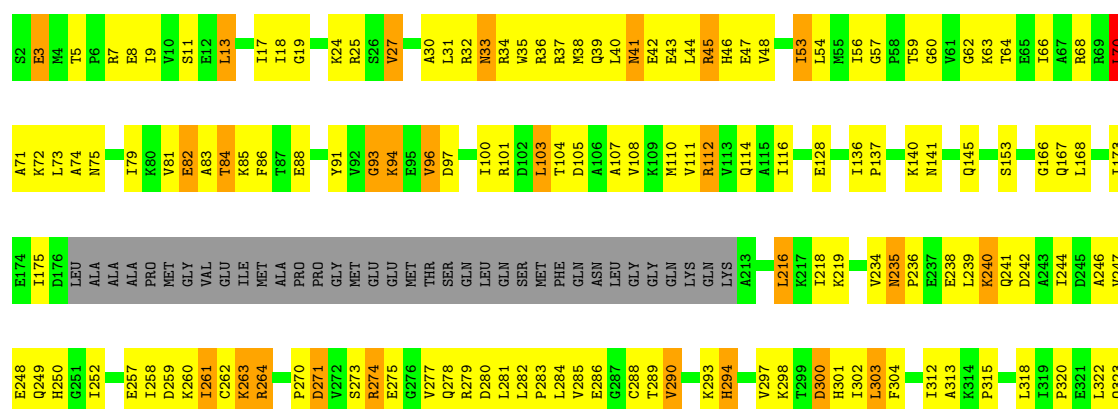
• Molecule 1: PROTEIN (HEAT SHOCK LOCUS U)

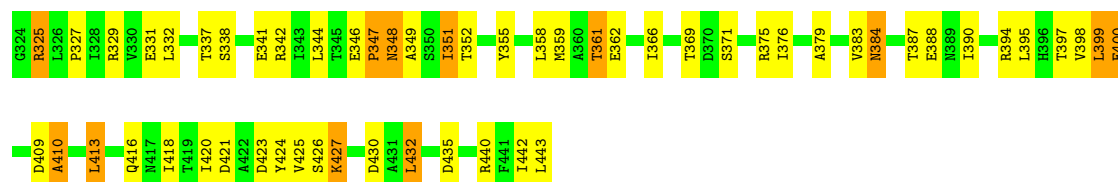
Chain C: 40% 43% 9% 8%



• Molecule 1: PROTEIN (HEAT SHOCK LOCUS U)

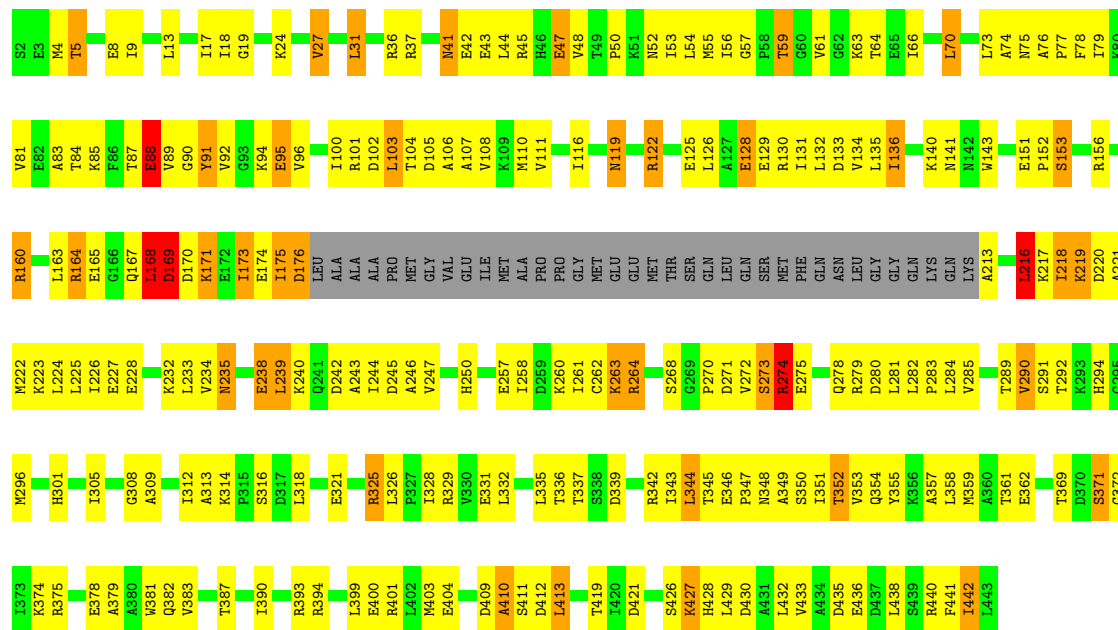
Chain D: 46% 37% 9% 8%





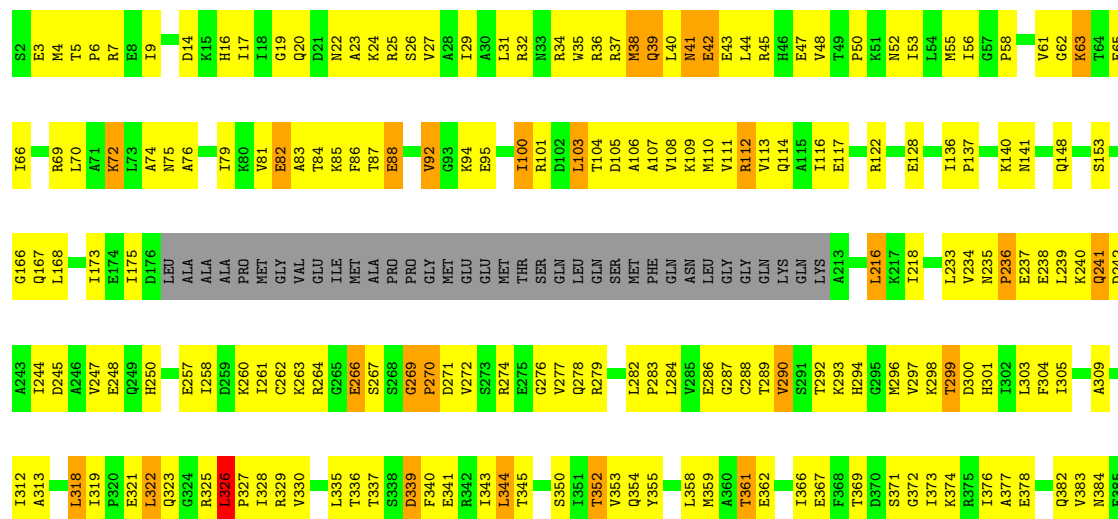
• Molecule 1: PROTEIN (HEAT SHOCK LOCUS U)

Chain E: 39% 43% 9% 8%



• Molecule 1: PROTEIN (HEAT SHOCK LOCUS U)

Chain F: 39% 45% 7% 8%



S386	T387	E388	N389	I390	G391	A392	R393	R394	L395	H396	T397	V398	L399	E400	R401	L402	M403	E404	E405	I406	S407	Y408	D409	A410	S411	Q416	N417	A422	V425	S426	K427	H428	L429	D430	A431	L432	V433	E436	S439	R440	F441	I442	L443
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4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	208.15Å 167.70Å 108.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 3.00	Depositor
% Data completeness (in resolution range)	96.8 (25.00-3.00)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	7.90	Depositor
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.294 , 0.342	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	19366	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, ATP, 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	A	0.54	0/3244	1.01	13/4373 (0.3%)
1	B	0.62	3/3244 (0.1%)	1.10	22/4373 (0.5%)
1	C	0.54	0/3244	0.98	16/4373 (0.4%)
1	D	0.53	0/3244	1.01	15/4373 (0.3%)
1	E	0.62	0/3244	1.08	17/4373 (0.4%)
1	F	0.55	0/3244	0.94	13/4373 (0.3%)
All	All	0.57	3/19464 (0.0%)	1.02	96/26238 (0.4%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	170	ASP	N-CA	7.62	1.56	1.46
1	B	169	ASP	CA-C	5.33	1.59	1.52
1	B	167	GLN	C-N	-5.21	1.26	1.33

All (96) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	130	ARG	CG-CD-NE	14.03	142.86	112.00
1	E	442	ILE	N-CA-C	11.85	122.00	111.81
1	B	109	LYS	CB-CG-CD	-10.66	86.78	111.30
1	B	442	ILE	N-CA-C	10.54	120.87	111.81
1	B	274	ARG	N-CA-C	-9.16	101.25	111.14
1	A	442	ILE	N-CA-C	8.76	119.34	111.81
1	E	274	ARG	N-CA-C	-8.73	101.71	111.14
1	D	442	ILE	N-CA-C	8.69	119.28	111.81
1	C	34	ARG	N-CA-C	-8.15	101.98	111.03
1	B	165	GLU	CB-CA-C	-8.09	94.33	110.42
1	F	34	ARG	N-CA-C	-8.05	102.20	110.97
1	C	442	ILE	N-CA-C	7.96	118.99	111.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	309	ALA	N-CA-C	-7.52	103.05	112.90
1	E	47	GLU	N-CA-C	-7.48	104.76	114.04
1	F	442	ILE	N-CA-C	7.05	118.91	112.29
1	E	165	GLU	CB-CA-C	-6.94	101.43	111.70
1	D	235	ASN	CA-C-N	6.93	126.63	119.56
1	D	235	ASN	C-N-CA	6.93	126.63	119.56
1	A	300	ASP	N-CA-C	6.86	120.54	111.75
1	A	376	ILE	N-CA-C	-6.84	103.67	110.30
1	B	96	VAL	N-CA-C	6.82	118.40	110.62
1	E	168	LEU	CA-C-N	-6.72	108.71	121.54
1	E	168	LEU	C-N-CA	-6.72	108.71	121.54
1	E	273	SER	N-CA-C	-6.65	105.30	113.41
1	E	171	LYS	N-CA-C	6.63	116.61	107.73
1	B	167	GLN	O-C-N	6.56	130.99	123.31
1	C	235	ASN	CA-C-N	6.47	127.92	119.84
1	C	235	ASN	C-N-CA	6.47	127.92	119.84
1	A	345	THR	N-CA-C	6.39	121.69	113.18
1	E	169	ASP	CA-CB-CG	-6.33	106.27	112.60
1	F	339	ASP	N-CA-C	-6.33	104.30	111.07
1	C	278	GLN	N-CA-C	-6.31	104.33	111.14
1	D	376	ILE	N-CA-C	-6.27	104.40	110.42
1	B	273	SER	N-CA-C	-6.27	105.77	113.41
1	A	274	ARG	N-CA-C	-6.26	103.76	111.40
1	B	355	TYR	N-CA-C	-6.17	104.47	111.14
1	F	309	ALA	N-CA-C	-6.12	104.89	112.90
1	B	165	GLU	O-C-N	6.11	130.72	122.59
1	B	168	LEU	O-C-N	-6.06	116.22	123.31
1	E	167	GLN	O-C-N	6.00	130.29	123.27
1	D	240	LYS	N-CA-C	-5.96	104.78	111.28
1	E	168	LEU	CB-CA-C	-5.94	97.63	109.33
1	D	300	ASP	N-CA-C	5.93	118.53	111.71
1	A	74	ALA	N-CA-C	-5.92	104.78	112.23
1	B	47	GLU	N-CA-C	-5.85	106.79	114.04
1	F	396	HIS	N-CA-C	5.79	117.28	110.97
1	D	84	THR	N-CA-C	5.77	118.52	111.82
1	C	318	LEU	N-CA-C	-5.71	102.45	110.50
1	A	428	HIS	N-CA-C	5.69	119.40	112.23
1	C	326	LEU	CA-C-N	5.68	125.04	118.85
1	C	326	LEU	C-N-CA	5.68	125.04	118.85
1	E	167	GLN	CB-CA-C	-5.68	101.05	110.14
1	C	294	HIS	N-CA-C	-5.67	105.77	114.16
1	B	101	ARG	N-CA-C	-5.66	104.80	110.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	278	GLN	N-CA-C	-5.66	105.03	111.14
1	F	318	LEU	N-CA-C	-5.58	102.63	110.50
1	F	95	GLU	N-CA-C	5.55	118.00	109.07
1	A	260	LYS	N-CA-C	-5.46	106.65	113.15
1	B	89	VAL	N-CA-C	-5.45	107.16	112.29
1	C	95	GLU	N-CA-C	5.45	118.37	109.59
1	B	169	ASP	O-C-N	-5.43	115.36	122.59
1	A	240	LYS	N-CA-C	-5.43	105.36	111.28
1	B	166	GLY	N-CA-C	5.42	126.01	113.18
1	C	360	ALA	N-CA-C	-5.40	105.95	112.54
1	D	348	ASN	N-CA-C	5.37	117.58	111.02
1	F	326	LEU	CA-C-N	5.35	124.98	119.05
1	F	326	LEU	C-N-CA	5.35	124.98	119.05
1	F	425	VAL	N-CA-C	5.34	115.97	110.36
1	D	274	ARG	N-CA-C	-5.34	104.88	111.40
1	E	89	VAL	N-CA-C	-5.34	107.27	112.29
1	A	246	ALA	N-CA-C	-5.34	104.89	111.40
1	C	62	GLY	N-CA-C	5.33	122.05	114.85
1	C	236	PRO	N-CA-C	5.20	123.19	112.47
1	E	91	TYR	CB-CA-C	-5.20	109.59	117.07
1	B	344	LEU	N-CA-C	5.19	116.94	111.28
1	B	372	GLY	N-CA-C	-5.19	106.47	112.50
1	D	74	ALA	N-CA-C	-5.19	105.69	112.23
1	B	169	ASP	N-CA-C	5.16	121.78	110.80
1	B	105	ASP	N-CA-C	-5.13	104.95	111.11
1	C	166	GLY	N-CA-C	5.13	117.12	110.45
1	D	70	LEU	N-CA-C	-5.13	105.38	111.69
1	A	70	LEU	N-CA-C	-5.13	105.76	112.23
1	A	166	GLY	N-CA-C	5.13	117.12	110.45
1	E	163	LEU	N-CA-C	5.13	116.56	110.97
1	A	318	LEU	N-CA-C	-5.11	103.65	110.55
1	F	276	GLY	N-CA-C	-5.11	107.98	114.16
1	B	160	ARG	NE-CZ-NH2	5.10	123.79	119.20
1	E	160	ARG	NE-CZ-NH2	5.10	123.79	119.20
1	D	57	GLY	CA-C-N	5.08	126.19	119.84
1	D	57	GLY	C-N-CA	5.08	126.19	119.84
1	F	166	GLY	N-CA-C	5.08	117.05	110.45
1	D	166	GLY	N-CA-C	5.07	117.04	110.45
1	E	164	ARG	NE-CZ-NH2	5.05	123.74	119.20
1	D	318	LEU	N-CA-C	-5.04	103.88	110.53
1	C	376	ILE	N-CA-C	-5.03	104.81	111.09
1	B	164	ARG	NE-CZ-NH2	5.00	123.70	119.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3205	0	3265	213	0
1	B	3205	0	3265	255	0
1	C	3205	0	3265	237	0
1	D	3205	0	3265	202	0
1	E	3205	0	3265	248	0
1	F	3205	0	3265	236	0
2	A	31	0	12	5	0
2	B	31	0	12	4	0
2	D	31	0	12	5	0
2	E	31	0	12	5	0
3	B	1	0	0	0	0
3	E	1	0	0	0	0
4	C	5	0	0	0	0
4	F	5	0	0	0	0
All	All	19366	0	19638	1325	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 40.

All (1325) 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:E:168:LEU:O	1:E:169:ASP:CB	1.77	1.20
1:B:169:ASP:O	1:B:170:ASP:CG	1.84	1.20
1:D:105:ASP:HB3	1:E:296:MET:HE1	1.32	1.12
1:C:312:ILE:HG13	1:C:313:ALA:H	1.20	1.07
1:E:168:LEU:O	1:E:169:ASP:HB2	1.28	1.06
1:B:169:ASP:O	1:B:170:ASP:OD1	1.74	1.06
1:A:17:ILE:HD13	1:A:66:ILE:HG12	1.42	1.02
1:D:27:VAL:CG1	1:D:70:LEU:HG	1.94	0.98
1:F:312:ILE:HG13	1:F:313:ALA:H	1.26	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:282:LEU:HD11	1:C:319:ILE:HD11	1.47	0.97
1:E:41:ASN:C	1:E:41:ASN:HD22	1.73	0.95
1:E:264:ARG:HB2	1:E:274:ARG:HH12	1.29	0.95
1:F:371:SER:HB2	1:F:422:ALA:HB2	1.46	0.95
1:D:41:ASN:HD21	1:D:44:LEU:H	1.12	0.94
1:D:68:ARG:O	1:D:72:LYS:HD3	1.67	0.94
1:E:168:LEU:O	1:E:169:ASP:CG	2.09	0.94
1:B:41:ASN:C	1:B:41:ASN:HD22	1.71	0.93
1:A:105:ASP:HB3	1:B:296:MET:HE1	1.50	0.93
1:B:136:ILE:HD11	1:B:159:PHE:HD2	1.32	0.92
1:B:261:ILE:HG21	1:B:278:GLN:HG3	1.49	0.92
1:F:344:LEU:HD13	1:F:395:LEU:HD13	1.50	0.92
1:A:41:ASN:HD21	1:A:44:LEU:H	1.15	0.91
1:F:37:ARG:HD2	1:F:48:VAL:HG23	1.50	0.91
1:B:264:ARG:HB2	1:B:274:ARG:HH12	1.36	0.90
1:D:17:ILE:HD13	1:D:66:ILE:HG12	1.55	0.89
1:C:108:VAL:HG21	1:C:294:HIS:CE1	2.08	0.89
1:A:110:MET:O	1:A:114:GLN:HB2	1.73	0.88
1:A:327:PRO:HA	1:F:397:THR:HG22	1.54	0.88
1:B:131:ILE:HG23	1:B:171:LYS:HE3	1.54	0.87
1:A:41:ASN:C	1:A:41:ASN:HD22	1.79	0.87
1:A:27:VAL:HG13	1:A:70:LEU:HG	1.57	0.87
1:A:3:GLU:HG2	1:A:35:TRP:HB2	1.57	0.86
1:D:27:VAL:HG13	1:D:70:LEU:HG	1.58	0.86
1:A:27:VAL:CG1	1:A:70:LEU:HG	2.06	0.85
1:C:344:LEU:HD13	1:C:395:LEU:HD13	1.57	0.85
1:B:344:LEU:CD1	1:B:351:ILE:HD11	2.07	0.85
1:D:3:GLU:HG2	1:D:35:TRP:HB2	1.59	0.85
1:B:261:ILE:HG22	1:B:278:GLN:HE21	1.42	0.84
1:B:84:THR:HA	1:B:87:THR:HG23	1.59	0.84
1:B:18:ILE:H	2:B:905:ATP:HN61	1.26	0.84
1:D:18:ILE:H	2:D:910:ATP:HN61	1.22	0.84
1:D:41:ASN:ND2	1:D:44:LEU:H	1.75	0.83
1:F:312:ILE:HG13	1:F:313:ALA:N	1.93	0.83
1:D:41:ASN:C	1:D:41:ASN:HD22	1.83	0.83
1:E:41:ASN:HD21	1:E:44:LEU:H	1.26	0.83
1:A:18:ILE:H	2:A:900:ATP:HN61	1.23	0.83
1:E:130:ARG:HH21	1:E:225:LEU:HD22	1.43	0.83
1:F:22:ASN:O	1:F:25:ARG:HG2	1.78	0.83
1:F:298:LYS:O	1:F:299:THR:HB	1.77	0.83
1:A:68:ARG:O	1:A:72:LYS:HD3	1.79	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:351:ILE:HD12	1:A:399:LEU:HD12	1.61	0.82
1:C:62:GLY:O	1:C:66:ILE:HG13	1.80	0.81
1:A:239:LEU:HA	1:A:242:ASP:HB2	1.62	0.81
1:C:397:THR:HG22	1:D:327:PRO:HA	1.62	0.81
1:B:101:ARG:HG3	1:B:101:ARG:HH11	1.44	0.81
1:C:240:LYS:O	1:C:244:ILE:HG12	1.81	0.81
1:D:108:VAL:HG21	1:D:294:HIS:CE1	2.16	0.81
1:E:122:ARG:NH2	1:E:126:LEU:HD21	1.95	0.81
1:A:239:LEU:HD12	1:A:240:LYS:N	1.96	0.80
1:C:298:LYS:O	1:C:299:THR:HB	1.81	0.80
1:C:312:ILE:HG13	1:C:313:ALA:N	1.93	0.80
1:C:371:SER:HB2	1:C:422:ALA:HB2	1.62	0.80
1:A:235:ASN:HB2	1:A:238:GLU:HB2	1.62	0.80
1:C:337:THR:O	1:C:341:GLU:HG3	1.82	0.79
1:E:84:THR:HA	1:E:87:THR:HG23	1.65	0.79
1:B:285:VAL:O	1:B:325:ARG:HD3	1.83	0.79
1:A:361:THR:HG21	1:B:36:ARG:HG2	1.63	0.79
1:B:17:ILE:HD13	1:B:66:ILE:HG12	1.66	0.78
1:A:282:LEU:O	1:A:286:GLU:HG3	1.83	0.78
1:B:172:GLU:HB3	1:B:215:LYS:HA	1.66	0.78
1:E:369:THR:HG22	1:E:371:SER:H	1.49	0.78
1:A:5:THR:OG1	1:A:8:GLU:HG3	1.84	0.78
1:E:344:LEU:CD1	1:E:351:ILE:HD11	2.14	0.78
1:D:325:ARG:HE	1:D:325:ARG:HA	1.49	0.78
1:C:257:GLU:HB2	1:C:260:LYS:HG2	1.64	0.77
1:D:298:LYS:HE3	1:D:300:ASP:OD1	1.84	0.77
1:F:416:GLN:C	1:F:417:ASN:HD22	1.92	0.77
1:B:261:ILE:CG2	1:B:278:GLN:HG3	2.15	0.77
1:E:175:ILE:HD13	1:E:224:LEU:HD11	1.66	0.77
1:A:41:ASN:ND2	1:A:44:LEU:H	1.82	0.77
1:C:37:ARG:HD2	1:C:48:VAL:HG23	1.65	0.77
1:F:235:ASN:HB3	1:F:238:GLU:HB2	1.68	0.76
1:A:108:VAL:HG21	1:A:294:HIS:CE1	2.20	0.76
1:A:288:CYS:O	1:A:298:LYS:HD2	1.85	0.76
1:C:5:THR:HB	1:C:6:PRO:HD2	1.68	0.76
1:D:239:LEU:HA	1:D:242:ASP:HB2	1.68	0.76
1:F:257:GLU:HB2	1:F:260:LYS:HG2	1.67	0.76
1:B:216:LEU:HD13	1:B:217:LYS:HG2	1.68	0.76
1:D:5:THR:OG1	1:D:8:GLU:HG3	1.86	0.76
1:D:239:LEU:HD12	1:D:240:LYS:N	2.00	0.76
1:D:325:ARG:HE	1:D:325:ARG:CA	1.99	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:88:GLU:OE2	1:D:93:GLY:HA3	1.85	0.75
1:B:168:LEU:O	1:B:169:ASP:CB	2.31	0.75
1:F:235:ASN:CB	1:F:238:GLU:HB2	2.16	0.75
1:F:303:LEU:HG	1:F:305:ILE:HD11	1.69	0.75
1:A:64:THR:HB	2:A:900:ATP:O2A	1.85	0.75
1:C:408:TYR:CE1	1:D:25:ARG:HG2	2.21	0.75
1:E:173:ILE:N	1:E:173:ILE:HD12	2.02	0.75
1:F:108:VAL:HG21	1:F:294:HIS:CE1	2.21	0.75
1:D:394:ARG:HH11	1:D:394:ARG:HG2	1.52	0.74
1:C:359:MET:HE2	1:C:406:ILE:HG22	1.69	0.74
1:B:41:ASN:HD21	1:B:44:LEU:H	1.32	0.74
1:C:383:VAL:HB	1:C:394:ARG:HD3	1.68	0.74
1:C:389:ASN:C	1:C:389:ASN:HD22	1.95	0.74
1:E:101:ARG:HD3	1:E:292:THR:HG22	1.70	0.74
1:A:62:GLY:O	1:A:66:ILE:HG13	1.86	0.74
1:A:100:ILE:HB	1:A:290:VAL:HG21	1.70	0.74
1:E:429:LEU:O	1:E:433:VAL:HG23	1.88	0.74
1:F:101:ARG:HH11	1:F:101:ARG:HG2	1.53	0.74
1:F:289:THR:O	1:F:290:VAL:HB	1.86	0.73
1:F:389:ASN:C	1:F:389:ASN:HD22	1.94	0.73
1:C:282:LEU:HB2	1:C:283:PRO:HD3	1.70	0.73
1:B:427:LYS:HD2	1:B:427:LYS:O	1.89	0.73
1:D:361:THR:HG21	1:E:36:ARG:HG2	1.70	0.73
1:A:387:THR:HG22	1:A:440:ARG:CZ	2.18	0.73
1:E:18:ILE:H	2:E:915:ATP:HN61	1.36	0.73
1:F:383:VAL:HB	1:F:394:ARG:HD3	1.69	0.73
1:C:84:THR:O	1:C:87:THR:HG22	1.89	0.73
1:F:42:GLU:H	1:F:42:GLU:CD	1.96	0.73
1:A:298:LYS:HE3	1:A:300:ASP:OD1	1.88	0.72
1:D:100:ILE:HB	1:D:290:VAL:HG21	1.71	0.72
1:D:18:ILE:H	2:D:910:ATP:N6	1.88	0.72
1:A:41:ASN:C	1:A:41:ASN:ND2	2.47	0.72
1:B:168:LEU:O	1:B:169:ASP:CG	2.33	0.72
1:B:151:GLU:HB3	1:B:152:PRO:HD3	1.71	0.72
1:E:216:LEU:HD13	1:E:217:LYS:HG2	1.70	0.72
1:C:108:VAL:HG13	1:C:239:LEU:HD21	1.71	0.72
1:E:119:ASN:C	1:E:119:ASN:HD22	1.98	0.72
1:D:288:CYS:O	1:D:298:LYS:HD2	1.90	0.72
1:C:288:CYS:O	1:C:298:LYS:O	2.07	0.72
1:D:282:LEU:O	1:D:286:GLU:HG3	1.90	0.71
1:F:240:LYS:O	1:F:244:ILE:HG12	1.89	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:GLY:O	1:A:24:LYS:HE3	1.90	0.71
1:B:427:LYS:HG3	1:B:428:HIS:CD2	2.25	0.71
1:E:101:ARG:HG3	1:E:101:ARG:HH11	1.56	0.71
1:E:261:ILE:HG22	1:E:261:ILE:O	1.89	0.71
1:C:339:ASP:O	1:C:343:ILE:HG13	1.91	0.71
1:C:393:ARG:HG3	1:C:393:ARG:HH11	1.54	0.71
1:B:171:LYS:HZ3	1:B:173:ILE:HD11	1.56	0.71
1:A:88:GLU:HB2	1:A:91:TYR:HB2	1.71	0.71
1:C:101:ARG:HG2	1:C:101:ARG:HH11	1.55	0.71
1:A:25:ARG:HG2	1:F:408:TYR:CE1	2.26	0.71
1:E:239:LEU:HD12	1:E:240:LYS:N	2.06	0.71
1:A:394:ARG:HG2	1:A:394:ARG:HH11	1.55	0.70
1:B:152:PRO:O	1:B:156:ARG:HB3	1.91	0.70
1:B:350:SER:O	1:B:354:GLN:HG3	1.91	0.70
1:D:100:ILE:HD11	1:D:284:LEU:HD22	1.73	0.70
1:F:84:THR:O	1:F:87:THR:HG22	1.91	0.70
1:C:22:ASN:O	1:C:25:ARG:HG2	1.90	0.70
1:E:312:ILE:HG13	1:E:313:ALA:H	1.55	0.70
1:F:62:GLY:O	1:F:66:ILE:HG13	1.90	0.70
1:B:222:MET:O	1:B:226:ILE:HG13	1.92	0.70
1:D:27:VAL:HG11	1:D:70:LEU:HG	1.72	0.70
1:A:88:GLU:OE2	1:A:93:GLY:HA3	1.91	0.70
1:C:402:LEU:HD11	1:C:425:VAL:HG22	1.73	0.70
1:F:359:MET:HE1	1:F:407:SER:HA	1.72	0.70
1:C:72:LYS:HE3	1:C:72:LYS:O	1.91	0.70
1:E:362:GLU:HG3	1:E:410:ALA:HB1	1.72	0.70
1:B:173:ILE:HD12	1:B:173:ILE:H	1.56	0.70
1:F:402:LEU:HD11	1:F:425:VAL:HG22	1.74	0.70
1:B:264:ARG:CB	1:B:274:ARG:HH12	2.04	0.70
1:D:387:THR:HG22	1:D:440:ARG:CZ	2.22	0.70
1:D:19:GLY:O	1:D:24:LYS:HE3	1.90	0.70
1:E:79:ILE:HD11	1:E:102:ASP:HB3	1.73	0.69
1:B:172:GLU:CB	1:B:215:LYS:HA	2.22	0.69
1:C:85:LYS:HE2	1:C:94:LYS:HE2	1.74	0.69
1:A:107:ALA:O	1:A:111:VAL:HG23	1.91	0.69
1:B:239:LEU:HD12	1:B:240:LYS:N	2.07	0.69
1:E:427:LYS:HD2	1:E:427:LYS:O	1.91	0.69
1:F:339:ASP:O	1:F:343:ILE:HG13	1.93	0.69
1:A:18:ILE:H	2:A:900:ATP:N6	1.91	0.69
1:A:359:MET:HE1	1:B:36:ARG:NH1	2.08	0.69
1:B:42:GLU:HA	1:B:45:ARG:NH1	2.08	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:303:LEU:HG	1:C:305:ILE:HD11	1.73	0.69
1:D:100:ILE:HG13	1:D:290:VAL:HG11	1.75	0.69
1:F:393:ARG:HG3	1:F:393:ARG:HH11	1.55	0.69
1:B:383:VAL:HG12	1:B:394:ARG:NH1	2.07	0.69
1:E:271:ASP:C	1:E:273:SER:H	2.00	0.69
1:B:41:ASN:C	1:B:41:ASN:ND2	2.45	0.69
1:C:344:LEU:CD1	1:C:395:LEU:HD13	2.23	0.69
1:D:88:GLU:HB2	1:D:91:TYR:HB2	1.74	0.69
1:E:18:ILE:HD13	1:E:347:PRO:HG3	1.73	0.69
1:E:261:ILE:HG21	1:E:278:GLN:HG3	1.74	0.69
1:C:42:GLU:CD	1:C:42:GLU:H	1.99	0.68
1:C:289:THR:O	1:C:290:VAL:HB	1.91	0.68
1:F:359:MET:HG3	1:F:366:ILE:HD11	1.74	0.68
1:B:79:ILE:HD11	1:B:102:ASP:HB3	1.76	0.68
1:A:30:ALA:HB2	1:A:53:ILE:HD11	1.74	0.68
1:E:103:LEU:HD13	1:E:247:VAL:HG13	1.76	0.68
1:B:344:LEU:HD12	1:B:351:ILE:HD11	1.76	0.68
1:E:5:THR:OG1	1:E:8:GLU:HG3	1.93	0.68
1:B:130:ARG:NH1	1:B:225:LEU:HD22	2.09	0.68
1:C:240:LYS:HD2	1:C:241:GLN:HG3	1.76	0.68
1:A:327:PRO:HA	1:F:397:THR:CG2	2.24	0.67
1:F:103:LEU:HD13	1:F:247:VAL:HG13	1.77	0.67
1:A:100:ILE:HD11	1:A:284:LEU:HD22	1.76	0.67
1:D:105:ASP:HB3	1:E:296:MET:CE	2.17	0.67
1:F:235:ASN:HB3	1:F:238:GLU:C	2.20	0.67
1:F:384:ASN:HD22	1:F:394:ARG:HH11	1.43	0.67
1:B:429:LEU:O	1:B:433:VAL:HG23	1.95	0.67
1:E:168:LEU:HB3	1:E:218:ILE:HB	1.76	0.67
1:F:335:LEU:HD22	1:F:339:ASP:HB3	1.77	0.67
1:A:100:ILE:HG13	1:A:290:VAL:HG11	1.76	0.67
1:D:375:ARG:HB3	1:D:425:VAL:HG11	1.77	0.67
1:E:369:THR:HG22	1:E:371:SER:N	2.10	0.67
1:F:359:MET:CE	1:F:406:ILE:HG22	2.24	0.67
1:B:173:ILE:HD12	1:B:173:ILE:N	2.09	0.67
1:D:64:THR:HB	2:D:910:ATP:O2A	1.95	0.67
1:C:55:MET:HE1	1:C:63:LYS:HA	1.75	0.67
1:D:108:VAL:HG21	1:D:294:HIS:ND1	2.10	0.67
1:F:337:THR:O	1:F:341:GLU:HG3	1.94	0.67
1:C:389:ASN:C	1:C:389:ASN:ND2	2.53	0.66
1:D:41:ASN:ND2	1:D:41:ASN:C	2.50	0.66
1:D:107:ALA:O	1:D:111:VAL:HG23	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:359:MET:HG3	1:F:366:ILE:CD1	2.24	0.66
1:B:271:ASP:C	1:B:273:SER:H	2.03	0.66
1:A:325:ARG:CA	1:A:325:ARG:HE	2.08	0.66
1:D:35:TRP:O	1:D:39:GLN:HG2	1.95	0.66
1:A:427:LYS:HD2	1:A:427:LYS:O	1.95	0.66
1:D:110:MET:O	1:D:114:GLN:HB2	1.95	0.66
1:F:282:LEU:HD11	1:F:319:ILE:HD11	1.76	0.66
1:B:132:LEU:HB2	1:B:156:ARG:NH1	2.11	0.66
1:E:383:VAL:HG12	1:E:394:ARG:NH1	2.09	0.66
1:B:27:VAL:HG22	1:B:70:LEU:HG	1.78	0.66
1:A:39:GLN:HB2	1:F:361:THR:OG1	1.96	0.66
1:E:41:ASN:C	1:E:41:ASN:ND2	2.46	0.66
1:D:349:ALA:HB3	1:E:47:GLU:HG3	1.76	0.65
1:E:427:LYS:HG3	1:E:428:HIS:CD2	2.31	0.65
1:B:128:GLU:OE2	1:B:164:ARG:HD3	1.96	0.65
1:B:312:ILE:HG13	1:B:313:ALA:H	1.60	0.65
1:C:106:ALA:HA	1:C:109:LYS:HE3	1.77	0.65
1:C:111:VAL:CG2	1:C:239:LEU:HD11	2.25	0.65
1:B:261:ILE:HG22	1:B:261:ILE:O	1.95	0.65
1:F:274:ARG:HA	1:F:277:VAL:HG23	1.78	0.65
1:B:136:ILE:CD1	1:B:156:ARG:HA	2.26	0.65
1:E:151:GLU:HB3	1:E:152:PRO:HD3	1.78	0.65
1:F:289:THR:HG22	1:F:290:VAL:H	1.62	0.65
1:F:85:LYS:HE2	1:F:94:LYS:HE2	1.79	0.65
1:F:389:ASN:C	1:F:389:ASN:ND2	2.52	0.65
1:E:285:VAL:O	1:E:325:ARG:HD3	1.97	0.65
1:F:374:LYS:O	1:F:378:GLU:HG3	1.96	0.65
1:B:85:LYS:HG2	1:B:85:LYS:O	1.96	0.64
1:B:131:ILE:HD12	1:B:222:MET:SD	2.37	0.64
1:E:350:SER:O	1:E:354:GLN:HG3	1.97	0.64
1:C:359:MET:HE1	1:C:407:SER:HA	1.79	0.64
1:F:240:LYS:HD2	1:F:241:GLN:HG3	1.79	0.64
1:A:9:ILE:HG23	1:A:73:LEU:HD21	1.79	0.64
1:B:151:GLU:C	1:B:153:SER:H	2.04	0.64
1:C:235:ASN:HB3	1:C:238:GLU:OE1	1.98	0.64
1:A:298:LYS:HE3	1:A:300:ASP:CG	2.22	0.64
1:B:170:ASP:O	1:B:171:LYS:HB3	1.97	0.64
1:C:55:MET:CE	1:C:63:LYS:HA	2.27	0.64
1:C:274:ARG:HA	1:C:277:VAL:HG23	1.78	0.64
1:D:41:ASN:HD21	1:D:44:LEU:HD12	1.63	0.64
1:E:17:ILE:HD13	1:E:66:ILE:HG12	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:54:LEU:HD23	1:D:329:ARG:HD2	1.80	0.64
1:D:298:LYS:HE3	1:D:300:ASP:CG	2.22	0.64
1:C:37:ARG:O	1:C:39:GLN:N	2.30	0.64
1:E:393:ARG:HH11	1:E:393:ARG:HG2	1.62	0.64
1:D:41:ASN:ND2	1:D:44:LEU:HD12	2.12	0.63
1:D:384:ASN:ND2	1:D:394:ARG:NH1	2.46	0.63
1:C:359:MET:CE	1:C:406:ILE:HG22	2.28	0.63
1:F:106:ALA:HA	1:F:109:LYS:HE3	1.80	0.63
1:D:32:ARG:O	1:D:36:ARG:HG3	1.98	0.63
1:B:18:ILE:HD13	1:B:347:PRO:HG3	1.80	0.63
1:E:312:ILE:HG13	1:E:313:ALA:N	2.13	0.63
1:F:390:ILE:HG21	1:F:393:ARG:HB2	1.81	0.63
1:B:369:THR:HG22	1:B:371:SER:H	1.64	0.63
1:E:9:ILE:HD13	1:E:31:LEU:HD12	1.80	0.63
1:E:173:ILE:HD12	1:E:173:ILE:H	1.62	0.63
1:B:336:THR:O	1:B:339:ASP:HB2	1.98	0.63
1:C:25:ARG:O	1:C:29:ILE:HG13	1.98	0.63
1:A:108:VAL:HG21	1:A:294:HIS:ND1	2.13	0.63
1:F:72:LYS:HE3	1:F:72:LYS:O	1.98	0.63
1:A:38:MET:HE3	1:A:38:MET:HA	1.81	0.62
1:D:427:LYS:HD2	1:D:427:LYS:C	2.24	0.62
1:F:100:ILE:HG21	1:F:290:VAL:HG11	1.80	0.62
1:A:3:GLU:HG2	1:A:35:TRP:CB	2.30	0.62
1:B:5:THR:HG23	1:B:8:GLU:CD	2.24	0.62
1:C:23:ALA:HA	1:C:330:VAL:HG21	1.80	0.62
1:C:74:ALA:O	1:C:75:ASN:C	2.43	0.62
1:B:42:GLU:HA	1:B:45:ARG:HH12	1.64	0.62
1:B:174:GLU:HA	1:B:213:ALA:N	2.14	0.62
1:B:362:GLU:HG3	1:B:410:ALA:HB1	1.80	0.62
1:C:335:LEU:HD22	1:C:339:ASP:HB3	1.80	0.62
1:E:27:VAL:HG22	1:E:70:LEU:HG	1.81	0.62
1:A:397:THR:HG21	1:A:443:LEU:C	2.25	0.62
1:E:261:ILE:CG2	1:E:278:GLN:HG3	2.30	0.62
1:F:65:GLU:OE1	1:F:69:ARG:NH2	2.29	0.62
1:D:384:ASN:HA	1:D:394:ARG:HH12	1.64	0.62
1:B:18:ILE:N	2:B:905:ATP:HN61	1.95	0.62
1:D:105:ASP:O	1:E:296:MET:HE2	1.99	0.62
1:B:106:ALA:O	1:B:110:MET:HG3	2.00	0.62
1:D:300:ASP:C	1:D:301:HIS:HD2	2.07	0.61
1:F:398:VAL:HG22	1:F:443:LEU:HD22	1.80	0.61
1:A:279:ARG:HG2	1:A:279:ARG:HH11	1.66	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:369:THR:HG22	1:B:371:SER:N	2.15	0.61
1:A:94:LYS:HE3	1:B:95:GLU:HG2	1.82	0.61
1:B:235:ASN:HB2	1:B:238:GLU:HB2	1.81	0.61
1:E:169:ASP:O	1:E:170:ASP:CG	2.43	0.61
1:F:240:LYS:HD2	1:F:241:GLN:N	2.14	0.61
1:E:270:PRO:HA	1:E:273:SER:HB3	1.82	0.61
1:F:282:LEU:O	1:F:286:GLU:HG2	2.01	0.61
1:F:344:LEU:CD1	1:F:395:LEU:HD13	2.27	0.61
1:B:393:ARG:HG2	1:B:393:ARG:HH11	1.65	0.61
1:C:107:ALA:O	1:C:110:MET:HB2	2.00	0.61
1:C:83:ALA:HB1	1:C:261:ILE:HD13	1.82	0.61
1:C:264:ARG:HB3	1:C:271:ASP:OD2	2.00	0.61
1:E:264:ARG:CB	1:E:274:ARG:HH12	2.09	0.61
1:A:387:THR:HA	1:A:440:ARG:NH2	2.16	0.61
1:D:358:LEU:O	1:D:361:THR:HB	2.01	0.61
1:A:270:PRO:O	1:A:273:SER:HB3	2.01	0.61
1:C:53:ILE:HG23	1:C:328:ILE:HG22	1.81	0.61
1:A:3:GLU:HG2	1:A:35:TRP:CD1	2.36	0.60
1:E:217:LYS:HE3	1:E:220:ASP:OD2	2.01	0.60
1:C:282:LEU:HD11	1:C:319:ILE:CD1	2.28	0.60
1:E:216:LEU:CD1	1:E:217:LYS:H	2.12	0.60
1:A:88:GLU:CB	1:A:91:TYR:HB2	2.32	0.60
1:B:55:MET:HE3	1:B:305:ILE:HG21	1.83	0.60
1:C:20:GLN:O	1:C:24:LYS:HG3	2.01	0.60
1:F:111:VAL:CG2	1:F:239:LEU:HD11	2.32	0.60
1:A:366:ILE:HD13	1:A:418:ILE:HB	1.82	0.60
1:E:222:MET:O	1:E:226:ILE:HG13	2.01	0.60
1:F:266:GLU:HG3	1:F:270:PRO:HG2	1.83	0.60
1:B:359:MET:HE2	1:B:359:MET:HA	1.83	0.60
1:C:441:PHE:HD1	1:D:56:ILE:HD13	1.66	0.60
1:D:40:LEU:CD1	1:D:48:VAL:HG21	2.32	0.60
1:B:27:VAL:CG2	1:B:70:LEU:HG	2.32	0.60
1:B:101:ARG:HG3	1:B:101:ARG:NH1	2.16	0.60
1:D:37:ARG:HG3	1:D:48:VAL:HG11	1.82	0.60
1:E:336:THR:O	1:E:339:ASP:HB2	2.01	0.60
1:B:103:LEU:HD13	1:B:247:VAL:HG13	1.82	0.60
1:C:289:THR:HG22	1:C:290:VAL:H	1.66	0.60
1:D:427:LYS:HD2	1:D:427:LYS:O	2.02	0.60
1:B:359:MET:HE1	1:C:36:ARG:NH1	2.16	0.60
1:D:33:ASN:HD22	1:D:36:ARG:HD2	1.66	0.60
1:D:238:GLU:O	1:D:241:GLN:N	2.34	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:63:LYS:HG2	1:E:332:LEU:HD22	1.84	0.60
1:E:122:ARG:HH22	1:E:126:LEU:HD21	1.65	0.60
1:B:161:LYS:HG3	1:B:165:GLU:HG3	1.84	0.60
1:C:266:GLU:HG3	1:C:270:PRO:HG2	1.82	0.60
1:C:397:THR:CG2	1:D:327:PRO:HA	2.32	0.60
1:C:416:GLN:C	1:C:417:ASN:HD22	2.09	0.60
1:E:59:THR:HG22	1:E:393:ARG:HE	1.67	0.60
1:A:56:ILE:HD13	1:F:441:PHE:HD1	1.66	0.59
1:D:38:MET:CE	1:D:45:ARG:HE	2.14	0.59
1:B:312:ILE:HG13	1:B:313:ALA:N	2.17	0.59
1:E:85:LYS:O	1:E:85:LYS:HG2	2.01	0.59
1:C:240:LYS:HD2	1:C:241:GLN:N	2.17	0.59
1:C:388:GLU:N	1:C:388:GLU:OE1	2.35	0.59
1:E:435:ASP:CG	1:E:438:LEU:HD23	2.27	0.59
1:D:359:MET:HE1	1:E:36:ARG:NH1	2.18	0.59
1:A:51:LYS:O	1:A:328:ILE:HD12	2.02	0.59
1:A:427:LYS:HD2	1:A:427:LYS:C	2.27	0.59
1:C:103:LEU:HD13	1:C:247:VAL:HG13	1.85	0.59
1:E:216:LEU:HD12	1:E:217:LYS:H	1.65	0.59
1:F:359:MET:HE2	1:F:406:ILE:HG22	1.84	0.59
1:A:413:LEU:HA	1:A:416:GLN:OE1	2.03	0.59
1:B:101:ARG:HD3	1:B:292:THR:HG22	1.83	0.59
1:E:131:ILE:HD12	1:E:222:MET:SD	2.42	0.59
1:E:271:ASP:C	1:E:273:SER:N	2.60	0.59
1:C:65:GLU:OE1	1:C:69:ARG:NH2	2.36	0.59
1:C:390:ILE:HG21	1:C:393:ARG:HB2	1.85	0.59
1:F:110:MET:O	1:F:114:GLN:HB2	2.02	0.59
1:F:304:PHE:C	1:F:305:ILE:HD12	2.27	0.59
1:B:413:LEU:O	1:B:416:GLN:HG3	2.02	0.59
1:D:361:THR:HG22	1:D:362:GLU:HG2	1.85	0.59
1:E:175:ILE:HG21	1:E:228:GLU:OE1	2.03	0.59
1:E:441:PHE:HD1	1:F:56:ILE:HD13	1.67	0.59
1:E:4:MET:HE2	1:E:9:ILE:HA	1.83	0.59
1:E:106:ALA:O	1:E:110:MET:HG3	2.02	0.59
1:C:111:VAL:HG21	1:C:239:LEU:HD11	1.83	0.59
1:B:132:LEU:HD11	1:B:160:ARG:HE	1.67	0.58
1:E:261:ILE:HG22	1:E:278:GLN:HE21	1.68	0.58
1:F:20:GLN:O	1:F:24:LYS:HG3	2.02	0.58
1:B:264:ARG:HB2	1:B:274:ARG:NH1	2.13	0.58
1:D:44:LEU:O	1:D:47:GLU:N	2.32	0.58
1:E:79:ILE:CD1	1:E:102:ASP:HB3	2.32	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:401:ARG:HA	1:E:404:GLU:HG3	1.86	0.58
1:C:372:GLY:O	1:C:376:ILE:HG13	2.03	0.58
1:B:383:VAL:CG1	1:B:394:ARG:NH1	2.66	0.58
1:D:88:GLU:CB	1:D:91:TYR:HB2	2.34	0.58
1:A:35:TRP:O	1:A:39:GLN:HG2	2.03	0.58
1:A:238:GLU:O	1:A:241:GLN:N	2.36	0.58
1:E:175:ILE:HG22	1:E:176:ASP:N	2.18	0.58
1:E:270:PRO:O	1:E:273:SER:HB3	2.03	0.58
1:F:55:MET:HE1	1:F:63:LYS:HA	1.86	0.58
1:F:352:THR:HG22	1:F:353:VAL:N	2.19	0.58
1:A:33:ASN:HD22	1:A:36:ARG:HD2	1.69	0.58
1:F:9:ILE:HD13	1:F:31:LEU:HD23	1.85	0.58
1:A:88:GLU:HB2	1:A:91:TYR:CB	2.34	0.58
1:A:384:ASN:ND2	1:A:390:ILE:H	2.02	0.58
1:E:131:ILE:O	1:E:134:VAL:HG12	2.04	0.58
1:E:152:PRO:O	1:E:156:ARG:HB3	2.04	0.58
1:E:27:VAL:CG2	1:E:70:LEU:HG	2.33	0.58
1:A:257:GLU:C	1:A:259:ASP:H	2.12	0.57
1:A:312:ILE:HG13	1:A:313:ALA:N	2.19	0.57
1:B:234:VAL:CG1	1:B:239:LEU:HD23	2.34	0.57
1:D:38:MET:HE3	1:D:38:MET:HA	1.85	0.57
1:E:262:CYS:SG	1:E:318:LEU:HD23	2.44	0.57
1:C:107:ALA:O	1:C:111:VAL:HG22	2.03	0.57
1:C:432:LEU:HG	1:C:442:ILE:HD11	1.86	0.57
1:B:270:PRO:HA	1:B:273:SER:HB3	1.86	0.57
1:C:266:GLU:HG2	1:C:271:ASP:HB2	1.85	0.57
1:E:42:GLU:HA	1:E:45:ARG:NH1	2.20	0.57
1:E:270:PRO:C	1:E:273:SER:HB3	2.30	0.57
1:E:41:ASN:ND2	1:E:44:LEU:H	1.99	0.57
1:F:282:LEU:HD12	1:F:322:LEU:HD13	1.87	0.57
1:E:101:ARG:NE	1:E:291:SER:O	2.35	0.57
1:E:122:ARG:HE	1:E:122:ARG:HA	1.70	0.57
1:F:402:LEU:HB2	1:F:429:LEU:HD21	1.86	0.57
1:F:240:LYS:C	1:F:242:ASP:H	2.11	0.57
1:F:345:THR:HG22	1:F:373:ILE:HD13	1.84	0.57
1:B:5:THR:OG1	1:B:8:GLU:HG3	2.05	0.57
1:E:174:GLU:HA	1:E:213:ALA:HA	1.86	0.57
1:F:345:THR:CG2	1:F:373:ILE:HD13	2.34	0.57
1:C:361:THR:OG1	1:D:39:GLN:HB2	2.05	0.57
1:E:289:THR:CG2	1:E:296:MET:HE3	2.35	0.57
1:E:342:ARG:HG2	1:E:346:GLU:OE1	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:5:THR:HB	1:F:6:PRO:HD2	1.85	0.57
1:A:105:ASP:O	1:B:296:MET:HE2	2.05	0.57
1:C:359:MET:HG3	1:C:366:ILE:CD1	2.35	0.57
1:B:289:THR:CG2	1:B:296:MET:HE3	2.35	0.57
1:F:83:ALA:HB1	1:F:261:ILE:HD13	1.87	0.57
1:B:9:ILE:HD13	1:B:31:LEU:HD12	1.87	0.56
1:B:152:PRO:O	1:B:153:SER:O	2.22	0.56
1:B:359:MET:HG3	1:B:366:ILE:HG13	1.87	0.56
1:B:435:ASP:CG	1:B:438:LEU:HD23	2.30	0.56
1:C:345:THR:CG2	1:C:373:ILE:HD13	2.35	0.56
1:F:372:GLY:O	1:F:376:ILE:HG13	2.05	0.56
1:D:270:PRO:O	1:D:273:SER:HB3	2.05	0.56
1:D:325:ARG:HA	1:D:325:ARG:NE	2.19	0.56
1:F:74:ALA:O	1:F:75:ASN:C	2.48	0.56
1:F:303:LEU:HG	1:F:305:ILE:CD1	2.34	0.56
1:B:354:GLN:OE1	1:C:48:VAL:HG12	2.06	0.56
1:C:72:LYS:HE3	1:C:72:LYS:C	2.30	0.56
1:C:85:LYS:CE	1:C:94:LYS:HE2	2.35	0.56
1:E:383:VAL:CG1	1:E:394:ARG:NH1	2.68	0.56
1:F:16:HIS:CD2	1:F:69:ARG:HD2	2.40	0.56
1:F:288:CYS:O	1:F:298:LYS:O	2.23	0.56
1:F:442:ILE:C	1:F:442:ILE:HD12	2.31	0.56
1:B:175:ILE:HD13	1:B:224:LEU:HD11	1.88	0.56
1:D:249:GLN:O	1:D:250:HIS:ND1	2.38	0.56
1:A:18:ILE:N	2:A:900:ATP:HN61	2.00	0.56
1:F:233:LEU:O	1:F:235:ASN:N	2.39	0.56
1:B:81:VAL:HG11	1:B:99:ILE:HG12	1.87	0.56
1:D:366:ILE:HD13	1:D:418:ILE:HB	1.87	0.56
1:D:387:THR:HA	1:D:440:ARG:NH2	2.20	0.56
1:A:249:GLN:O	1:A:250:HIS:ND1	2.39	0.56
1:A:375:ARG:HB3	1:A:425:VAL:HG11	1.87	0.56
1:F:350:SER:O	1:F:354:GLN:HG3	2.06	0.56
1:B:270:PRO:O	1:B:273:SER:HB3	2.06	0.56
1:C:75:ASN:HD22	1:C:114:GLN:HE22	1.54	0.56
1:D:384:ASN:HD22	1:D:394:ARG:HH12	1.53	0.56
1:B:95:GLU:OE1	1:B:95:GLU:HA	2.06	0.56
1:B:132:LEU:HD13	1:B:156:ARG:HG3	1.89	0.56
1:B:217:LYS:HG3	1:B:220:ASP:HB2	1.87	0.56
1:B:384:ASN:HD21	1:B:394:ARG:HB2	1.69	0.56
1:E:125:GLU:O	1:E:128:GLU:HG3	2.06	0.56
1:E:358:LEU:O	1:E:361:THR:HB	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:389:ASN:HD22	1:F:390:ILE:N	2.03	0.56
1:B:136:ILE:HD11	1:B:159:PHE:CD2	2.25	0.55
1:E:17:ILE:HD13	1:E:66:ILE:CG1	2.35	0.55
1:B:79:ILE:CD1	1:B:102:ASP:HB3	2.37	0.55
1:D:359:MET:HE2	1:D:359:MET:HA	1.88	0.55
1:E:105:ASP:HB3	1:F:296:MET:HE1	1.88	0.55
1:E:132:LEU:HB2	1:E:156:ARG:NH1	2.22	0.55
1:A:41:ASN:HD21	1:A:44:LEU:HD12	1.70	0.55
1:E:261:ILE:HG23	1:E:274:ARG:O	2.05	0.55
1:F:264:ARG:HB3	1:F:271:ASP:OD2	2.07	0.55
1:C:9:ILE:HD13	1:C:31:LEU:HD23	1.89	0.55
1:C:19:GLY:O	1:C:24:LYS:HE3	2.07	0.55
1:F:72:LYS:HE3	1:F:72:LYS:C	2.31	0.55
1:B:41:ASN:ND2	1:B:44:LEU:H	2.04	0.55
1:D:100:ILE:CB	1:D:290:VAL:HG21	2.37	0.55
1:E:107:ALA:O	1:E:111:VAL:HG23	2.06	0.55
1:F:417:ASN:HD22	1:F:417:ASN:N	2.01	0.55
1:B:229:GLU:O	1:B:233:LEU:HD23	2.07	0.55
1:B:349:ALA:CB	1:C:47:GLU:HG3	2.37	0.55
1:C:3:GLU:HG3	1:C:35:TRP:HB2	1.89	0.55
1:F:429:LEU:O	1:F:433:VAL:HG23	2.07	0.55
1:C:41:ASN:HD21	1:C:43:GLU:HB3	1.72	0.55
1:D:62:GLY:O	1:D:66:ILE:HG13	2.07	0.55
1:A:300:ASP:C	1:A:301:HIS:HD2	2.15	0.55
1:C:390:ILE:HG21	1:C:393:ARG:HG3	1.88	0.55
1:E:383:VAL:HB	1:E:394:ARG:HH11	1.72	0.55
1:F:41:ASN:HD21	1:F:43:GLU:HB3	1.72	0.55
1:A:41:ASN:ND2	1:A:44:LEU:HD12	2.21	0.54
1:A:278:GLN:OE1	1:A:319:ILE:HG12	2.07	0.54
1:B:383:VAL:HB	1:B:394:ARG:HH11	1.72	0.54
1:E:168:LEU:HB2	1:E:169:ASP:OD2	2.06	0.54
1:A:271:ASP:HA	1:A:274:ARG:HH11	1.70	0.54
1:B:346:GLU:O	1:B:347:PRO:C	2.49	0.54
1:F:100:ILE:CG2	1:F:290:VAL:HG11	2.37	0.54
1:F:384:ASN:ND2	1:F:394:ARG:HH11	2.03	0.54
1:D:38:MET:HE3	1:D:45:ARG:HE	1.70	0.54
1:A:34:ARG:HH12	1:A:250:HIS:C	2.16	0.54
1:A:82:GLU:O	1:A:84:THR:N	2.39	0.54
1:B:134:VAL:HG11	1:B:171:LYS:HD2	1.89	0.54
1:B:442:ILE:HG23	1:C:329:ARG:HB2	1.89	0.54
1:D:279:ARG:HH11	1:D:279:ARG:HG2	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:GLN:HB2	1:F:361:THR:HG1	1.73	0.54
1:C:111:VAL:O	1:C:114:GLN:HB3	2.08	0.54
1:C:436:GLU:O	1:C:439:SER:HB2	2.08	0.54
1:D:63:LYS:HG2	1:D:332:LEU:HD22	1.89	0.54
1:D:88:GLU:HB2	1:D:91:TYR:CB	2.38	0.54
1:E:383:VAL:HG12	1:E:394:ARG:HH12	1.72	0.54
1:A:100:ILE:CB	1:A:290:VAL:HG21	2.37	0.54
1:B:18:ILE:CD1	1:B:347:PRO:HG3	2.37	0.54
1:B:355:TYR:CZ	1:B:403:MET:HB2	2.42	0.54
1:C:5:THR:HB	1:C:6:PRO:CD	2.37	0.54
1:C:41:ASN:ND2	1:C:43:GLU:H	2.05	0.54
1:C:75:ASN:CG	1:C:75:ASN:O	2.51	0.54
1:D:355:TYR:OH	1:D:400:GLU:HG2	2.07	0.54
1:E:50:PRO:O	1:E:325:ARG:NH2	2.36	0.54
1:B:161:LYS:O	1:B:165:GLU:HG3	2.07	0.54
1:C:81:VAL:HG11	1:C:86:PHE:CE2	2.42	0.54
1:C:361:THR:HG21	1:D:36:ARG:CB	2.38	0.54
1:D:71:ALA:HA	1:D:252:ILE:HD12	1.90	0.54
1:E:128:GLU:OE2	1:E:164:ARG:HD3	2.06	0.54
1:A:4:MET:HB2	1:A:8:GLU:OE1	2.07	0.54
1:A:84:THR:O	1:A:86:PHE:N	2.41	0.54
1:D:63:LYS:HG2	1:D:332:LEU:HD13	1.90	0.54
1:D:384:ASN:ND2	1:D:390:ILE:H	2.06	0.54
1:F:81:VAL:HG11	1:F:86:PHE:CE2	2.43	0.54
1:F:266:GLU:HG2	1:F:271:ASP:HB2	1.89	0.54
1:F:305:ILE:HD12	1:F:305:ILE:N	2.22	0.54
1:D:270:PRO:HB2	1:D:274:ARG:HH12	1.73	0.54
1:F:284:LEU:HA	1:F:299:THR:HG21	1.90	0.54
1:F:402:LEU:HD23	1:F:403:MET:SD	2.48	0.54
1:A:384:ASN:ND2	1:A:394:ARG:NH1	2.55	0.54
1:E:261:ILE:O	1:E:261:ILE:CG2	2.56	0.54
1:F:48:VAL:HG23	1:F:48:VAL:O	2.08	0.54
1:B:151:GLU:C	1:B:153:SER:N	2.64	0.53
1:D:94:LYS:HE3	1:E:95:GLU:HG2	1.90	0.53
1:F:288:CYS:SG	1:F:289:THR:N	2.81	0.53
1:B:59:THR:HG22	1:B:393:ARG:HE	1.72	0.53
1:F:108:VAL:O	1:F:112:ARG:HB2	2.07	0.53
1:B:100:ILE:HB	1:B:290:VAL:HG21	1.89	0.53
1:E:239:LEU:HD12	1:E:239:LEU:C	2.34	0.53
1:F:101:ARG:HG2	1:F:101:ARG:NH1	2.20	0.53
1:F:361:THR:HG22	1:F:362:GLU:HG2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:LEU:O	1:A:47:GLU:N	2.37	0.53
1:C:240:LYS:CD	1:C:241:GLN:HG3	2.37	0.53
1:E:55:MET:HE3	1:E:305:ILE:HG21	1.90	0.53
1:E:244:ILE:O	1:E:245:ASP:C	2.50	0.53
1:E:335:LEU:HD13	1:E:343:ILE:HD11	1.90	0.53
1:E:379:ALA:O	1:E:383:VAL:HG23	2.08	0.53
1:A:40:LEU:CD1	1:A:48:VAL:HG21	2.39	0.53
1:A:249:GLN:C	1:A:250:HIS:ND1	2.66	0.53
1:C:398:VAL:HG22	1:C:443:LEU:HD22	1.90	0.53
1:D:248:GLU:HG3	1:D:297:VAL:HG13	1.91	0.53
1:B:107:ALA:O	1:B:111:VAL:HG23	2.09	0.53
1:A:36:ARG:CB	1:F:361:THR:HG21	2.39	0.53
1:C:335:LEU:HD22	1:C:339:ASP:CB	2.38	0.53
1:F:100:ILE:HG21	1:F:290:VAL:CG1	2.38	0.53
1:A:275:GLU:O	1:A:279:ARG:HG3	2.08	0.53
1:B:271:ASP:C	1:B:273:SER:N	2.63	0.53
1:D:79:ILE:CG2	1:D:103:LEU:HG	2.38	0.53
1:E:101:ARG:HG3	1:E:101:ARG:NH1	2.24	0.53
1:F:282:LEU:HB2	1:F:283:PRO:HD3	1.90	0.53
1:A:384:ASN:HD21	1:A:390:ILE:H	1.54	0.53
1:B:361:THR:HG23	1:C:39:GLN:HG2	1.89	0.53
1:A:18:ILE:N	2:A:900:ATP:N6	2.56	0.53
1:D:300:ASP:C	1:D:301:HIS:CD2	2.86	0.53
1:F:388:GLU:N	1:F:388:GLU:OE1	2.42	0.53
1:A:3:GLU:HG2	1:A:35:TRP:CG	2.45	0.52
1:A:325:ARG:HE	1:A:325:ARG:HA	1.72	0.52
1:B:109:LYS:O	1:B:113:VAL:HG23	2.10	0.52
1:D:312:ILE:HG13	1:D:313:ALA:N	2.23	0.52
1:E:132:LEU:HD23	1:E:135:LEU:HD12	1.91	0.52
1:C:240:LYS:NZ	1:C:241:GLN:HG3	2.23	0.52
1:C:292:THR:C	1:C:294:HIS:H	2.16	0.52
1:D:349:ALA:CB	1:E:47:GLU:HG3	2.39	0.52
1:E:18:ILE:N	2:E:915:ATP:HN61	2.06	0.52
1:F:240:LYS:CD	1:F:241:GLN:HG3	2.39	0.52
1:C:272:VAL:C	1:C:274:ARG:N	2.66	0.52
1:E:95:GLU:HA	1:E:95:GLU:OE1	2.09	0.52
1:B:55:MET:HE3	1:B:305:ILE:CG2	2.39	0.52
1:E:325:ARG:HE	1:E:325:ARG:HA	1.75	0.52
1:F:244:ILE:HG13	1:F:245:ASP:N	2.24	0.52
1:A:36:ARG:HG2	1:F:361:THR:HG21	1.92	0.52
1:C:374:LYS:O	1:C:378:GLU:HG3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:3:GLU:HG3	1:F:35:TRP:HB2	1.91	0.52
1:F:240:LYS:O	1:F:242:ASP:N	2.39	0.52
1:F:390:ILE:HG21	1:F:393:ARG:CB	2.40	0.52
1:F:394:ARG:HG3	1:F:394:ARG:O	2.08	0.52
1:A:361:THR:HG22	1:A:362:GLU:HG2	1.91	0.52
1:B:73:LEU:C	1:B:73:LEU:HD12	2.35	0.52
1:C:303:LEU:HG	1:C:305:ILE:CD1	2.40	0.52
1:E:74:ALA:O	1:E:75:ASN:C	2.53	0.52
1:F:111:VAL:HG21	1:F:239:LEU:HD11	1.92	0.52
1:F:257:GLU:HB2	1:F:260:LYS:CG	2.38	0.52
1:D:246:ALA:O	1:D:250:HIS:HB2	2.09	0.52
1:E:100:ILE:HB	1:E:290:VAL:HG21	1.90	0.52
1:B:261:ILE:HG22	1:B:278:GLN:NE2	2.20	0.52
1:D:30:ALA:HB2	1:D:53:ILE:HD11	1.90	0.52
1:F:436:GLU:O	1:F:439:SER:HB2	2.09	0.52
1:A:248:GLU:OE1	1:A:297:VAL:HA	2.10	0.52
1:B:441:PHE:HD1	1:C:56:ILE:HD13	1.75	0.52
1:C:44:LEU:O	1:C:45:ARG:C	2.53	0.52
1:D:248:GLU:OE1	1:D:297:VAL:HA	2.10	0.52
1:A:32:ARG:O	1:A:36:ARG:HG3	2.09	0.52
1:C:240:LYS:C	1:C:242:ASP:H	2.17	0.52
1:C:282:LEU:O	1:C:286:GLU:HG2	2.10	0.52
1:C:402:LEU:HD23	1:C:403:MET:SD	2.50	0.52
1:E:126:LEU:HA	1:E:129:GLU:HB3	1.92	0.52
1:A:409:ASP:O	1:A:410:ALA:C	2.54	0.51
1:B:4:MET:HE2	1:B:9:ILE:HA	1.91	0.51
1:B:101:ARG:HH11	1:B:101:ARG:CG	2.20	0.51
1:B:384:ASN:HD22	1:B:394:ARG:NH1	2.07	0.51
1:A:37:ARG:NE	1:A:301:HIS:CE1	2.78	0.51
1:C:108:VAL:HG21	1:C:294:HIS:ND1	2.24	0.51
1:D:18:ILE:N	2:D:910:ATP:N6	2.55	0.51
1:D:37:ARG:HB3	1:D:37:ARG:NH1	2.26	0.51
1:F:272:VAL:C	1:F:274:ARG:N	2.67	0.51
1:F:390:ILE:CG2	1:F:393:ARG:HB2	2.41	0.51
1:B:104:THR:HA	1:B:247:VAL:HG21	1.91	0.51
1:E:41:ASN:O	1:E:45:ARG:HB2	2.10	0.51
1:E:128:GLU:N	1:E:222:MET:HE1	2.25	0.51
1:E:279:ARG:HG2	1:E:279:ARG:HH11	1.75	0.51
1:A:44:LEU:O	1:A:45:ARG:C	2.53	0.51
1:A:361:THR:CG2	1:B:36:ARG:HG2	2.38	0.51
1:B:272:VAL:HG12	1:B:272:VAL:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:103:LEU:O	1:C:106:ALA:HB3	2.11	0.51
1:C:361:THR:HG21	1:D:36:ARG:HB3	1.92	0.51
1:A:38:MET:HE3	1:A:45:ARG:HE	1.74	0.51
1:A:246:ALA:O	1:A:250:HIS:HB2	2.10	0.51
1:B:261:ILE:CG2	1:B:261:ILE:O	2.57	0.51
1:B:280:ASP:O	1:B:283:PRO:HD2	2.10	0.51
1:D:249:GLN:C	1:D:250:HIS:ND1	2.68	0.51
1:D:421:ASP:OD2	1:D:423:ASP:HB2	2.10	0.51
1:A:294:HIS:N	1:A:294:HIS:CD2	2.79	0.51
1:A:432:LEU:N	1:A:432:LEU:CD1	2.73	0.51
1:B:17:ILE:O	1:B:24:LYS:NZ	2.40	0.51
1:C:104:THR:O	1:C:105:ASP:C	2.52	0.51
1:E:168:LEU:N	1:E:168:LEU:HD13	2.26	0.51
1:E:401:ARG:CZ	1:E:442:ILE:HG21	2.41	0.51
1:F:43:GLU:O	1:F:47:GLU:HG2	2.10	0.51
1:F:282:LEU:CD1	1:F:319:ILE:HD11	2.39	0.51
1:B:54:LEU:HD23	1:B:329:ARG:HD2	1.91	0.51
1:C:108:VAL:O	1:C:112:ARG:HB2	2.11	0.51
1:F:107:ALA:O	1:F:110:MET:HB2	2.11	0.51
1:A:63:LYS:HG2	1:A:332:LEU:HD13	1.93	0.51
1:A:361:THR:HG21	1:B:36:ARG:CG	2.37	0.51
1:D:96:VAL:O	1:D:97:ASP:C	2.54	0.51
1:E:17:ILE:CD1	1:E:66:ILE:HA	2.41	0.51
1:F:292:THR:C	1:F:294:HIS:H	2.18	0.51
1:B:37:ARG:O	1:B:45:ARG:HG3	2.10	0.51
1:B:126:LEU:HA	1:B:129:GLU:HB3	1.93	0.51
1:B:243:ALA:O	1:B:246:ALA:HB3	2.11	0.51
1:D:88:GLU:CD	1:D:93:GLY:HA3	2.36	0.51
1:D:351:ILE:HD12	1:D:399:LEU:HD12	1.92	0.51
1:E:169:ASP:O	1:E:170:ASP:OD1	2.28	0.51
1:C:441:PHE:HA	1:D:315:PRO:HG2	1.93	0.50
1:C:442:ILE:C	1:C:442:ILE:HD12	2.36	0.50
1:D:37:ARG:HB3	1:D:37:ARG:HH11	1.76	0.50
1:D:247:VAL:HG12	1:D:302:ILE:HD11	1.93	0.50
1:D:384:ASN:HD21	1:D:390:ILE:H	1.59	0.50
1:F:267:SER:HB2	1:F:270:PRO:CG	2.42	0.50
1:A:302:ILE:HB	1:A:304:PHE:HE1	1.76	0.50
1:B:41:ASN:O	1:B:45:ARG:HB2	2.12	0.50
1:C:53:ILE:HG23	1:C:328:ILE:CG2	2.40	0.50
1:C:264:ARG:HB2	1:C:274:ARG:HH11	1.76	0.50
1:E:272:VAL:HG12	1:E:272:VAL:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:349:ALA:CB	1:F:47:GLU:HG3	2.42	0.50
1:F:393:ARG:HH11	1:F:393:ARG:CG	2.23	0.50
1:A:37:ARG:HH11	1:A:37:ARG:HB3	1.76	0.50
1:B:435:ASP:HB3	1:B:438:LEU:HB2	1.94	0.50
1:E:264:ARG:HB2	1:E:274:ARG:NH1	2.12	0.50
1:E:442:ILE:HG23	1:F:329:ARG:HB2	1.93	0.50
1:F:240:LYS:NZ	1:F:241:GLN:HG3	2.25	0.50
1:B:84:THR:HG22	1:B:87:THR:HG21	1.94	0.50
1:F:358:LEU:O	1:F:361:THR:HB	2.11	0.50
1:F:384:ASN:ND2	1:F:394:ARG:HD2	2.27	0.50
1:D:337:THR:O	1:D:341:GLU:HG3	2.11	0.50
1:F:17:ILE:HB	1:F:24:LYS:HZ1	1.76	0.50
1:F:52:ASN:HB2	1:F:325:ARG:O	2.11	0.50
1:B:383:VAL:HG12	1:B:394:ARG:HH12	1.74	0.50
1:D:82:GLU:O	1:D:84:THR:N	2.44	0.50
1:E:96:VAL:HG12	1:E:284:LEU:HD11	1.93	0.50
1:E:224:LEU:HD13	1:E:224:LEU:C	2.37	0.50
1:E:268:SER:O	1:E:272:VAL:HG23	2.12	0.50
1:F:336:THR:O	1:F:339:ASP:HB2	2.11	0.50
1:D:44:LEU:O	1:D:45:ARG:C	2.55	0.50
1:D:240:LYS:HE2	1:D:294:HIS:O	2.11	0.50
1:E:168:LEU:HD23	1:E:219:LYS:CB	2.41	0.50
1:E:223:LYS:O	1:E:227:GLU:HG2	2.11	0.50
1:F:355:TYR:CE1	1:F:359:MET:HG3	2.46	0.50
1:D:257:GLU:HB3	1:D:260:LYS:CG	2.42	0.50
1:F:267:SER:HB2	1:F:270:PRO:CD	2.42	0.50
1:A:20:GLN:O	1:A:24:LYS:HG3	2.11	0.49
1:A:41:ASN:HD21	1:A:44:LEU:N	1.96	0.49
1:A:64:THR:HG21	1:A:68:ARG:HH11	1.76	0.49
1:B:131:ILE:O	1:B:134:VAL:HG12	2.11	0.49
1:B:168:LEU:HB3	1:B:219:LYS:HB2	1.94	0.49
1:C:355:TYR:CE1	1:C:359:MET:HG3	2.46	0.49
1:D:394:ARG:HG2	1:D:394:ARG:NH1	2.25	0.49
1:E:354:GLN:OE1	1:F:48:VAL:HG12	2.11	0.49
1:B:244:ILE:O	1:B:245:ASP:C	2.54	0.49
1:B:270:PRO:C	1:B:273:SER:HB3	2.36	0.49
1:C:282:LEU:HD12	1:C:322:LEU:HD13	1.94	0.49
1:D:369:THR:HG22	1:D:371:SER:H	1.77	0.49
1:E:168:LEU:N	1:E:168:LEU:CD1	2.76	0.49
1:F:55:MET:CE	1:F:63:LYS:HA	2.41	0.49
1:A:38:MET:CE	1:A:45:ARG:HE	2.25	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:54:LEU:HB2	1:E:326:LEU:HD13	1.93	0.49
1:F:369:THR:O	1:F:372:GLY:N	2.45	0.49
1:B:393:ARG:HG2	1:B:393:ARG:NH1	2.27	0.49
1:E:104:THR:HA	1:E:247:VAL:HG21	1.93	0.49
1:E:107:ALA:CB	1:E:247:VAL:HG23	2.42	0.49
1:B:233:LEU:N	1:B:233:LEU:HD22	2.28	0.49
1:D:280:ASP:O	1:D:283:PRO:HD2	2.12	0.49
1:E:332:LEU:N	1:E:332:LEU:HD12	2.27	0.49
1:A:37:ARG:HB3	1:A:37:ARG:NH1	2.27	0.49
1:A:79:ILE:CG2	1:A:103:LEU:HG	2.43	0.49
1:A:394:ARG:O	1:A:398:VAL:HG23	2.13	0.49
1:D:37:ARG:NE	1:D:301:HIS:CE1	2.80	0.49
1:D:294:HIS:N	1:D:294:HIS:CD2	2.79	0.49
1:D:394:ARG:HH11	1:D:394:ARG:CG	2.24	0.49
1:F:270:PRO:C	1:F:272:VAL:H	2.19	0.49
1:B:16:HIS:C	1:B:17:ILE:HG13	2.37	0.49
1:D:3:GLU:HG2	1:D:35:TRP:CB	2.37	0.49
1:E:175:ILE:HG22	1:E:176:ASP:H	1.77	0.49
1:F:350:SER:OG	1:F:353:VAL:HG23	2.12	0.49
1:B:17:ILE:HD13	1:B:66:ILE:CG1	2.41	0.49
1:C:16:HIS:CD2	1:C:69:ARG:HD2	2.47	0.49
1:C:393:ARG:HH11	1:C:393:ARG:CG	2.23	0.49
1:B:358:LEU:O	1:B:361:THR:HB	2.12	0.49
1:E:234:VAL:CG1	1:E:239:LEU:HD23	2.43	0.49
1:F:107:ALA:O	1:F:111:VAL:HG22	2.13	0.49
1:B:408:TYR:CD1	1:C:6:PRO:HB3	2.48	0.49
1:C:389:ASN:HD22	1:C:390:ILE:N	2.10	0.49
1:D:234:VAL:O	1:D:236:PRO:HD3	2.12	0.49
1:F:27:VAL:HB	1:F:70:LEU:HG	1.94	0.49
1:F:270:PRO:HG2	1:F:271:ASP:H	1.77	0.49
1:B:175:ILE:CD1	1:B:224:LEU:HD11	2.43	0.48
1:B:239:LEU:HD12	1:B:239:LEU:C	2.37	0.48
1:B:410:ALA:O	1:B:413:LEU:N	2.41	0.48
1:C:234:VAL:HG12	1:C:235:ASN:N	2.28	0.48
1:D:3:GLU:HG2	1:D:35:TRP:CD1	2.48	0.48
1:A:257:GLU:C	1:A:259:ASP:N	2.67	0.48
1:B:426:SER:HB3	1:B:430:ASP:OD1	2.12	0.48
1:E:346:GLU:O	1:E:347:PRO:C	2.55	0.48
1:A:100:ILE:O	1:A:104:THR:OG1	2.26	0.48
1:B:21:ASP:OD2	1:B:25:ARG:NH1	2.47	0.48
1:B:369:THR:HB	1:B:421:ASP:HA	1.93	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:325:ARG:CA	1:D:325:ARG:NE	2.74	0.48
1:A:17:ILE:HD11	1:A:66:ILE:HA	1.95	0.48
1:A:41:ASN:O	1:A:45:ARG:HG3	2.14	0.48
1:B:357:ALA:HB1	1:C:40:LEU:HD22	1.96	0.48
1:C:30:ALA:HB2	1:C:53:ILE:HD11	1.94	0.48
1:D:100:ILE:O	1:D:104:THR:OG1	2.28	0.48
1:D:293:LYS:C	1:D:294:HIS:HD2	2.21	0.48
1:F:103:LEU:O	1:F:106:ALA:HB3	2.13	0.48
1:B:83:ALA:HB1	1:B:261:ILE:HD11	1.93	0.48
1:F:104:THR:O	1:F:105:ASP:C	2.56	0.48
1:A:104:THR:O	1:A:108:VAL:HG23	2.13	0.48
1:A:282:LEU:C	1:A:286:GLU:HG3	2.38	0.48
1:B:84:THR:HA	1:B:87:THR:CG2	2.36	0.48
1:C:288:CYS:SG	1:C:289:THR:N	2.86	0.48
1:C:390:ILE:HG21	1:C:393:ARG:CB	2.43	0.48
1:C:394:ARG:O	1:C:394:ARG:HG3	2.12	0.48
1:D:17:ILE:HD11	1:D:66:ILE:HA	1.94	0.48
1:A:302:ILE:HB	1:A:304:PHE:CE1	2.48	0.48
1:B:136:ILE:CD1	1:B:159:PHE:HD2	2.16	0.48
1:B:369:THR:HG22	1:B:372:GLY:H	1.79	0.48
1:C:32:ARG:O	1:C:36:ARG:HG3	2.13	0.48
1:C:41:ASN:HD21	1:C:43:GLU:CB	2.26	0.48
1:A:394:ARG:HG2	1:A:394:ARG:NH1	2.25	0.48
1:D:70:LEU:O	1:D:70:LEU:HD22	2.13	0.48
1:E:9:ILE:CD1	1:E:31:LEU:HD12	2.44	0.48
1:A:84:THR:O	1:A:85:LYS:C	2.56	0.48
1:B:375:ARG:HB3	1:B:425:VAL:HG11	1.96	0.48
1:B:401:ARG:CZ	1:B:442:ILE:HG21	2.43	0.48
1:D:342:ARG:O	1:D:347:PRO:HD3	2.13	0.48
1:A:384:ASN:HD22	1:A:384:ASN:HA	1.48	0.48
1:B:23:ALA:HB2	1:B:332:LEU:HG	1.96	0.48
1:C:104:THR:O	1:C:107:ALA:N	2.47	0.48
1:C:299:THR:O	1:C:299:THR:CG2	2.62	0.48
1:C:359:MET:HG3	1:C:366:ILE:HD11	1.95	0.48
1:D:33:ASN:ND2	1:D:36:ARG:HD2	2.29	0.48
1:C:81:VAL:HG11	1:C:86:PHE:HE2	1.79	0.47
1:C:366:ILE:HD13	1:C:403:MET:CE	2.45	0.47
1:D:244:ILE:HG23	1:D:297:VAL:HG22	1.96	0.47
1:D:397:THR:HG21	1:D:443:LEU:C	2.39	0.47
1:E:111:VAL:HG21	1:E:243:ALA:HA	1.95	0.47
1:F:19:GLY:O	1:F:24:LYS:HE3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:108:VAL:HG13	1:F:239:LEU:HD21	1.95	0.47
1:A:71:ALA:HA	1:A:252:ILE:HD12	1.96	0.47
1:A:280:ASP:O	1:A:283:PRO:HD2	2.13	0.47
1:B:427:LYS:HD2	1:B:427:LYS:C	2.39	0.47
1:C:369:THR:OG1	1:C:421:ASP:HA	2.14	0.47
1:C:409:ASP:O	1:C:410:ALA:C	2.57	0.47
1:D:384:ASN:HD22	1:D:394:ARG:NH1	2.10	0.47
1:D:409:ASP:O	1:D:410:ALA:C	2.57	0.47
1:E:133:ASP:OD1	1:E:156:ARG:NH2	2.47	0.47
1:F:282:LEU:CB	1:F:283:PRO:HD3	2.44	0.47
1:A:325:ARG:HA	1:A:325:ARG:NE	2.30	0.47
1:C:32:ARG:HG3	1:C:36:ARG:HE	1.79	0.47
1:C:44:LEU:O	1:C:47:GLU:N	2.47	0.47
1:D:59:THR:HG22	1:D:60:GLY:N	2.30	0.47
1:E:175:ILE:CD1	1:E:224:LEU:HD11	2.39	0.47
1:E:393:ARG:HG2	1:E:393:ARG:NH1	2.27	0.47
1:E:400:GLU:HG3	1:F:327:PRO:HB2	1.97	0.47
1:F:81:VAL:HG11	1:F:86:PHE:HE2	1.79	0.47
1:C:284:LEU:HA	1:C:299:THR:HG21	1.96	0.47
1:C:361:THR:HG22	1:C:362:GLU:HG2	1.97	0.47
1:F:282:LEU:HD21	1:F:321:GLU:HG2	1.96	0.47
1:A:73:LEU:HD12	1:A:73:LEU:C	2.39	0.47
1:B:381:TRP:CD1	1:B:381:TRP:C	2.92	0.47
1:C:43:GLU:O	1:C:47:GLU:HG2	2.14	0.47
1:D:427:LYS:C	1:D:427:LYS:CD	2.86	0.47
1:E:108:VAL:HG21	1:E:294:HIS:CE1	2.49	0.47
1:E:128:GLU:C	1:E:130:ARG:H	2.21	0.47
1:F:5:THR:O	1:F:6:PRO:C	2.57	0.47
1:F:75:ASN:O	1:F:75:ASN:CG	2.56	0.47
1:D:79:ILE:HG22	1:D:103:LEU:HG	1.96	0.47
1:E:168:LEU:C	1:E:169:ASP:CG	2.78	0.47
1:E:216:LEU:CD1	1:E:217:LYS:HG2	2.39	0.47
1:A:248:GLU:O	1:A:301:HIS:HB2	2.14	0.47
1:A:277:VAL:O	1:A:281:LEU:HB2	2.15	0.47
1:B:77:PRO:HB2	1:B:103:LEU:CD2	2.43	0.47
1:B:151:GLU:O	1:B:153:SER:N	2.48	0.47
1:B:289:THR:HG21	1:B:296:MET:HE3	1.95	0.47
1:C:52:ASN:HB2	1:C:325:ARG:O	2.15	0.47
1:C:282:LEU:O	1:C:283:PRO:C	2.58	0.47
1:C:341:GLU:HG2	1:C:377:ALA:CB	2.44	0.47
1:D:271:ASP:O	1:D:275:GLU:HG3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:361:THR:HG21	1:E:36:ARG:CG	2.42	0.47
1:D:361:THR:CG2	1:E:36:ARG:HG2	2.44	0.47
1:E:128:GLU:C	1:E:130:ARG:N	2.72	0.47
1:F:38:MET:O	1:F:39:GLN:HB2	2.14	0.47
1:F:108:VAL:HG21	1:F:294:HIS:ND1	2.30	0.47
1:A:96:VAL:CG1	1:A:284:LEU:HD11	2.45	0.47
1:A:300:ASP:C	1:A:301:HIS:CD2	2.93	0.47
1:B:103:LEU:O	1:B:106:ALA:HB3	2.15	0.47
1:B:245:ASP:O	1:B:246:ALA:C	2.58	0.47
1:B:374:LYS:O	1:B:378:GLU:HG3	2.14	0.47
1:C:41:ASN:OD1	1:C:44:LEU:HG	2.15	0.47
1:C:390:ILE:HG23	1:D:320:PRO:HB3	1.96	0.47
1:D:303:LEU:HD12	1:D:304:PHE:N	2.29	0.47
1:D:384:ASN:ND2	1:D:394:ARG:HH12	2.11	0.47
1:E:345:THR:HG22	1:E:352:THR:HG21	1.97	0.47
1:A:64:THR:CG2	1:A:68:ARG:HH11	2.28	0.47
1:B:234:VAL:HG11	1:B:239:LEU:HD23	1.97	0.47
1:D:384:ASN:CA	1:D:394:ARG:HH12	2.25	0.47
1:E:84:THR:HA	1:E:87:THR:CG2	2.40	0.47
1:F:25:ARG:O	1:F:29:ILE:HG13	2.14	0.47
1:A:7:ARG:NH2	1:F:409:ASP:OD1	2.47	0.47
1:B:136:ILE:HD12	1:B:156:ARG:HA	1.97	0.47
1:B:241:GLN:NE2	1:B:245:ASP:OD2	2.48	0.47
1:C:305:ILE:HD12	1:C:305:ILE:N	2.30	0.47
1:C:374:LYS:O	1:C:377:ALA:HB3	2.15	0.47
1:E:56:ILE:HA	1:E:308:GLY:O	2.16	0.47
1:E:107:ALA:HB2	1:E:247:VAL:HG23	1.96	0.47
1:F:81:VAL:HG12	1:F:82:GLU:N	2.30	0.47
1:F:344:LEU:O	1:F:350:SER:HB2	2.15	0.47
1:A:94:LYS:HE3	1:B:95:GLU:CG	2.45	0.46
1:C:101:ARG:HG2	1:C:101:ARG:NH1	2.26	0.46
1:C:329:ARG:HG3	1:C:329:ARG:HH11	1.80	0.46
1:D:44:LEU:O	1:D:46:HIS:N	2.48	0.46
1:D:72:LYS:N	1:D:72:LYS:CD	2.78	0.46
1:E:280:ASP:O	1:E:283:PRO:HD2	2.15	0.46
1:F:23:ALA:HA	1:F:330:VAL:HG21	1.97	0.46
1:F:279:ARG:HG2	1:F:279:ARG:HH11	1.80	0.46
1:A:17:ILE:CD1	1:A:66:ILE:HA	2.45	0.46
1:A:31:LEU:HG	1:A:70:LEU:HD21	1.96	0.46
1:C:304:PHE:C	1:C:305:ILE:HD12	2.40	0.46
1:C:402:LEU:HB2	1:C:429:LEU:HD21	1.95	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:335:LEU:HD11	2:E:915:ATP:C2	2.50	0.46
1:E:411:SER:C	1:E:413:LEU:H	2.24	0.46
1:F:112:ARG:C	1:F:114:GLN:N	2.71	0.46
1:A:314:LYS:O	1:A:315:PRO:C	2.56	0.46
1:B:432:LEU:HD23	1:B:443:LEU:HD21	1.98	0.46
1:C:43:GLU:O	1:C:43:GLU:HG2	2.15	0.46
1:D:239:LEU:HD12	1:D:240:LYS:H	1.76	0.46
1:E:359:MET:HA	1:E:359:MET:HE2	1.98	0.46
1:A:96:VAL:HG13	1:A:99:ILE:HD12	1.96	0.46
1:B:435:ASP:OD1	1:B:438:LEU:HD23	2.15	0.46
1:C:75:ASN:ND2	1:C:114:GLN:HE22	2.14	0.46
1:E:19:GLY:O	1:E:24:LYS:HE3	2.16	0.46
1:F:37:ARG:O	1:F:39:GLN:N	2.47	0.46
1:A:84:THR:C	1:A:86:PHE:N	2.71	0.46
1:A:93:GLY:O	1:A:94:LYS:C	2.59	0.46
1:B:101:ARG:NH1	1:B:101:ARG:CG	2.78	0.46
1:B:325:ARG:HE	1:B:325:ARG:HA	1.79	0.46
1:D:271:ASP:HA	1:D:274:ARG:HH11	1.81	0.46
1:E:55:MET:HE3	1:E:305:ILE:CG2	2.45	0.46
1:E:83:ALA:HB1	1:E:261:ILE:HD11	1.98	0.46
1:E:382:GLN:HG2	1:E:433:VAL:CG1	2.45	0.46
1:A:351:ILE:HA	1:A:354:GLN:HB2	1.97	0.46
1:B:76:ALA:HB1	1:B:250:HIS:O	2.16	0.46
1:B:96:VAL:HG12	1:B:284:LEU:HD11	1.97	0.46
1:B:130:ARG:HH12	1:B:225:LEU:CD2	2.29	0.46
1:B:401:ARG:HD2	1:B:432:LEU:HD21	1.98	0.46
1:A:372:GLY:O	1:A:376:ILE:HG13	2.16	0.46
1:A:427:LYS:C	1:A:427:LYS:CD	2.86	0.46
1:C:18:ILE:HG21	1:C:339:ASP:HA	1.97	0.46
1:C:22:ASN:O	1:C:23:ALA:C	2.59	0.46
1:D:240:LYS:HE3	1:D:294:HIS:HA	1.97	0.46
1:D:383:VAL:O	1:D:387:THR:OG1	2.33	0.46
1:E:104:THR:O	1:E:105:ASP:C	2.58	0.46
1:E:270:PRO:CA	1:E:273:SER:HB3	2.46	0.46
1:F:355:TYR:CE1	1:F:403:MET:HE3	2.51	0.46
1:C:409:ASP:OD1	1:D:7:ARG:NH2	2.49	0.46
1:F:267:SER:O	1:F:270:PRO:HD2	2.16	0.46
1:A:54:LEU:HD23	1:A:329:ARG:HD2	1.98	0.46
1:B:168:LEU:O	1:B:169:ASP:HB2	2.12	0.46
1:C:57:GLY:O	1:C:309:ALA:HA	2.16	0.46
1:C:358:LEU:O	1:C:361:THR:HB	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:359:MET:HG3	1:C:366:ILE:HD12	1.98	0.46
1:E:44:LEU:O	1:E:48:VAL:HG23	2.16	0.46
1:F:41:ASN:OD1	1:F:44:LEU:HG	2.15	0.46
1:F:104:THR:HA	1:F:247:VAL:HG21	1.98	0.46
1:F:235:ASN:HB3	1:F:238:GLU:CB	2.43	0.46
1:B:346:GLU:O	1:B:347:PRO:O	2.33	0.46
1:D:17:ILE:CD1	1:D:66:ILE:HA	2.46	0.46
1:D:275:GLU:O	1:D:279:ARG:HG3	2.16	0.46
1:E:132:LEU:O	1:E:136:ILE:HD12	2.15	0.46
1:E:235:ASN:HB2	1:E:238:GLU:HB2	1.98	0.46
1:A:244:ILE:HG23	1:A:297:VAL:HG22	1.97	0.45
1:A:349:ALA:HB3	1:B:47:GLU:HG3	1.96	0.45
1:B:348:ASN:O	1:B:350:SER:N	2.45	0.45
1:C:359:MET:CE	1:C:410:ALA:HB2	2.46	0.45
1:F:293:LYS:O	1:F:293:LYS:HG3	2.15	0.45
1:B:153:SER:C	1:B:155:ALA:N	2.74	0.45
1:C:17:ILE:HB	1:C:24:LYS:NZ	2.31	0.45
1:E:130:ARG:NH2	1:E:225:LEU:HD22	2.22	0.45
1:F:282:LEU:CD1	1:F:322:LEU:HD13	2.47	0.45
1:A:65:GLU:OE2	1:A:65:GLU:HA	2.15	0.45
1:A:114:GLN:O	1:A:118:LYS:HB2	2.16	0.45
1:A:257:GLU:HB3	1:A:260:LYS:CG	2.46	0.45
1:A:344:LEU:O	1:A:352:THR:HB	2.16	0.45
1:E:128:GLU:CA	1:E:222:MET:HE1	2.46	0.45
1:F:53:ILE:O	1:F:305:ILE:HA	2.16	0.45
1:F:426:SER:HB3	1:F:430:ASP:OD1	2.15	0.45
1:A:234:VAL:O	1:A:236:PRO:HD3	2.16	0.45
1:B:431:ALA:O	1:B:434:ALA:HB3	2.16	0.45
1:C:25:ARG:HG2	1:C:26:SER:N	2.32	0.45
1:D:96:VAL:CG1	1:D:284:LEU:HD11	2.46	0.45
1:E:130:ARG:HA	1:E:133:ASP:OD2	2.16	0.45
1:F:4:MET:O	1:F:32:ARG:NH1	2.50	0.45
1:F:406:ILE:O	1:F:410:ALA:N	2.49	0.45
1:B:44:LEU:O	1:B:48:VAL:HG23	2.16	0.45
1:B:171:LYS:NZ	1:B:218:ILE:HG12	2.32	0.45
1:C:100:ILE:HG21	1:C:290:VAL:HG11	1.97	0.45
1:D:300:ASP:O	1:D:301:HIS:HD2	2.00	0.45
1:F:248:GLU:OE1	1:F:297:VAL:HA	2.16	0.45
1:B:130:ARG:HH12	1:B:225:LEU:HD22	1.80	0.45
1:B:274:ARG:NH1	1:B:274:ARG:HG3	2.31	0.45
1:C:257:GLU:HB2	1:C:260:LYS:CG	2.43	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:94:LYS:HE3	1:E:95:GLU:CG	2.47	0.45
1:D:289:THR:OG1	1:D:298:LYS:HD3	2.17	0.45
1:E:52:ASN:HB2	1:E:325:ARG:O	2.17	0.45
1:A:3:GLU:CG	1:A:35:TRP:HB2	2.38	0.45
1:A:312:ILE:HG13	1:A:313:ALA:H	1.82	0.45
1:B:130:ARG:NH1	1:B:225:LEU:CD2	2.79	0.45
1:B:132:LEU:HD13	1:B:156:ARG:CG	2.47	0.45
1:B:432:LEU:HB3	1:B:443:LEU:HD11	1.97	0.45
1:C:104:THR:HA	1:C:247:VAL:HG21	1.99	0.45
1:D:413:LEU:O	1:D:416:GLN:HG3	2.17	0.45
1:E:281:LEU:O	1:E:285:VAL:HG22	2.17	0.45
1:F:258:ILE:HG13	1:F:261:ILE:HD11	1.99	0.45
1:F:335:LEU:HD22	1:F:339:ASP:CB	2.45	0.45
1:F:341:GLU:HG2	1:F:377:ALA:CB	2.47	0.45
1:B:241:GLN:OE1	1:B:241:GLN:HA	2.17	0.45
1:B:332:LEU:N	1:B:332:LEU:HD12	2.31	0.45
1:B:403:MET:CE	1:B:406:ILE:HD12	2.46	0.45
1:C:81:VAL:HG12	1:C:82:GLU:N	2.32	0.45
1:D:400:GLU:OE1	1:E:328:ILE:HG12	2.17	0.45
1:E:57:GLY:O	1:E:309:ALA:HA	2.17	0.45
1:F:337:THR:O	1:F:341:GLU:N	2.44	0.45
1:F:359:MET:CE	1:F:410:ALA:HB2	2.47	0.45
1:A:25:ARG:HG3	1:A:25:ARG:HH11	1.81	0.45
1:A:359:MET:HE2	1:A:359:MET:HA	1.99	0.45
1:B:128:GLU:HA	1:B:222:MET:HE1	1.97	0.45
1:C:350:SER:OG	1:C:353:VAL:HG23	2.17	0.45
1:D:101:ARG:HG3	1:D:101:ARG:HH11	1.82	0.45
1:E:37:ARG:O	1:E:45:ARG:HG3	2.17	0.45
1:F:44:LEU:O	1:F:45:ARG:C	2.60	0.45
1:F:108:VAL:HA	1:F:111:VAL:HG22	1.98	0.45
1:F:274:ARG:HA	1:F:277:VAL:CG2	2.47	0.45
1:A:27:VAL:HG11	1:A:70:LEU:HG	1.92	0.45
1:A:73:LEU:HD12	1:A:74:ALA:N	2.32	0.45
1:B:382:GLN:HG2	1:B:433:VAL:CG1	2.47	0.45
1:C:109:LYS:O	1:C:113:VAL:HG23	2.16	0.45
1:A:235:ASN:CB	1:A:238:GLU:HB2	2.42	0.44
1:A:335:LEU:HD13	1:A:343:ILE:HD11	1.98	0.44
1:A:336:THR:O	1:A:339:ASP:HB2	2.17	0.44
1:C:371:SER:CB	1:C:422:ALA:HB2	2.39	0.44
1:C:393:ARG:HG3	1:C:393:ARG:NH1	2.28	0.44
1:E:88:GLU:HB2	1:E:91:TYR:H	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:171:LYS:O	1:E:216:LEU:N	2.50	0.44
1:E:409:ASP:HB3	1:E:413:LEU:HD22	1.99	0.44
1:F:35:TRP:O	1:F:35:TRP:CG	2.69	0.44
1:E:73:LEU:C	1:E:73:LEU:HD12	2.42	0.44
1:E:221:ALA:O	1:E:225:LEU:HG	2.18	0.44
1:F:41:ASN:ND2	1:F:43:GLU:H	2.14	0.44
1:F:112:ARG:C	1:F:114:GLN:H	2.25	0.44
1:A:259:ASP:C	1:A:261:ILE:H	2.25	0.44
1:A:261:ILE:HG23	1:A:278:GLN:HE21	1.83	0.44
1:B:216:LEU:CD1	1:B:217:LYS:H	2.30	0.44
1:C:356:LYS:HB2	1:C:356:LYS:HE3	1.72	0.44
1:C:390:ILE:CG2	1:C:393:ARG:HB2	2.48	0.44
1:E:78:PHE:CG	1:E:79:ILE:N	2.86	0.44
1:B:4:MET:HE1	1:B:12:GLU:HG2	1.99	0.44
1:C:52:ASN:CB	1:C:326:LEU:HD12	2.48	0.44
1:C:345:THR:HG22	1:C:373:ILE:HD13	2.00	0.44
1:D:18:ILE:N	2:D:910:ATP:HN61	2.01	0.44
1:D:257:GLU:C	1:D:259:ASP:H	2.25	0.44
1:E:409:ASP:O	1:E:411:SER:N	2.50	0.44
1:F:37:ARG:HD2	1:F:48:VAL:CG2	2.35	0.44
1:A:271:ASP:O	1:A:275:GLU:HG3	2.17	0.44
1:D:432:LEU:CD1	1:D:432:LEU:N	2.81	0.44
1:E:151:GLU:C	1:E:153:SER:N	2.76	0.44
1:E:173:ILE:N	1:E:173:ILE:CD1	2.73	0.44
1:A:262:CYS:O	1:A:263:LYS:C	2.60	0.44
1:A:325:ARG:CA	1:A:325:ARG:NE	2.79	0.44
1:A:379:ALA:O	1:A:383:VAL:HG23	2.18	0.44
1:C:350:SER:O	1:C:354:GLN:HG3	2.17	0.44
1:D:13:LEU:HD12	1:D:13:LEU:HA	1.78	0.44
1:A:241:GLN:NE2	1:A:245:ASP:OD2	2.51	0.44
1:A:261:ILE:HG21	1:A:278:GLN:HG2	2.00	0.44
1:A:358:LEU:O	1:A:361:THR:HB	2.16	0.44
1:B:50:PRO:O	1:B:325:ARG:NH2	2.43	0.44
1:B:216:LEU:HD13	1:B:217:LYS:CG	2.45	0.44
1:B:216:LEU:HD12	1:B:221:ALA:HB2	2.00	0.44
1:C:270:PRO:C	1:C:272:VAL:H	2.26	0.44
1:E:435:ASP:OD2	1:E:438:LEU:HD23	2.18	0.44
1:A:282:LEU:O	1:A:286:GLU:N	2.48	0.44
1:B:410:ALA:O	1:B:411:SER:C	2.61	0.44
1:D:84:THR:C	1:D:86:PHE:N	2.75	0.44
1:D:293:LYS:C	1:D:294:HIS:CD2	2.95	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:359:MET:HE2	1:D:359:MET:CA	2.48	0.44
1:E:374:LYS:O	1:E:378:GLU:HG3	2.18	0.44
1:B:19:GLY:O	1:B:24:LYS:HE3	2.17	0.44
1:E:56:ILE:O	1:E:331:GLU:HA	2.18	0.44
1:E:239:LEU:C	1:E:239:LEU:CD1	2.90	0.44
1:F:248:GLU:OE1	1:F:298:LYS:N	2.48	0.44
1:F:289:THR:O	1:F:290:VAL:CB	2.59	0.44
1:A:36:ARG:CG	1:F:361:THR:HG21	2.47	0.43
1:C:48:VAL:HG23	1:C:48:VAL:O	2.18	0.43
1:C:282:LEU:HB2	1:C:283:PRO:CD	2.45	0.43
1:D:293:LYS:HB2	1:D:294:HIS:HD2	1.83	0.43
1:D:379:ALA:O	1:D:383:VAL:HG23	2.18	0.43
1:E:410:ALA:O	1:E:413:LEU:HB2	2.18	0.43
1:F:58:PRO:O	1:F:61:VAL:HG22	2.17	0.43
1:A:56:ILE:HB	1:A:331:GLU:HG2	2.00	0.43
1:B:70:LEU:C	1:B:70:LEU:HD13	2.44	0.43
1:B:169:ASP:C	1:B:170:ASP:CG	2.73	0.43
1:C:25:ARG:CG	1:C:26:SER:N	2.80	0.43
1:C:100:ILE:HG21	1:C:290:VAL:CG1	2.48	0.43
1:D:68:ARG:O	1:D:71:ALA:HB3	2.19	0.43
1:E:37:ARG:NH2	1:E:301:HIS:CE1	2.86	0.43
1:E:42:GLU:HA	1:E:45:ARG:HH12	1.82	0.43
1:F:248:GLU:CD	1:F:298:LYS:H	2.24	0.43
1:F:262:CYS:SG	1:F:318:LEU:HD23	2.58	0.43
1:F:322:LEU:O	1:F:323:GLN:C	2.61	0.43
1:B:274:ARG:HG3	1:B:274:ARG:HH11	1.84	0.43
1:C:29:ILE:HG22	1:C:33:ASN:OD1	2.19	0.43
1:C:32:ARG:HG3	1:C:36:ARG:NE	2.33	0.43
1:D:8:GLU:O	1:D:11:SER:HB2	2.19	0.43
1:D:18:ILE:HG23	1:D:342:ARG:CZ	2.48	0.43
1:D:262:CYS:O	1:D:263:LYS:C	2.61	0.43
1:A:72:LYS:N	1:A:72:LYS:CD	2.81	0.43
1:A:384:ASN:HA	1:A:394:ARG:HH12	1.82	0.43
1:C:346:GLU:HB2	1:C:347:PRO:HD3	1.99	0.43
1:D:394:ARG:O	1:D:398:VAL:HG23	2.18	0.43
1:E:233:LEU:N	1:E:233:LEU:HD22	2.34	0.43
1:E:355:TYR:CZ	1:E:403:MET:HB2	2.54	0.43
1:F:14:ASP:HA	1:F:24:LYS:NZ	2.34	0.43
1:F:85:LYS:O	1:F:88:GLU:HG2	2.17	0.43
1:F:266:GLU:CG	1:F:270:PRO:HG2	2.47	0.43
1:F:312:ILE:CG1	1:F:313:ALA:N	2.72	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:LYS:HE3	1:F:400:GLU:OE1	2.19	0.43
1:B:54:LEU:HB3	1:B:329:ARG:HD3	2.01	0.43
1:B:384:ASN:HD22	1:B:394:ARG:CZ	2.31	0.43
1:C:4:MET:O	1:C:32:ARG:NH1	2.51	0.43
1:C:264:ARG:HB2	1:C:274:ARG:NH1	2.34	0.43
1:F:108:VAL:HA	1:F:111:VAL:CG2	2.48	0.43
1:B:57:GLY:O	1:B:309:ALA:HA	2.18	0.43
1:B:130:ARG:HH11	1:B:130:ARG:HD2	1.66	0.43
1:B:133:ASP:OD1	1:B:156:ARG:NH2	2.51	0.43
1:E:131:ILE:O	1:E:135:LEU:HG	2.19	0.43
1:F:237:GLU:C	1:F:239:LEU:H	2.27	0.43
1:F:401:ARG:HD3	1:F:432:LEU:HD21	2.01	0.43
1:A:64:THR:HG21	1:A:68:ARG:NH1	2.33	0.43
1:B:155:ALA:O	1:B:158:ALA:HB3	2.19	0.43
1:B:261:ILE:HG23	1:B:274:ARG:O	2.18	0.43
1:C:279:ARG:HG2	1:C:279:ARG:HH11	1.82	0.43
1:E:77:PRO:HB2	1:E:103:LEU:HD22	2.00	0.43
1:E:344:LEU:HD12	1:E:351:ILE:HD11	1.93	0.43
1:A:18:ILE:HG23	1:A:342:ARG:CZ	2.49	0.43
1:B:136:ILE:O	1:B:137:PRO:C	2.61	0.43
1:B:153:SER:O	1:B:155:ALA:N	2.52	0.43
1:E:17:ILE:CD1	1:E:66:ILE:HG12	2.46	0.43
1:F:354:GLN:HE21	1:F:354:GLN:HB3	1.66	0.43
1:F:426:SER:O	1:F:430:ASP:CG	2.62	0.43
1:A:37:ARG:CZ	1:A:301:HIS:CE1	3.02	0.43
1:B:107:ALA:CB	1:B:247:VAL:HG23	2.47	0.43
1:B:107:ALA:HB2	1:B:247:VAL:HG23	2.00	0.43
1:B:132:LEU:HD12	1:B:156:ARG:HH11	1.84	0.43
1:C:44:LEU:O	1:C:46:HIS:N	2.52	0.43
1:C:319:ILE:O	1:C:319:ILE:HG13	2.18	0.43
1:C:361:THR:HG21	1:D:36:ARG:HG2	2.01	0.43
1:D:56:ILE:O	1:D:331:GLU:HA	2.19	0.43
1:D:271:ASP:OD2	1:D:271:ASP:N	2.51	0.43
1:D:426:SER:HB3	1:D:430:ASP:OD1	2.18	0.43
1:E:63:LYS:HB2	2:E:915:ATP:O1B	2.19	0.43
1:E:270:PRO:HA	1:E:273:SER:CB	2.49	0.43
1:F:409:ASP:O	1:F:410:ALA:C	2.61	0.43
1:A:23:ALA:HA	1:A:330:VAL:HG21	2.01	0.43
1:B:96:VAL:HG13	1:B:99:ILE:HD12	2.01	0.43
1:C:12:GLU:HG3	1:C:73:LEU:HD11	2.01	0.43
1:C:38:MET:O	1:C:39:GLN:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:244:ILE:HG13	1:C:245:ASP:OD2	2.19	0.43
1:D:93:GLY:O	1:D:94:LYS:C	2.60	0.43
1:F:390:ILE:HG21	1:F:393:ARG:HG3	2.00	0.43
1:A:93:GLY:O	1:A:94:LYS:O	2.37	0.42
1:B:124:GLU:O	1:B:127:ALA:HB3	2.19	0.42
1:C:355:TYR:O	1:C:356:LYS:C	2.62	0.42
1:C:393:ARG:CG	1:C:393:ARG:NH1	2.80	0.42
1:D:84:THR:O	1:D:85:LYS:C	2.60	0.42
1:E:282:LEU:HD23	1:E:282:LEU:HA	1.81	0.42
1:E:427:LYS:HE3	1:E:428:HIS:NE2	2.34	0.42
1:E:436:GLU:O	1:E:440:ARG:HG3	2.19	0.42
1:F:240:LYS:C	1:F:242:ASP:N	2.75	0.42
1:C:84:THR:HG23	1:C:260:LYS:HB2	2.01	0.42
1:D:72:LYS:O	1:D:75:ASN:N	2.52	0.42
1:E:103:LEU:O	1:E:106:ALA:HB3	2.19	0.42
1:E:359:MET:HE1	1:F:36:ARG:NH1	2.34	0.42
1:F:240:LYS:HE3	1:F:240:LYS:HB3	1.95	0.42
1:F:373:ILE:O	1:F:374:LYS:C	2.62	0.42
1:A:83:ALA:HB1	1:A:261:ILE:HD12	2.01	0.42
1:B:62:GLY:O	1:B:66:ILE:HG13	2.18	0.42
1:D:257:GLU:C	1:D:259:ASP:N	2.76	0.42
1:D:277:VAL:O	1:D:281:LEU:HB2	2.19	0.42
1:D:413:LEU:HA	1:D:416:GLN:OE1	2.20	0.42
1:E:279:ARG:O	1:E:282:LEU:HB2	2.19	0.42
1:F:236:PRO:HG2	1:F:237:GLU:H	1.84	0.42
1:A:57:GLY:O	1:A:309:ALA:HA	2.20	0.42
1:B:88:GLU:HB2	1:B:91:TYR:H	1.84	0.42
1:C:423:ASP:OD2	1:C:423:ASP:N	2.53	0.42
1:D:9:ILE:HG23	1:D:73:LEU:HD21	2.01	0.42
1:E:78:PHE:O	1:E:79:ILE:HB	2.19	0.42
1:E:84:THR:HG22	1:E:87:THR:HG21	2.01	0.42
1:E:101:ARG:HH11	1:E:101:ARG:CG	2.30	0.42
1:E:262:CYS:O	1:E:263:LYS:C	2.62	0.42
1:F:272:VAL:C	1:F:274:ARG:H	2.27	0.42
1:F:274:ARG:O	1:F:277:VAL:N	2.46	0.42
1:A:34:ARG:CZ	1:A:250:HIS:HA	2.49	0.42
1:A:87:THR:HG22	1:A:277:VAL:HG21	2.02	0.42
1:A:384:ASN:HD22	1:A:394:ARG:HH12	1.68	0.42
1:B:23:ALA:HA	1:B:330:VAL:HG21	2.00	0.42
1:B:128:GLU:CA	1:B:222:MET:HE1	2.49	0.42
1:B:279:ARG:HH11	1:B:279:ARG:HG2	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:174:GLU:C	1:E:175:ILE:HG13	2.45	0.42
1:E:381:TRP:CD1	1:E:381:TRP:C	2.97	0.42
1:A:348:ASN:HD22	1:A:348:ASN:HA	1.58	0.42
1:B:224:LEU:O	1:B:228:GLU:HG3	2.20	0.42
1:B:346:GLU:C	1:B:347:PRO:O	2.62	0.42
1:C:366:ILE:C	1:C:367:GLU:HG2	2.44	0.42
1:E:169:ASP:OD2	1:E:219:LYS:HB2	2.20	0.42
1:E:312:ILE:CG1	1:E:313:ALA:H	2.30	0.42
1:B:70:LEU:HD13	1:B:70:LEU:O	2.20	0.42
1:C:262:CYS:O	1:C:263:LYS:C	2.62	0.42
1:C:329:ARG:HG3	1:C:329:ARG:NH1	2.35	0.42
1:D:257:GLU:HB3	1:D:260:LYS:HG2	2.00	0.42
1:D:285:VAL:HG12	1:D:304:PHE:CG	2.54	0.42
1:D:395:LEU:O	1:D:399:LEU:HB2	2.20	0.42
1:F:282:LEU:HD21	1:F:321:GLU:CG	2.49	0.42
1:A:38:MET:HA	1:A:38:MET:CE	2.47	0.42
1:A:79:ILE:HG22	1:A:103:LEU:HG	2.02	0.42
1:B:172:GLU:O	1:B:172:GLU:HG3	2.20	0.42
1:C:20:GLN:HA	1:C:20:GLN:OE1	2.20	0.42
1:D:84:THR:O	1:D:86:PHE:N	2.52	0.42
1:D:346:GLU:O	1:D:347:PRO:C	2.63	0.42
1:E:96:VAL:HG11	1:E:281:LEU:HD23	2.02	0.42
1:E:369:THR:HG22	1:E:372:GLY:H	1.85	0.42
1:E:412:ASP:OD2	1:F:7:ARG:NH1	2.53	0.42
1:F:25:ARG:CG	1:F:26:SER:N	2.83	0.42
1:B:344:LEU:HD11	1:B:351:ILE:HD11	1.96	0.42
1:C:57:GLY:O	1:C:63:LYS:NZ	2.46	0.42
1:C:314:LYS:C	1:C:316:SER:N	2.77	0.42
1:B:61:VAL:C	2:B:905:ATP:C2	2.98	0.42
1:B:231:ALA:O	1:B:233:LEU:N	2.53	0.42
1:C:432:LEU:HB3	1:C:443:LEU:HD12	2.00	0.42
1:D:322:LEU:O	1:D:323:GLN:C	2.63	0.42
1:E:217:LYS:HG3	1:E:220:ASP:OD2	2.19	0.42
1:E:321:GLU:CD	1:E:321:GLU:H	2.28	0.42
1:F:74:ALA:O	1:F:75:ASN:OD1	2.37	0.42
1:F:393:ARG:CG	1:F:393:ARG:NH1	2.81	0.42
1:A:82:GLU:O	1:A:83:ALA:C	2.63	0.41
1:B:78:PHE:CG	1:B:79:ILE:N	2.88	0.41
1:B:218:ILE:C	1:B:220:ASP:N	2.76	0.41
1:C:236:PRO:C	1:C:238:GLU:H	2.28	0.41
1:D:31:LEU:HG	1:D:70:LEU:HD21	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:349:ALA:HB1	1:E:44:LEU:HD23	2.01	0.41
1:E:76:ALA:HB1	1:E:250:HIS:O	2.20	0.41
1:E:427:LYS:HD2	1:E:427:LYS:C	2.44	0.41
1:F:359:MET:HE1	1:F:410:ALA:HB2	2.02	0.41
1:A:44:LEU:O	1:A:46:HIS:N	2.53	0.41
1:B:224:LEU:HD13	1:B:224:LEU:C	2.45	0.41
1:B:262:CYS:O	1:B:263:LYS:C	2.63	0.41
1:C:5:THR:O	1:C:9:ILE:HG13	2.20	0.41
1:C:379:ALA:O	1:C:383:VAL:HG23	2.20	0.41
1:E:53:ILE:HG22	1:E:54:LEU:N	2.35	0.41
1:F:366:ILE:C	1:F:367:GLU:HG2	2.45	0.41
1:A:33:ASN:ND2	1:A:36:ARG:HD2	2.32	0.41
1:A:239:LEU:HD12	1:A:240:LYS:H	1.78	0.41
1:C:29:ILE:O	1:C:30:ALA:C	2.63	0.41
1:C:272:VAL:C	1:C:274:ARG:H	2.28	0.41
1:C:275:GLU:O	1:C:278:GLN:HB2	2.21	0.41
1:A:72:LYS:O	1:A:75:ASN:N	2.53	0.41
1:B:30:ALA:O	1:B:31:LEU:C	2.64	0.41
1:C:240:LYS:HZ2	1:C:241:GLN:HG3	1.84	0.41
1:A:29:ILE:O	1:A:30:ALA:C	2.63	0.41
1:A:33:ASN:HD22	1:A:33:ASN:HA	1.50	0.41
1:A:279:ARG:HG2	1:A:279:ARG:NH1	2.32	0.41
1:B:18:ILE:N	2:B:905:ATP:N6	2.65	0.41
1:B:257:GLU:HB3	1:B:260:LYS:HG3	2.02	0.41
1:C:50:PRO:HD3	1:C:301:HIS:HB3	2.02	0.41
1:E:42:GLU:O	1:E:43:GLU:C	2.64	0.41
1:E:357:ALA:HB1	1:F:40:LEU:HD22	2.01	0.41
1:F:76:ALA:HB1	1:F:250:HIS:O	2.20	0.41
1:F:264:ARG:HB2	1:F:274:ARG:NH1	2.35	0.41
1:A:111:VAL:O	1:A:112:ARG:C	2.64	0.41
1:B:66:ILE:O	1:B:70:LEU:HB2	2.21	0.41
1:B:409:ASP:O	1:B:411:SER:N	2.54	0.41
1:C:100:ILE:CD1	1:C:284:LEU:HD22	2.50	0.41
1:C:352:THR:HG22	1:C:353:VAL:N	2.35	0.41
1:C:373:ILE:O	1:C:374:LYS:C	2.60	0.41
1:C:441:PHE:CD1	1:D:56:ILE:HG21	2.55	0.41
1:E:412:ASP:O	1:E:413:LEU:HD12	2.21	0.41
1:F:53:ILE:HG23	1:F:328:ILE:HG22	2.01	0.41
1:F:340:PHE:CE1	1:F:392:ALA:HA	2.56	0.41
1:F:384:ASN:HD22	1:F:394:ARG:NH1	2.13	0.41
1:A:31:LEU:HD23	1:A:31:LEU:HA	1.91	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:LYS:HE3	1:A:300:ASP:OD2	2.21	0.41
1:B:384:ASN:ND2	1:B:394:ARG:CZ	2.83	0.41
1:C:390:ILE:CG2	1:C:393:ARG:HG3	2.49	0.41
1:E:369:THR:HB	1:E:421:ASP:HA	2.02	0.41
1:F:384:ASN:HD22	1:F:384:ASN:HA	1.64	0.41
1:C:108:VAL:HA	1:C:111:VAL:CG2	2.51	0.41
1:C:266:GLU:CG	1:C:270:PRO:HG2	2.47	0.41
1:C:282:LEU:C	1:C:286:GLU:HG2	2.46	0.41
1:D:34:ARG:HH12	1:D:250:HIS:C	2.29	0.41
1:D:111:VAL:O	1:D:112:ARG:C	2.63	0.41
1:E:54:LEU:HB3	1:E:329:ARG:HD3	2.03	0.41
1:E:239:LEU:HA	1:E:242:ASP:HB2	2.01	0.41
1:E:243:ALA:O	1:E:246:ALA:HB3	2.20	0.41
1:E:261:ILE:HG23	1:E:274:ARG:C	2.45	0.41
1:F:52:ASN:HB3	1:F:326:LEU:HD12	2.03	0.41
1:F:69:ARG:HD3	1:F:69:ARG:HA	1.93	0.41
1:F:113:VAL:O	1:F:113:VAL:HG12	2.21	0.41
1:A:36:ARG:HB3	1:F:361:THR:HG21	2.03	0.41
1:A:96:VAL:O	1:A:97:ASP:C	2.63	0.41
1:A:355:TYR:CE2	1:A:399:LEU:HD13	2.55	0.41
1:A:369:THR:HG22	1:A:371:SER:H	1.84	0.41
1:B:134:VAL:HG13	1:B:135:LEU:N	2.35	0.41
1:B:384:ASN:ND2	1:B:394:ARG:HB2	2.35	0.41
1:B:401:ARG:HA	1:B:404:GLU:HG3	2.01	0.41
1:C:27:VAL:HB	1:C:70:LEU:HG	2.02	0.41
1:C:55:MET:HE2	1:C:63:LYS:HA	2.02	0.41
1:C:390:ILE:HG21	1:C:393:ARG:CG	2.51	0.41
1:C:403:MET:O	1:C:404:GLU:C	2.64	0.41
1:D:41:ASN:HD21	1:D:44:LEU:N	1.95	0.41
1:D:42:GLU:O	1:D:43:GLU:C	2.64	0.41
1:D:293:LYS:HB2	1:D:294:HIS:CD2	2.56	0.41
1:E:4:MET:HE3	1:E:8:GLU:CB	2.51	0.41
1:E:119:ASN:C	1:E:119:ASN:ND2	2.70	0.41
1:E:175:ILE:CG2	1:E:176:ASP:H	2.32	0.41
1:E:350:SER:OG	1:E:353:VAL:HG23	2.21	0.41
1:E:387:THR:HA	1:E:440:ARG:NH2	2.36	0.41
1:E:390:ILE:O	1:E:393:ARG:HD2	2.21	0.41
1:F:269:GLY:N	1:F:270:PRO:HD2	2.36	0.41
1:F:282:LEU:HD21	1:F:321:GLU:CB	2.51	0.41
1:F:390:ILE:HG22	1:F:391:GLY:N	2.36	0.41
1:A:13:LEU:HD12	1:A:13:LEU:HA	1.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:ARG:O	1:A:29:ILE:HG13	2.20	0.41
1:B:17:ILE:CD1	1:B:66:ILE:HA	2.51	0.41
1:B:37:ARG:NH2	1:B:301:HIS:CE1	2.89	0.41
1:B:111:VAL:HG21	1:B:243:ALA:HA	2.01	0.41
1:C:267:SER:HB2	1:C:270:PRO:CG	2.51	0.41
1:D:420:ILE:HD12	1:D:424:TYR:CE1	2.56	0.41
1:E:275:GLU:O	1:E:278:GLN:HB2	2.21	0.41
1:F:50:PRO:HD3	1:F:301:HIS:HB3	2.02	0.41
1:A:315:PRO:HG2	1:F:441:PHE:HA	2.04	0.40
1:B:220:ASP:C	1:B:222:MET:N	2.78	0.40
1:B:229:GLU:O	1:B:233:LEU:CD2	2.69	0.40
1:B:282:LEU:HD23	1:B:282:LEU:HA	1.90	0.40
1:B:361:THR:HG22	1:C:35:TRP:CH2	2.56	0.40
1:C:62:GLY:O	1:C:63:LYS:C	2.64	0.40
1:C:270:PRO:HG2	1:C:271:ASP:H	1.85	0.40
1:C:359:MET:HE1	1:C:410:ALA:HB2	2.03	0.40
1:D:72:LYS:N	1:D:72:LYS:HD2	2.36	0.40
1:E:426:SER:HB3	1:E:430:ASP:OD1	2.21	0.40
1:F:403:MET:O	1:F:404:GLU:C	2.64	0.40
1:C:247:VAL:HG11	1:C:297:VAL:HG11	2.03	0.40
1:C:274:ARG:HA	1:C:277:VAL:CG2	2.48	0.40
1:C:354:GLN:O	1:C:357:ALA:HB3	2.21	0.40
1:E:314:LYS:C	1:E:316:SER:N	2.77	0.40
1:A:110:MET:O	1:A:114:GLN:N	2.54	0.40
1:A:358:LEU:C	1:A:359:MET:HE2	2.46	0.40
1:A:384:ASN:HD22	1:A:394:ARG:NH1	2.19	0.40
1:A:413:LEU:O	1:A:416:GLN:HG3	2.21	0.40
1:C:11:SER:O	1:C:14:ASP:HB2	2.21	0.40
1:C:234:VAL:CG1	1:C:235:ASN:N	2.84	0.40
1:D:261:ILE:HG21	1:D:278:GLN:HG2	2.03	0.40
1:F:271:ASP:O	1:F:271:ASP:CG	2.65	0.40
1:C:44:LEU:C	1:C:46:HIS:N	2.78	0.40
1:D:56:ILE:HB	1:D:331:GLU:HG2	2.03	0.40
1:E:53:ILE:HD13	1:E:328:ILE:HB	2.03	0.40
1:E:61:VAL:C	2:E:915:ATP:C2	3.00	0.40
1:E:77:PRO:HB2	1:E:103:LEU:CD2	2.50	0.40
1:F:85:LYS:CE	1:F:94:LYS:HE2	2.49	0.40
1:A:100:ILE:H	1:A:100:ILE:HG12	1.74	0.40
1:A:432:LEU:N	1:A:432:LEU:HD12	2.36	0.40
1:C:277:VAL:O	1:C:281:LEU:HG	2.21	0.40
1:C:284:LEU:O	1:C:299:THR:CG2	2.70	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:361:THR:HG21	1:D:36:ARG:CG	2.52	0.40
1:D:37:ARG:CZ	1:D:301:HIS:CE1	3.05	0.40
1:E:257:GLU:HB3	1:E:260:LYS:HG3	2.04	0.40
1:F:355:TYR:CE1	1:F:359:MET:CG	3.05	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	402/442 (91%)	328 (82%)	50 (12%)	24 (6%)	1	7
1	B	402/442 (91%)	337 (84%)	46 (11%)	19 (5%)	2	11
1	C	402/442 (91%)	297 (74%)	72 (18%)	33 (8%)	0	3
1	D	402/442 (91%)	332 (83%)	49 (12%)	21 (5%)	1	9
1	E	402/442 (91%)	343 (85%)	42 (10%)	17 (4%)	2	13
1	F	402/442 (91%)	305 (76%)	68 (17%)	29 (7%)	1	4
All	All	2412/2652 (91%)	1942 (80%)	327 (14%)	143 (6%)	1	7

All (143) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	83	ALA
1	A	116	ILE
1	A	167	GLN
1	A	264	ARG
1	B	153	SER
1	B	169	ASP
1	B	171	LYS
1	B	263	LYS

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Mol	Chain	Res	Type
1	B	264	ARG
1	B	348	ASN
1	C	39	GLN
1	C	167	GLN
1	C	236	PRO
1	C	263	LYS
1	C	290	VAL
1	D	83	ALA
1	D	167	GLN
1	D	264	ARG
1	E	169	ASP
1	E	263	LYS
1	E	264	ARG
1	E	348	ASN
1	E	410	ALA
1	F	39	GLN
1	F	116	ILE
1	F	167	GLN
1	F	234	VAL
1	F	263	LYS
1	F	290	VAL
1	F	387	THR
1	A	85	LYS
1	A	93	GLY
1	A	94	LYS
1	A	96	VAL
1	A	153	SER
1	A	263	LYS
1	A	410	ALA
1	B	175	ILE
1	B	232	LYS
1	B	410	ALA
1	C	4	MET
1	C	63	LYS
1	C	88	GLU
1	C	92	VAL
1	C	116	ILE
1	C	153	SER
1	C	175	ILE
1	C	266	GLU
1	C	287	GLY
1	C	390	ILE

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Mol	Chain	Res	Type
1	C	428	HIS
1	D	93	GLY
1	D	94	LYS
1	D	96	VAL
1	D	116	ILE
1	D	153	SER
1	D	175	ILE
1	D	263	LYS
1	D	410	ALA
1	E	116	ILE
1	E	175	ILE
1	E	232	LYS
1	F	63	LYS
1	F	92	VAL
1	F	153	SER
1	F	175	ILE
1	F	236	PRO
1	F	266	GLU
1	F	386	SER
1	F	390	ILE
1	F	428	HIS
1	A	112	ARG
1	A	141	ASN
1	A	175	ILE
1	B	143	TRP
1	C	140	LYS
1	C	141	ASN
1	C	387	THR
1	D	45	ARG
1	D	141	ASN
1	E	140	LYS
1	E	141	ASN
1	E	143	TRP
1	E	216	LEU
1	F	88	GLU
1	F	141	ASN
1	F	241	GLN
1	F	287	GLY
1	A	140	LYS
1	A	435	ASP
1	B	88	GLU
1	B	90	GLY

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Mol	Chain	Res	Type
1	B	141	ASN
1	B	170	ASP
1	C	112	ARG
1	C	117	GLU
1	C	168	LEU
1	C	218	ILE
1	C	241	GLN
1	D	140	LYS
1	D	168	LEU
1	D	435	ASP
1	E	90	GLY
1	E	153	SER
1	F	117	GLU
1	F	270	PRO
1	A	45	ARG
1	A	168	LEU
1	A	216	LEU
1	C	38	MET
1	C	137	PRO
1	C	216	LEU
1	C	269	GLY
1	C	270	PRO
1	C	386	SER
1	C	410	ALA
1	D	112	ARG
1	D	347	PRO
1	E	88	GLU
1	F	112	ARG
1	F	137	PRO
1	F	140	LYS
1	F	168	LEU
1	F	216	LEU
1	F	269	GLY
1	A	137	PRO
1	A	143	TRP
1	A	247	VAL
1	B	154	ALA
1	B	218	ILE
1	C	247	VAL
1	D	137	PRO
1	D	216	LEU
1	A	50	PRO

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Mol	Chain	Res	Type
1	A	218	ILE
1	B	92	VAL
1	B	116	ILE
1	C	226	ILE
1	E	218	ILE
1	F	218	ILE
1	B	347	PRO
1	D	218	ILE
1	E	92	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	348/376 (93%)	311 (89%)	37 (11%)	6	27
1	B	348/376 (93%)	310 (89%)	38 (11%)	6	26
1	C	348/376 (93%)	314 (90%)	34 (10%)	7	30
1	D	348/376 (93%)	310 (89%)	38 (11%)	6	26
1	E	348/376 (93%)	308 (88%)	40 (12%)	5	24
1	F	348/376 (93%)	319 (92%)	29 (8%)	10	37
All	All	2088/2256 (93%)	1872 (90%)	216 (10%)	7	28

All (216) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLU
1	A	13	LEU
1	A	27	VAL
1	A	31	LEU
1	A	33	ASN
1	A	41	ASN
1	A	53	ILE
1	A	70	LEU
1	A	81	VAL

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Mol	Chain	Res	Type
1	A	82	GLU
1	A	84	THR
1	A	103	LEU
1	A	114	GLN
1	A	128	GLU
1	A	136	ILE
1	A	148	GLN
1	A	171	LYS
1	A	173	ILE
1	A	216	LEU
1	A	235	ASN
1	A	258	ILE
1	A	261	ILE
1	A	264	ARG
1	A	281	LEU
1	A	294	HIS
1	A	325	ARG
1	A	338	SER
1	A	344	LEU
1	A	348	ASN
1	A	351	ILE
1	A	352	THR
1	A	361	THR
1	A	384	ASN
1	A	388	GLU
1	A	399	LEU
1	A	400	GLU
1	A	432	LEU
1	B	5	THR
1	B	6	PRO
1	B	13	LEU
1	B	27	VAL
1	B	31	LEU
1	B	41	ASN
1	B	59	THR
1	B	64	THR
1	B	79	ILE
1	B	81	VAL
1	B	88	GLU
1	B	94	LYS
1	B	95	GLU
1	B	97	ASP

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Mol	Chain	Res	Type
1	B	103	LEU
1	B	122	ARG
1	B	128	GLU
1	B	140	LYS
1	B	145	GLN
1	B	148	GLN
1	B	168	LEU
1	B	173	ILE
1	B	216	LEU
1	B	233	LEU
1	B	238	GLU
1	B	239	LEU
1	B	258	ILE
1	B	274	ARG
1	B	290	VAL
1	B	344	LEU
1	B	352	THR
1	B	370	ASP
1	B	375	ARG
1	B	397	THR
1	B	399	LEU
1	B	413	LEU
1	B	419	THR
1	B	427	LYS
1	C	38	MET
1	C	41	ASN
1	C	42	GLU
1	C	72	LYS
1	C	79	ILE
1	C	82	GLU
1	C	92	VAL
1	C	103	LEU
1	C	128	GLU
1	C	136	ILE
1	C	140	LYS
1	C	145	GLN
1	C	148	GLN
1	C	173	ILE
1	C	216	LEU
1	C	233	LEU
1	C	236	PRO
1	C	289	THR

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Mol	Chain	Res	Type
1	C	299	THR
1	C	300	ASP
1	C	322	LEU
1	C	326	LEU
1	C	344	LEU
1	C	352	THR
1	C	361	THR
1	C	382	GLN
1	C	388	GLU
1	C	389	ASN
1	C	394	ARG
1	C	402	LEU
1	C	405	GLU
1	C	411	SER
1	C	423	ASP
1	C	427	LYS
1	D	3	GLU
1	D	13	LEU
1	D	27	VAL
1	D	33	ASN
1	D	41	ASN
1	D	53	ILE
1	D	70	LEU
1	D	81	VAL
1	D	82	GLU
1	D	103	LEU
1	D	128	GLU
1	D	136	ILE
1	D	145	GLN
1	D	173	ILE
1	D	216	LEU
1	D	219	LYS
1	D	235	ASN
1	D	258	ILE
1	D	261	ILE
1	D	264	ARG
1	D	271	ASP
1	D	290	VAL
1	D	294	HIS
1	D	303	LEU
1	D	325	ARG
1	D	338	SER

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Mol	Chain	Res	Type
1	D	344	LEU
1	D	348	ASN
1	D	351	ILE
1	D	352	THR
1	D	361	THR
1	D	384	ASN
1	D	388	GLU
1	D	399	LEU
1	D	400	GLU
1	D	413	LEU
1	D	427	LYS
1	D	432	LEU
1	E	5	THR
1	E	13	LEU
1	E	27	VAL
1	E	31	LEU
1	E	41	ASN
1	E	59	THR
1	E	64	THR
1	E	70	LEU
1	E	81	VAL
1	E	88	GLU
1	E	94	LYS
1	E	95	GLU
1	E	103	LEU
1	E	119	ASN
1	E	122	ARG
1	E	128	GLU
1	E	136	ILE
1	E	160	ARG
1	E	168	LEU
1	E	173	ILE
1	E	176	ASP
1	E	216	LEU
1	E	219	LYS
1	E	235	ASN
1	E	238	GLU
1	E	239	LEU
1	E	258	ILE
1	E	274	ARG
1	E	290	VAL
1	E	325	ARG

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Mol	Chain	Res	Type
1	E	337	THR
1	E	344	LEU
1	E	352	THR
1	E	371	SER
1	E	375	ARG
1	E	399	LEU
1	E	413	LEU
1	E	419	THR
1	E	427	LYS
1	E	432	LEU
1	F	38	MET
1	F	41	ASN
1	F	42	GLU
1	F	72	LYS
1	F	79	ILE
1	F	82	GLU
1	F	92	VAL
1	F	100	ILE
1	F	103	LEU
1	F	122	ARG
1	F	128	GLU
1	F	136	ILE
1	F	148	GLN
1	F	173	ILE
1	F	216	LEU
1	F	299	THR
1	F	300	ASP
1	F	322	LEU
1	F	326	LEU
1	F	344	LEU
1	F	352	THR
1	F	361	THR
1	F	382	GLN
1	F	388	GLU
1	F	400	GLU
1	F	402	LEU
1	F	405	GLU
1	F	411	SER
1	F	427	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (69) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	HIS
1	A	22	ASN
1	A	33	ASN
1	A	41	ASN
1	A	75	ASN
1	A	249	GLN
1	A	294	HIS
1	A	348	ASN
1	A	365	ASN
1	A	384	ASN
1	A	428	HIS
1	B	22	ASN
1	B	41	ASN
1	B	119	ASN
1	B	150	GLN
1	B	249	GLN
1	B	250	HIS
1	B	348	ASN
1	B	365	ASN
1	B	384	ASN
1	B	416	GLN
1	B	428	HIS
1	C	16	HIS
1	C	22	ASN
1	C	39	GLN
1	C	114	GLN
1	C	241	GLN
1	C	250	HIS
1	C	294	HIS
1	C	311	GLN
1	C	348	ASN
1	C	354	GLN
1	C	389	ASN
1	C	416	GLN
1	C	417	ASN
1	D	22	ASN
1	D	33	ASN
1	D	41	ASN
1	D	119	ASN
1	D	249	GLN
1	D	294	HIS
1	D	301	HIS
1	D	333	GLN

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Mol	Chain	Res	Type
1	D	348	ASN
1	D	384	ASN
1	E	16	HIS
1	E	22	ASN
1	E	41	ASN
1	E	119	ASN
1	E	249	GLN
1	E	250	HIS
1	E	333	GLN
1	E	348	ASN
1	E	365	ASN
1	E	384	ASN
1	E	416	GLN
1	E	428	HIS
1	F	16	HIS
1	F	22	ASN
1	F	39	GLN
1	F	75	ASN
1	F	250	HIS
1	F	311	GLN
1	F	354	GLN
1	F	365	ASN
1	F	384	ASN
1	F	389	ASN
1	F	416	GLN
1	F	417	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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
4	SO4	F	805	-	4,4,4	1.79	2 (50%)	6,6,6	0.92	0
2	ATP	A	900	-	32,33,33	0.77	0	48,52,52	0.68	1 (2%)
4	SO4	C	800	-	4,4,4	1.86	2 (50%)	6,6,6	0.89	0
2	ATP	D	910	-	32,33,33	0.79	1 (3%)	48,52,52	0.64	0
2	ATP	B	905	3	32,33,33	0.77	0	48,52,52	0.67	0
2	ATP	E	915	3	32,33,33	0.78	0	48,52,52	0.68	0

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	ATP	D	910	-	-	4/22/38/38	0/3/3/3
2	ATP	E	915	3	-	4/22/38/38	0/3/3/3
2	ATP	B	905	3	-	4/22/38/38	0/3/3/3
2	ATP	A	900	-	-	5/22/38/38	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	800	SO4	O1-S	3.05	1.62	1.44
4	F	805	SO4	O1-S	2.93	1.62	1.44
2	D	910	ATP	PB-O3B	2.14	1.61	1.59
4	C	800	SO4	O3-S	-2.09	1.31	1.48
4	F	805	SO4	O3-S	-2.00	1.31	1.48

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	A	900	ATP	O3A-PB-O1B	-2.05	104.54	110.70

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	900	ATP	O4'-C1'-N9-C8
2	A	900	ATP	O4'-C1'-N9-C4
2	B	905	ATP	O4'-C1'-N9-C8
2	B	905	ATP	O4'-C1'-N9-C4
2	D	910	ATP	O4'-C1'-N9-C8
2	D	910	ATP	O4'-C1'-N9-C4
2	E	915	ATP	O4'-C1'-N9-C8
2	E	915	ATP	O4'-C1'-N9-C4
2	A	900	ATP	O4'-C4'-C5'-O5'
2	A	900	ATP	C3'-C4'-C5'-O5'
2	D	910	ATP	O4'-C4'-C5'-O5'
2	B	905	ATP	O4'-C4'-C5'-O5'
2	D	910	ATP	C3'-C4'-C5'-O5'
2	E	915	ATP	O4'-C4'-C5'-O5'
2	E	915	ATP	C3'-C4'-C5'-O5'
2	B	905	ATP	C3'-C4'-C5'-O5'
2	A	900	ATP	C5'-O5'-PA-O1A

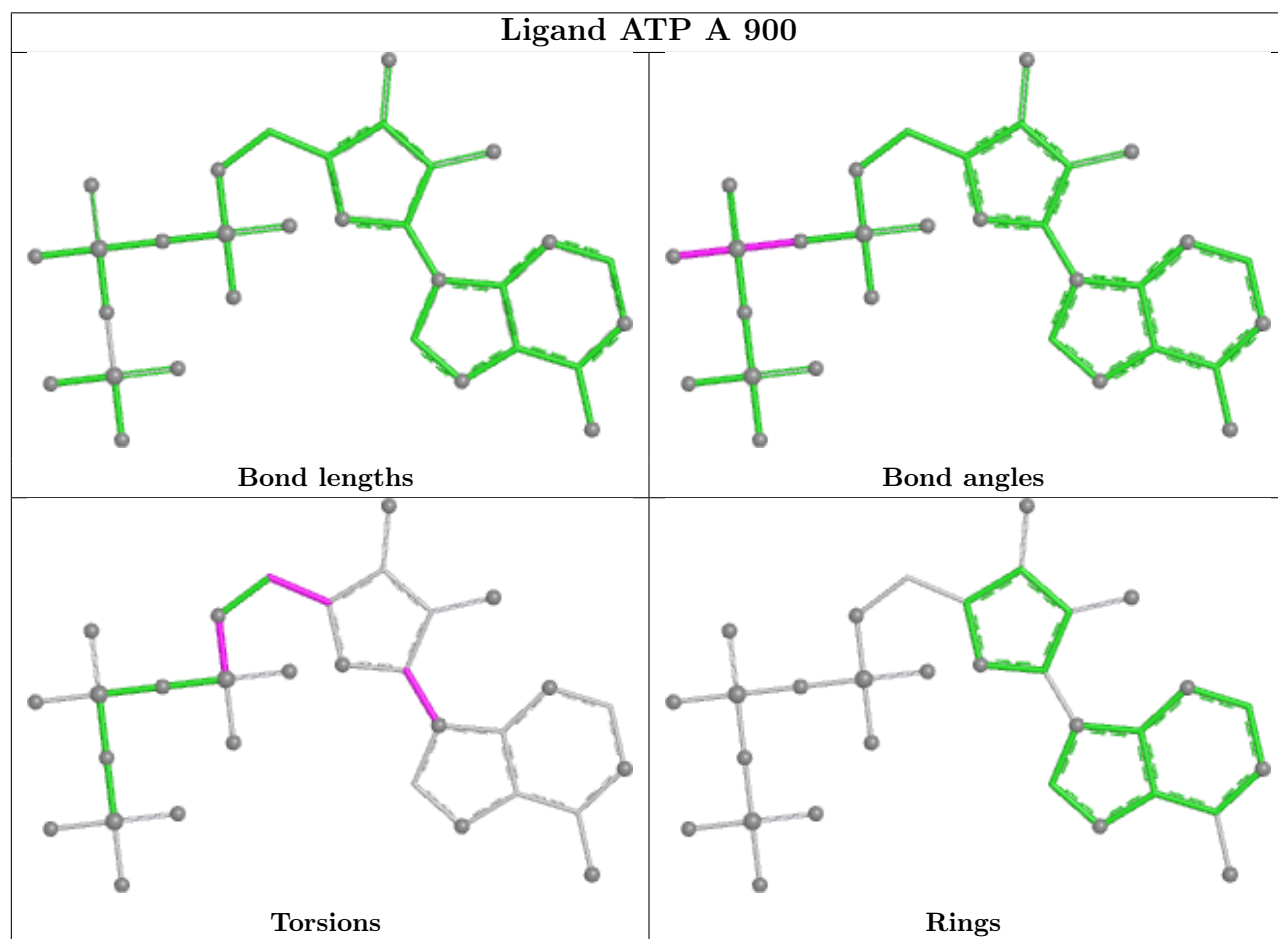
There are no ring outliers.

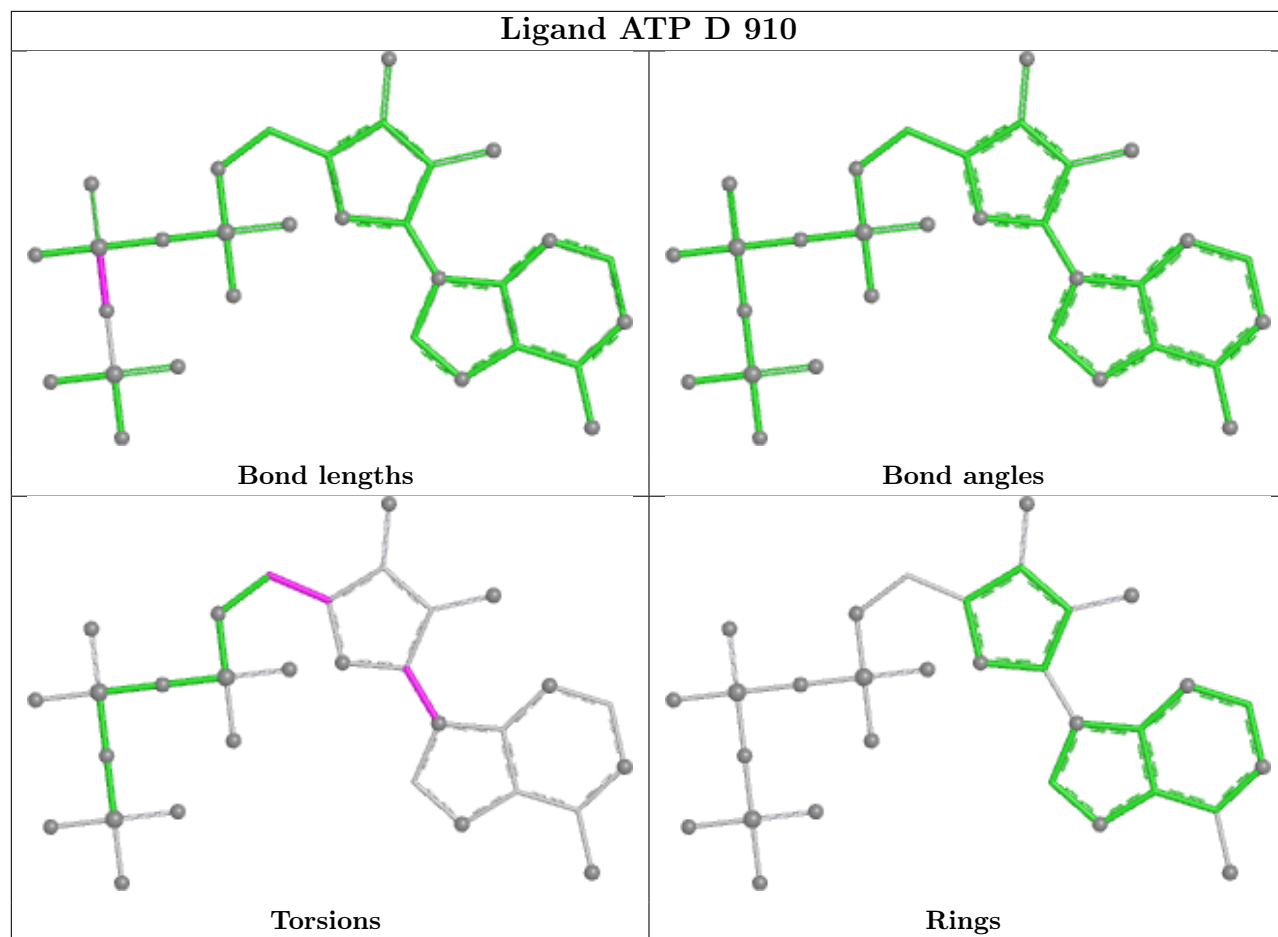
4 monomers are involved in 19 short contacts:

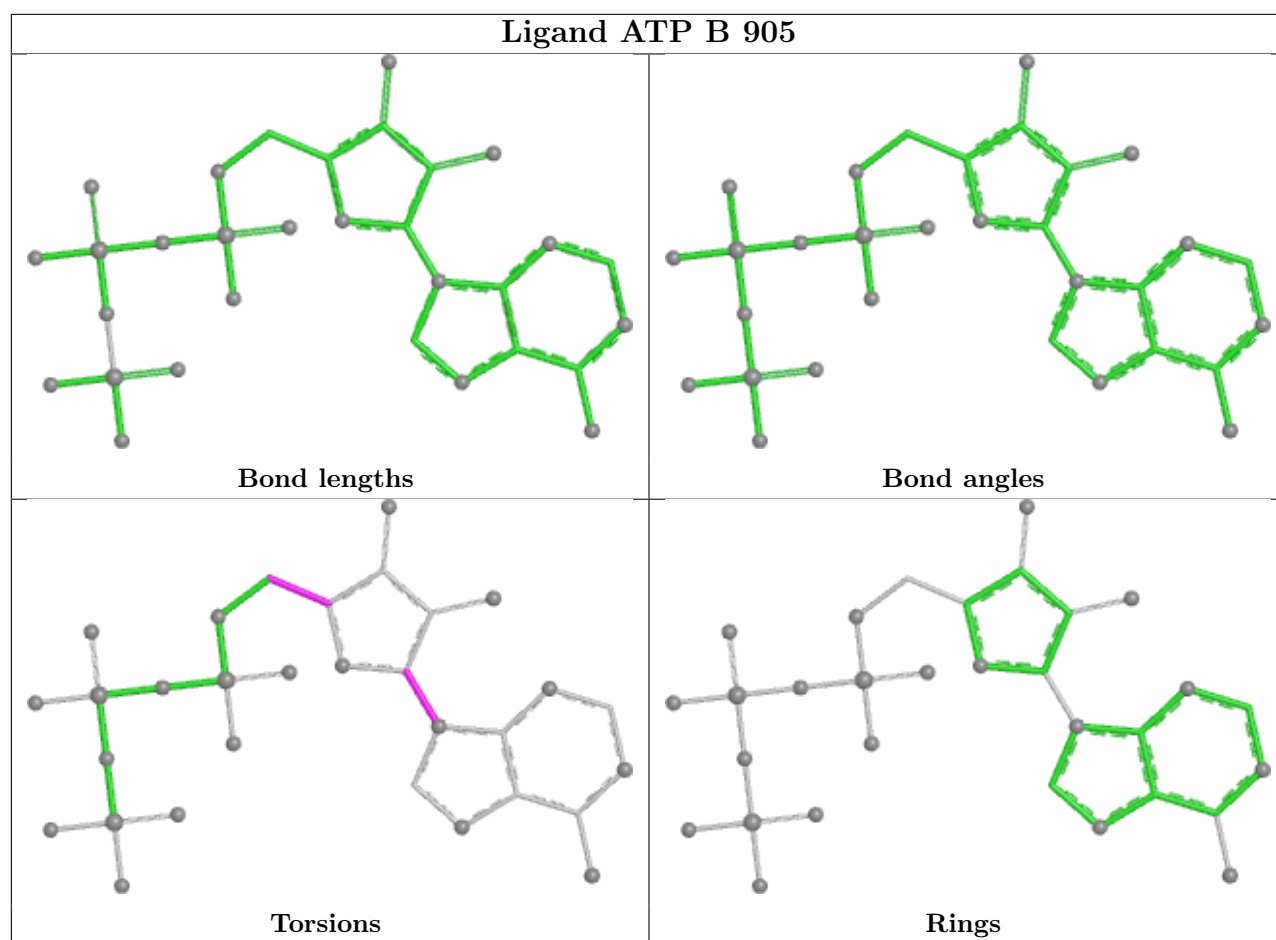
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	900	ATP	5	0
2	D	910	ATP	5	0
2	B	905	ATP	4	0
2	E	915	ATP	5	0

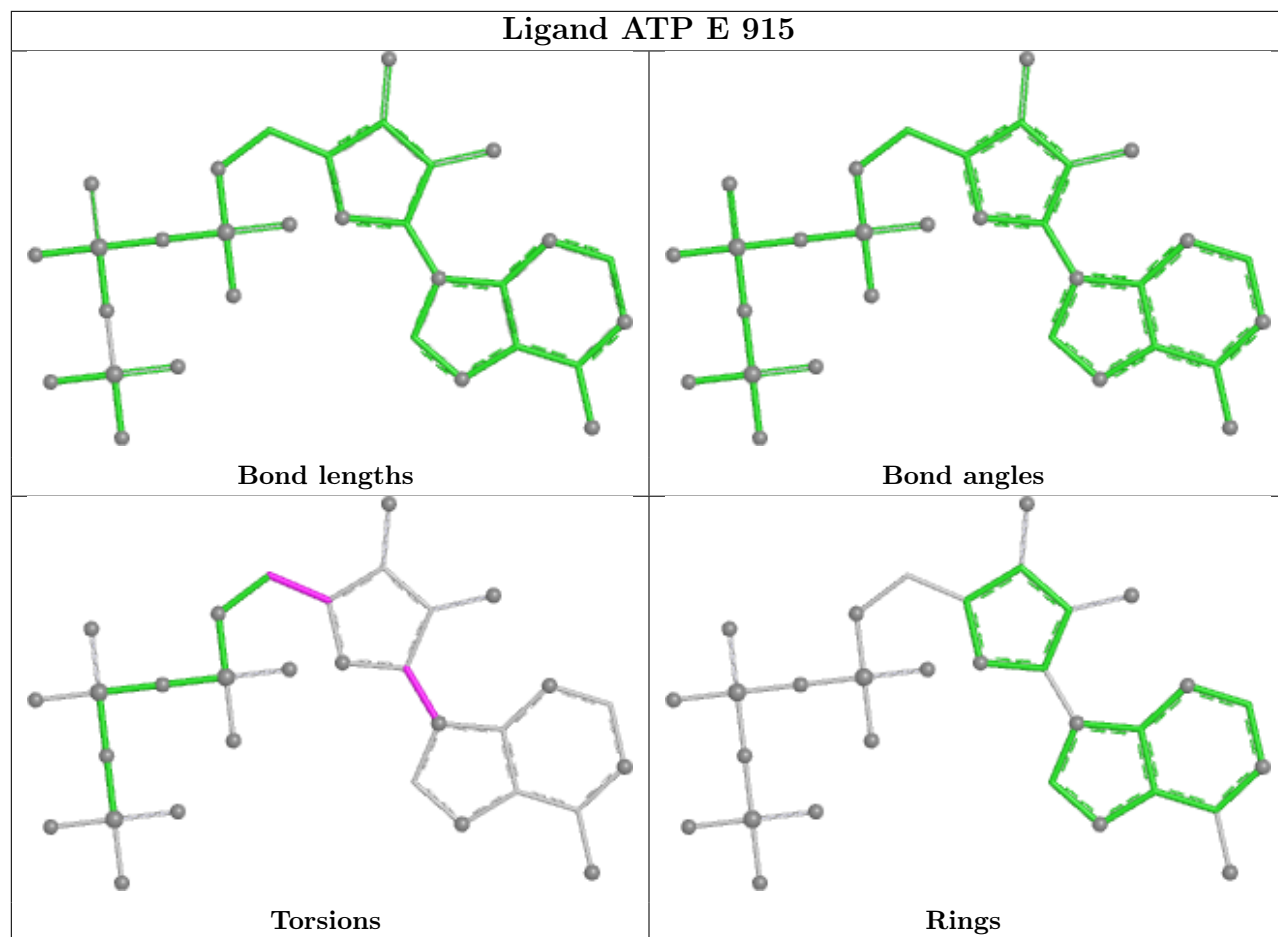
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.