



Full wwPDB X-ray Structure Validation Report ⓘ

May 7, 2026 – 11:37 AM EDT

PDB ID : 4M9Z / pdb_00004m9z
Title : Crystal structure of CED-4 bound CED-3 fragment
Authors : Huang, W.J.; Jinag, T.Y.; Choi, W.Y.; Wang, J.W.; Shi, Y.G.
Deposited on : 2013-08-15
Resolution : 3.40 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:

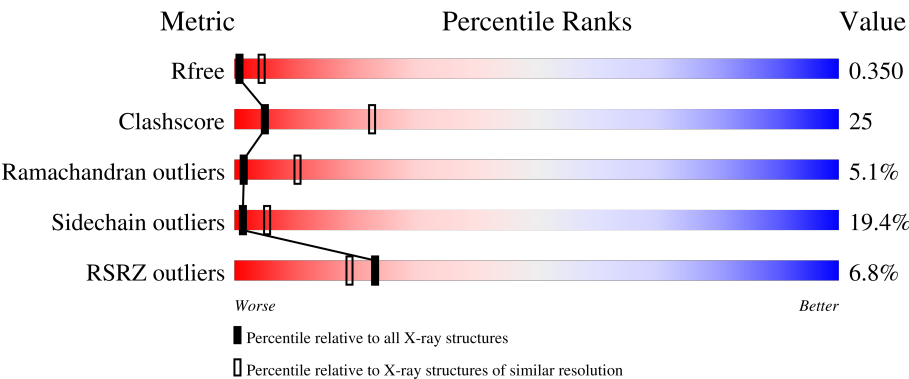
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal 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 i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	549	4% 41% 38% 12% 7%
1	B	549	7% 38% 41% 12% 7%
1	C	549	6% 40% 39% 12% 7%
1	D	549	7% 39% 40% 13% 7%
2	E	8	12% 38% 25% 12% 25%

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Mol	Chain	Length	Quality of chain
2	F	8	 25% 25% 12% 38%
2	G	8	 50% 12% 38%
2	H	8	 12% 50% 12% 25%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 16687 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 Cell death protein 4.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	510	Total	C	N	O	S	Se	0	0	0
			4088	2606	685	767	13	17			
1	B	511	Total	C	N	O	S	Se	0	0	0
			4094	2607	688	768	14	17			
1	C	510	Total	C	N	O	S	Se	0	0	0
			4088	2606	685	767	13	17			
1	D	511	Total	C	N	O	S	Se	0	0	0
			4094	2607	688	768	14	17			

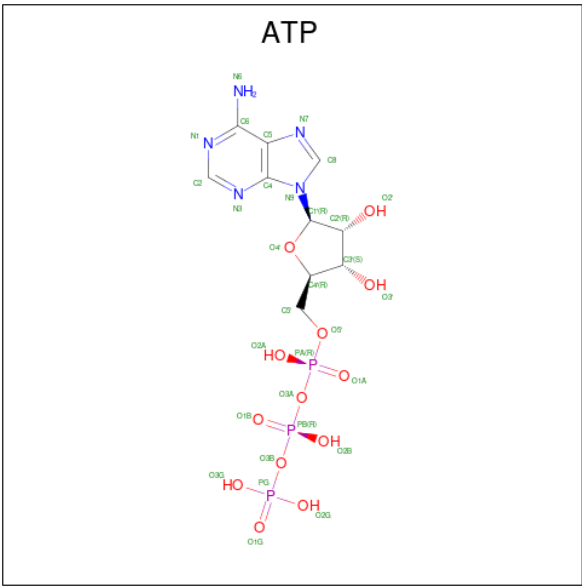
- Molecule 2 is a protein called CED-3 fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	6	Total	C	N	O	Se	0	0	0
			50	35	7	7	1			
2	F	5	Total	C	N	O	Se	0	0	0
			46	33	6	6	1			
2	G	5	Total	C	N	O	Se	0	0	0
			46	33	6	6	1			
2	H	6	Total	C	N	O	Se	0	0	0
			53	38	7	7	1			

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

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

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).

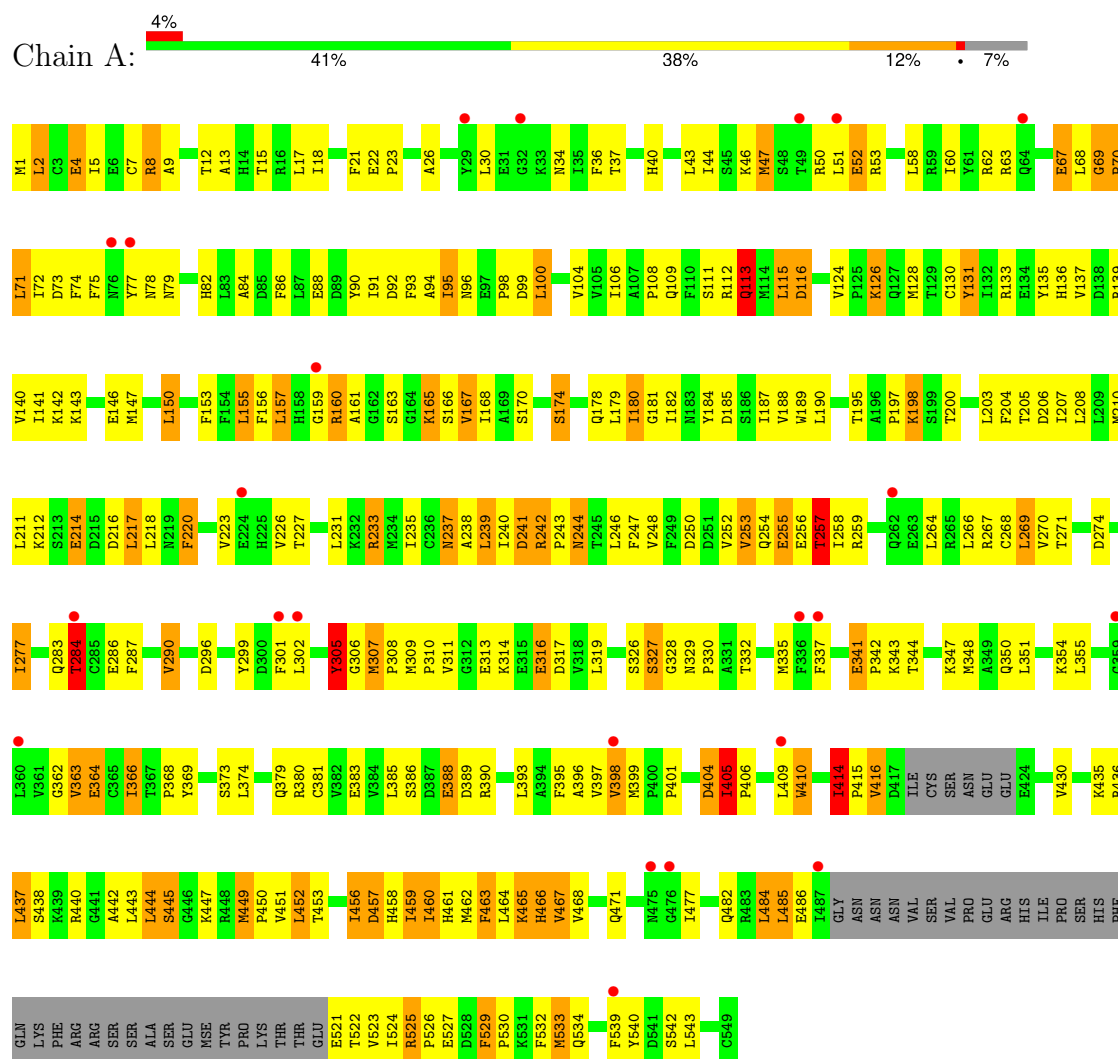


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
4	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
4	C	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
4	D	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

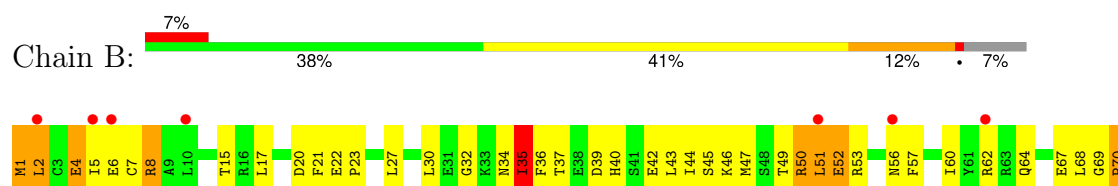
3 Residue-property plots

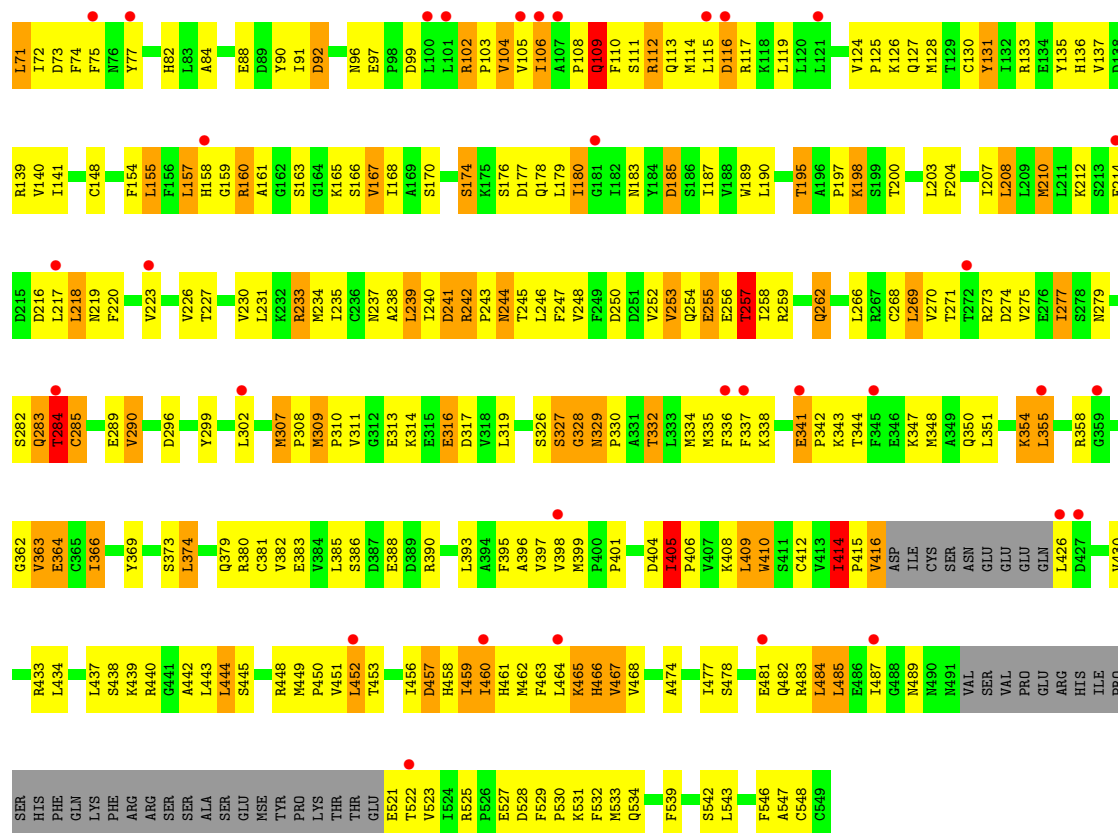
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Cell death protein 4

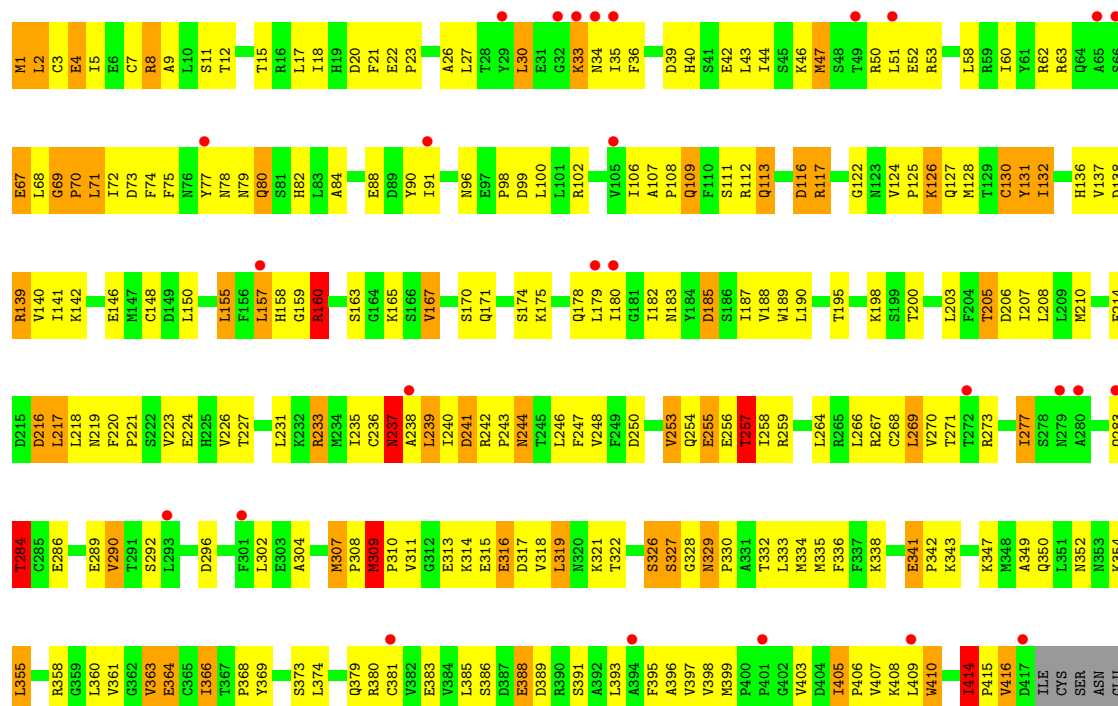
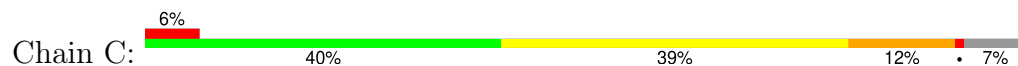


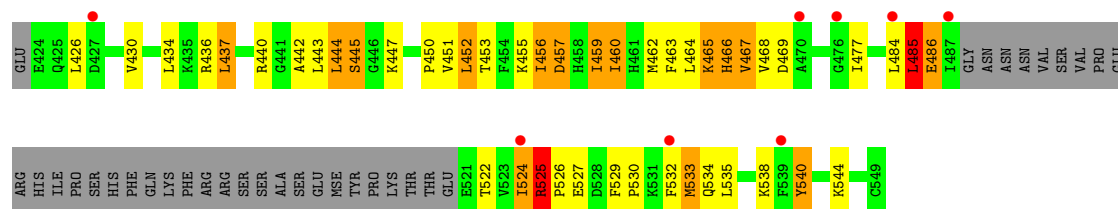
• Molecule 1: Cell death protein 4



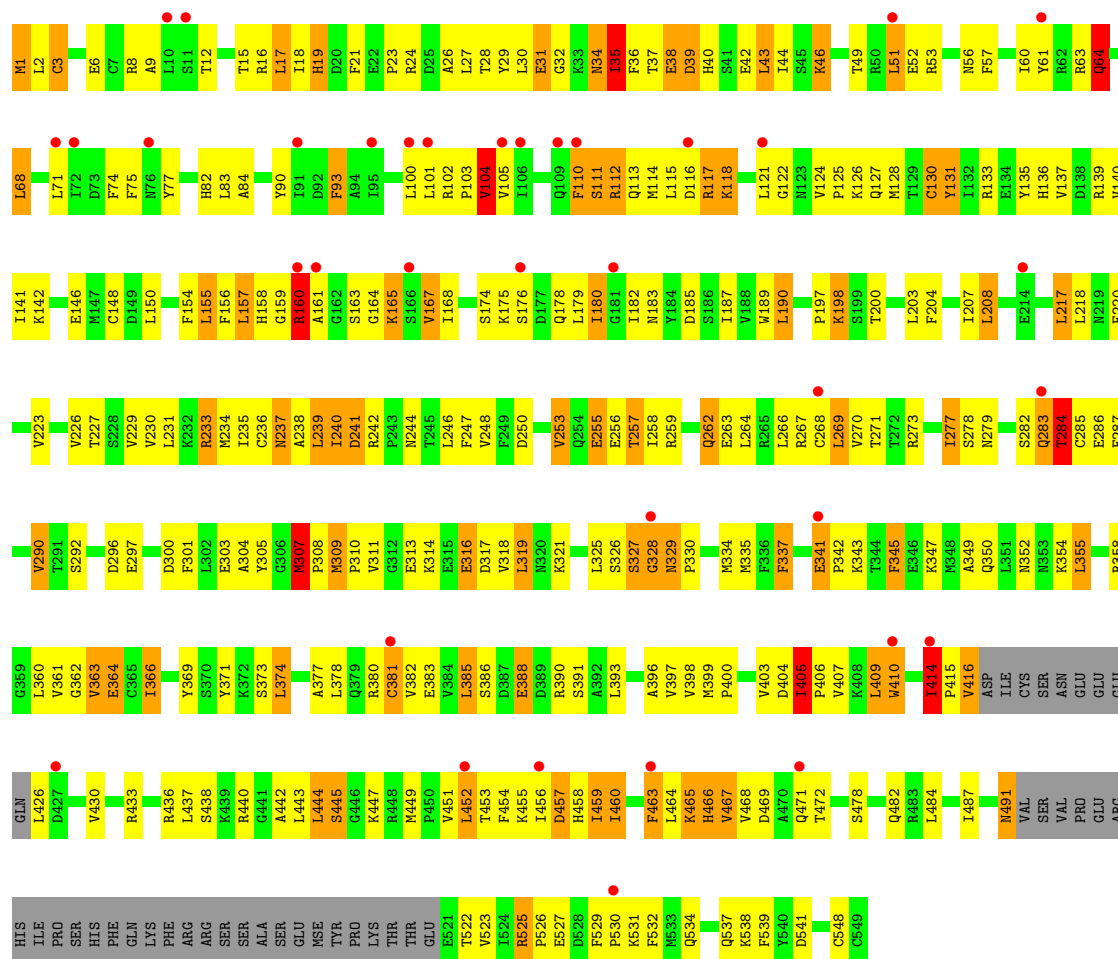


• Molecule 1: Cell death protein 4

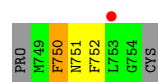




- Molecule 1: Cell death protein 4



- Molecule 2: CED-3 fragment



- Molecule 2: CED-3 fragment

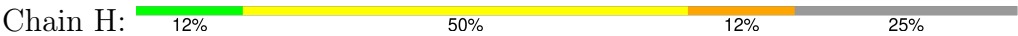




● Molecule 2: CED-3 fragment



● Molecule 2: CED-3 fragment



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	178.17Å 178.17Å 200.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.21 – 3.40 43.21 – 3.40	Depositor EDS
% Data completeness (in resolution range)	57.0 (43.21-3.40) 78.5 (43.21-3.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.63 (at 3.40Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.270 , 0.327 0.293 , 0.350	Depositor DCC
R_{free} test set	1759 reflections (3.92%)	wwPDB-VP
Wilson B-factor (Å ²)	81.5	Xtriage
Anisotropy	0.321	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 101.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	16687	wwPDB-VP
Average B, all atoms (Å ²)	121.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 27.22 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.2870e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: 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	0.46	0/4147	1.14	30/5577 (0.5%)
1	B	0.46	1/4153 (0.0%)	1.13	26/5585 (0.5%)
1	C	0.43	0/4147	1.09	24/5577 (0.4%)
1	D	0.43	0/4153	1.14	31/5585 (0.6%)
2	E	0.30	0/51	0.61	0/66
2	F	0.29	0/47	0.50	0/61
2	G	0.24	0/47	0.73	0/61
2	H	0.43	0/54	0.99	0/69
All	All	0.45	1/16799 (0.0%)	1.12	111/22581 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	285	CYS	CA-C	-5.08	1.46	1.52

All (111) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	405	ILE	CA-C-N	9.33	129.90	119.92
1	A	405	ILE	C-N-CA	9.33	129.90	119.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	484	LEU	N-CA-C	-9.15	101.25	111.14
1	D	414	ILE	CA-C-N	9.15	130.35	120.11
1	D	414	ILE	C-N-CA	9.15	130.35	120.11
1	B	484	LEU	N-CA-C	-8.85	101.58	111.14
1	B	307	MSE	CA-C-N	8.43	128.93	119.83
1	B	307	MSE	C-N-CA	8.43	128.93	119.83
1	C	217	LEU	N-CA-C	-8.11	103.59	113.97
1	D	329	ASN	CA-C-N	7.86	129.67	119.84
1	D	329	ASN	C-N-CA	7.86	129.67	119.84
1	A	540	TYR	N-CA-C	-7.76	104.42	114.04
1	D	112	ARG	N-CA-C	-7.61	104.89	114.56
1	D	117	ARG	N-CA-C	7.43	119.07	110.97
1	D	309	MSE	CA-C-N	7.43	128.43	120.11
1	D	309	MSE	C-N-CA	7.43	128.43	120.11
1	B	329	ASN	CA-C-N	7.35	129.03	119.84
1	B	329	ASN	C-N-CA	7.35	129.03	119.84
1	D	118	LYS	N-CA-C	7.30	119.32	111.36
1	D	405	ILE	CA-C-N	7.06	127.47	119.92
1	D	405	ILE	C-N-CA	7.06	127.47	119.92
1	D	242	ARG	CA-C-N	6.85	126.77	119.85
1	D	242	ARG	C-N-CA	6.85	126.77	119.85
1	B	405	ILE	CA-C-N	6.78	127.17	119.92
1	B	405	ILE	C-N-CA	6.78	127.17	119.92
1	A	69	GLY	CA-C-N	6.63	128.13	119.84
1	A	69	GLY	C-N-CA	6.63	128.13	119.84
1	C	329	ASN	CA-C-N	6.62	128.12	119.84
1	C	329	ASN	C-N-CA	6.62	128.12	119.84
1	D	268	CYS	N-CA-C	6.62	119.69	108.90
1	D	19	HIS	N-CA-C	6.59	118.46	111.28
1	B	92	ASP	N-CA-C	-6.45	103.76	111.69
1	D	253	VAL	N-CA-C	6.45	117.12	111.90
1	C	71	LEU	N-CA-C	-6.40	106.69	114.75
1	A	217	LEU	N-CA-C	-6.37	105.82	113.97
1	B	285	CYS	N-CA-C	-6.37	98.82	109.07
1	A	34	ASN	N-CA-C	6.36	118.35	110.91
1	A	329	ASN	CA-C-N	6.36	127.79	119.84
1	A	329	ASN	C-N-CA	6.36	127.79	119.84
1	C	214	GLU	N-CA-C	-6.35	104.36	111.28
1	C	525	ARG	CA-C-N	6.30	126.26	120.03
1	C	525	ARG	C-N-CA	6.30	126.26	120.03
1	B	393	LEU	N-CA-C	-6.29	104.43	111.28
1	B	242	ARG	CA-C-N	6.29	126.20	119.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	242	ARG	C-N-CA	6.29	126.20	119.85
1	C	69	GLY	CA-C-N	6.26	127.66	119.84
1	C	69	GLY	C-N-CA	6.26	127.66	119.84
1	D	458	HIS	N-CA-C	6.24	119.37	111.69
1	A	71	LEU	N-CA-C	-6.23	106.90	114.75
1	A	306	GLY	N-CA-C	6.16	127.79	113.18
1	A	307	MSE	N-CA-C	6.13	117.39	109.72
1	B	195	THR	N-CA-C	-6.07	105.78	113.43
1	C	309	MSE	CA-C-N	6.04	126.87	120.11
1	C	309	MSE	C-N-CA	6.04	126.87	120.11
1	C	4	GLU	N-CA-C	5.98	117.88	111.36
1	A	268	CYS	N-CA-C	5.96	118.61	108.90
1	C	216	ASP	N-CA-C	-5.87	103.82	111.74
1	D	217	LEU	N-CA-C	-5.81	106.53	113.97
1	D	283	GLN	N-CA-C	5.81	116.07	107.88
1	C	414	ILE	CA-C-N	5.80	126.61	120.11
1	C	414	ILE	C-N-CA	5.80	126.61	120.11
1	B	45	SER	N-CA-C	5.79	119.52	112.23
1	A	242	ARG	CA-C-N	5.77	125.68	119.85
1	A	242	ARG	C-N-CA	5.77	125.68	119.85
1	D	165	LYS	N-CA-C	5.74	117.21	111.07
1	C	540	TYR	N-CA-C	-5.72	106.94	114.04
1	B	489	ASN	N-CA-C	-5.71	106.44	113.41
1	A	165	LYS	N-CA-C	5.70	117.17	111.07
1	A	220	PHE	CA-C-N	5.70	127.08	120.98
1	A	220	PHE	C-N-CA	5.70	127.08	120.98
1	B	309	MSE	CA-C-N	5.64	126.43	120.11
1	B	309	MSE	C-N-CA	5.64	126.43	120.11
1	A	414	ILE	CA-C-N	5.63	126.42	120.11
1	A	414	ILE	C-N-CA	5.63	126.42	120.11
1	C	268	CYS	N-CA-C	5.62	118.06	108.90
1	A	529	PHE	CA-C-N	-5.60	112.84	119.84
1	A	529	PHE	C-N-CA	-5.60	112.84	119.84
1	C	34	ASN	N-CA-C	5.59	117.45	110.91
1	D	463	PHE	N-CA-C	-5.57	104.85	111.03
1	C	117	ARG	N-CA-C	5.54	117.76	111.11
1	B	7	CYS	N-CA-C	-5.52	105.66	112.90
1	B	448	ARG	N-CA-C	5.47	116.93	110.97
1	C	5	ILE	N-CA-C	5.45	115.65	110.53
1	D	385	LEU	N-CA-C	5.43	118.10	110.23
1	D	190	LEU	N-CA-C	5.42	117.92	108.76
1	D	16	ARG	N-CA-C	5.40	119.83	111.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	463	PHE	N-CA-C	-5.40	105.03	111.03
1	A	5	ILE	N-CA-C	5.39	115.60	110.53
1	D	307	MSE	CA-C-N	5.37	125.63	119.83
1	D	307	MSE	C-N-CA	5.37	125.63	119.83
1	B	50	ARG	N-CA-C	-5.36	105.48	112.23
1	A	124	VAL	CA-C-N	5.32	125.30	120.03
1	A	124	VAL	C-N-CA	5.32	125.30	120.03
1	B	268	CYS	N-CA-C	5.31	117.55	108.90
1	B	414	ILE	CA-C-N	5.27	126.01	120.11
1	B	414	ILE	C-N-CA	5.27	126.01	120.11
1	B	214	GLU	N-CA-C	-5.26	105.55	111.28
1	C	30	LEU	N-CA-C	5.24	119.80	113.41
1	B	439	LYS	N-CA-C	5.21	118.42	111.75
1	C	80	GLN	N-CA-C	-5.18	105.20	112.25
1	A	305	TYR	CA-C-O	-5.14	114.71	121.86
1	B	71	LEU	N-CA-C	-5.13	108.28	114.75
1	C	205	THR	N-CA-C	-5.13	105.86	111.82
1	D	240	ILE	N-CA-C	-5.10	105.74	110.53
1	C	253	VAL	N-CA-C	5.08	116.02	111.90
1	A	214	GLU	N-CA-C	-5.06	105.77	111.28
1	D	523	VAL	N-CA-C	5.05	115.39	108.12
1	D	400	PRO	CA-C-N	-5.04	114.79	120.14
1	D	400	PRO	C-N-CA	-5.04	114.79	120.14
1	A	150	LEU	N-CA-C	5.03	116.78	110.24
1	A	113	GLN	N-CA-C	-5.01	107.84	114.31

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	305	TYR	Mainchain
1	B	284	THR	Mainchain

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	4088	0	4108	209	1
1	B	4094	0	4116	202	1
1	C	4088	0	4108	215	0
1	D	4094	0	4116	214	0
2	E	50	0	46	1	0
2	F	46	0	43	2	0
2	G	46	0	43	4	0
2	H	53	0	51	5	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	31	0	12	2	0
4	B	31	0	12	2	0
4	C	31	0	12	1	0
4	D	31	0	12	3	0
All	All	16687	0	16679	831	1

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 25.

All (831) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:484:LEU:HD11	1:A:533:MSE:HG2	1.46	0.98
1:C:63:ARG:HE	1:C:240:ILE:HD11	1.31	0.93
1:D:160:ARG:HG3	1:D:160:ARG:HH11	1.34	0.91
1:C:160:ARG:HG3	1:C:160:ARG:HH11	1.35	0.91
1:D:364:GLU:HB2	1:D:373:SER:HB3	1.55	0.89
1:A:63:ARG:HE	1:A:240:ILE:HD11	1.38	0.88
1:C:47:MSE:HB2	1:C:53:ARG:HG3	1.57	0.87
1:A:302:LEU:HD13	1:A:319:LEU:HD11	1.59	0.82
1:A:47:MSE:HB2	1:A:53:ARG:HG3	1.60	0.82
1:B:20:ASP:OD2	1:B:82:HIS:NE2	2.12	0.82
1:D:160:ARG:NH1	1:D:442:ALA:O	2.15	0.79
1:D:150:LEU:O	1:D:267:ARG:NH1	2.18	0.77
1:C:364:GLU:HB2	1:C:373:SER:HB3	1.67	0.77
1:D:366:ILE:HD12	1:D:366:ILE:H	1.50	0.76
1:A:366:ILE:HD12	1:A:366:ILE:H	1.51	0.76
1:B:75:PHE:HB2	1:B:84:ALA:HB2	1.67	0.75
1:B:366:ILE:HD12	1:B:366:ILE:H	1.52	0.75
1:C:366:ILE:HD12	1:C:366:ILE:H	1.50	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:449:MSE:HE2	1:A:450:PRO:HA	1.69	0.74
1:C:75:PHE:HB2	1:C:84:ALA:HB2	1.70	0.74
1:A:364:GLU:HB2	1:A:373:SER:HB3	1.70	0.73
1:B:165:LYS:HG2	1:B:290:VAL:HG21	1.70	0.73
1:D:347:LYS:NZ	1:D:350:GLN:OE1	2.22	0.73
1:B:190:LEU:HD12	1:B:207:ILE:HG13	1.71	0.73
1:D:122:GLY:HA3	1:D:187:ILE:HG23	1.71	0.73
1:A:165:LYS:HG2	1:A:290:VAL:HG21	1.69	0.72
1:C:160:ARG:NH1	1:C:442:ALA:O	2.23	0.71
1:B:406:PRO:HA	1:B:453:THR:HG22	1.73	0.71
1:C:128:MSE:HE3	1:D:282:SER:HB2	1.72	0.71
1:C:90:TYR:HE2	1:C:106:ILE:HD11	1.55	0.71
1:D:443:LEU:HD23	1:D:460:ILE:HD11	1.71	0.71
1:A:8:ARG:HG2	1:A:8:ARG:HH11	1.55	0.70
1:B:302:LEU:HD13	1:B:319:LEU:HD11	1.72	0.70
1:A:75:PHE:HB2	1:A:84:ALA:HB2	1.73	0.70
1:B:135:TYR:O	1:B:139:ARG:HG3	1.90	0.70
1:B:102:ARG:HG3	1:B:103:PRO:HD2	1.73	0.69
1:C:406:PRO:HA	1:C:453:THR:HG22	1.74	0.69
1:C:128:MSE:HG3	1:C:167:VAL:HG22	1.73	0.69
1:A:386:SER:OG	1:A:388:GLU:OE1	2.10	0.69
1:A:96:ASN:O	1:A:98:PRO:HD3	1.91	0.69
1:B:478:SER:O	1:B:482:GLN:HG2	1.93	0.69
1:C:8:ARG:HG2	1:C:8:ARG:HH11	1.58	0.68
1:A:397:VAL:HG21	1:A:468:VAL:HG21	1.75	0.68
1:C:58:LEU:O	1:C:62:ARG:HG3	1.93	0.68
1:D:341:GLU:HB2	1:D:342:PRO:HD3	1.75	0.68
1:D:478:SER:O	1:D:482:GLN:HG2	1.93	0.68
2:H:750:PHE:HB3	2:H:753:LEU:HD21	1.73	0.68
1:C:96:ASN:O	1:C:98:PRO:HD3	1.94	0.67
1:D:255:GLU:O	1:D:259:ARG:HG3	1.94	0.67
1:B:230:VAL:HG12	1:B:234:MSE:HE2	1.76	0.67
1:C:195:THR:O	1:C:254:GLN:NE2	2.28	0.67
1:C:484:LEU:HD11	1:C:533:MSE:HG2	1.76	0.66
1:D:8:ARG:NH1	1:D:90:TYR:OH	2.27	0.66
1:D:230:VAL:HG12	1:D:234:MSE:HE2	1.78	0.66
1:C:67:GLU:HG2	1:C:69:GLY:H	1.59	0.66
1:B:216:ASP:OD1	1:B:219:ASN:ND2	2.26	0.66
1:C:255:GLU:O	1:C:259:ARG:HG3	1.97	0.65
1:D:165:LYS:HG2	1:D:290:VAL:HG21	1.78	0.65
1:A:327:SER:HB2	1:A:459:ILE:HG13	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:126:LYS:NZ	1:D:283:GLN:HG3	2.12	0.65
1:C:355:LEU:HD21	1:C:363:VAL:HB	1.77	0.65
1:B:92:ASP:OD1	1:B:96:ASN:ND2	2.29	0.64
1:D:463:PHE:O	1:D:467:VAL:HB	1.98	0.64
1:A:447:LYS:HE2	1:A:450:PRO:HD2	1.78	0.64
1:A:90:TYR:HE2	1:A:106:ILE:HD11	1.61	0.64
1:C:165:LYS:HG2	1:C:290:VAL:HG21	1.79	0.64
1:C:314:LYS:HA	1:C:317:ASP:HB2	1.79	0.64
1:C:397:VAL:HG21	1:C:468:VAL:HG21	1.80	0.64
1:A:214:GLU:HG3	1:B:237:ASN:CG	2.23	0.64
1:D:40:HIS:HA	1:D:43:LEU:HB2	1.80	0.64
1:C:141:ILE:HD12	1:C:175:LYS:HE2	1.80	0.63
1:C:309:MSE:HG3	1:C:310:PRO:HD2	1.81	0.63
1:C:335:MSE:HE3	1:C:364:GLU:HA	1.80	0.63
1:A:190:LEU:HD12	1:A:207:ILE:HG13	1.79	0.63
1:D:43:LEU:O	1:D:56:ASN:ND2	2.32	0.63
1:B:385:LEU:O	1:B:390:ARG:NH2	2.32	0.62
1:B:397:VAL:HG21	1:B:468:VAL:HG21	1.80	0.62
1:B:464:LEU:O	1:B:468:VAL:HG22	1.98	0.62
1:A:67:GLU:HG2	1:A:69:GLY:H	1.64	0.62
1:C:136:HIS:O	1:C:140:VAL:HG23	2.00	0.62
1:B:335:MSE:HE3	1:B:364:GLU:HA	1.81	0.62
1:A:128:MSE:HE3	1:B:282:SER:HB2	1.81	0.62
1:C:160:ARG:HB3	1:C:163:SER:HB3	1.81	0.62
1:D:128:MSE:HG3	1:D:167:VAL:HG22	1.81	0.62
1:A:200:THR:HG21	1:A:256:GLU:HB3	1.81	0.62
1:D:469:ASP:HB2	2:H:749:MSE:HA	1.80	0.62
1:D:49:THR:HG22	1:D:51:LEU:H	1.63	0.62
1:D:309:MSE:HG3	1:D:310:PRO:HD2	1.82	0.62
1:B:487:ILE:O	1:B:531:LYS:NZ	2.33	0.61
1:C:341:GLU:HB2	1:C:342:PRO:HD3	1.82	0.61
1:A:464:LEU:O	1:A:468:VAL:HG22	2.00	0.61
1:C:525:ARG:HD3	1:C:535:LEU:HD13	1.82	0.61
1:A:58:LEU:O	1:A:62:ARG:HG3	2.00	0.61
1:B:8:ARG:HH11	1:B:8:ARG:HG2	1.63	0.61
1:B:47:MSE:HB2	1:B:53:ARG:HG3	1.81	0.61
1:C:44:ILE:O	1:C:53:ARG:HG2	2.00	0.61
1:D:141:ILE:HD12	1:D:175:LYS:HE3	1.81	0.61
1:C:190:LEU:HD12	1:C:207:ILE:HG13	1.82	0.61
1:A:452:LEU:HD23	1:A:452:LEU:H	1.64	0.61
1:B:379:GLN:O	1:B:383:GLU:HG3	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:452:LEU:H	1:C:452:LEU:HD23	1.65	0.61
1:C:464:LEU:O	1:C:468:VAL:HG22	2.01	0.61
1:D:114:MSE:HG3	1:D:180:ILE:HB	1.83	0.61
1:A:92:ASP:OD1	1:A:96:ASN:ND2	2.29	0.61
1:D:397:VAL:HG21	1:D:468:VAL:HG21	1.82	0.60
1:B:299:TYR:HA	1:B:302:LEU:HD12	1.83	0.60
1:B:395:PHE:HD2	1:B:415:PRO:HD3	1.66	0.60
1:C:469:ASP:HB2	2:G:749:MSE:HA	1.83	0.60
1:C:327:SER:HB2	1:C:459:ILE:HG13	1.83	0.60
1:D:464:LEU:O	1:D:468:VAL:HG22	2.02	0.60
1:B:204:PHE:HA	1:B:207:ILE:HD12	1.83	0.60
1:B:341:GLU:HB2	1:B:342:PRO:HD3	1.83	0.60
1:C:160:ARG:HG3	1:C:160:ARG:NH1	2.13	0.60
1:A:128:MSE:HG3	1:A:167:VAL:HG22	1.83	0.60
1:A:410:TRP:HD1	1:A:414:ILE:HD13	1.66	0.60
1:B:47:MSE:HE3	1:B:52:GLU:HG2	1.84	0.60
1:D:160:ARG:HG3	1:D:160:ARG:NH1	2.10	0.60
1:A:341:GLU:HB2	1:A:342:PRO:HD3	1.84	0.60
1:C:178:GLN:HA	1:C:182:ILE:HB	1.83	0.60
1:D:386:SER:OG	1:D:388:GLU:OE1	2.20	0.60
1:C:447:LYS:HE2	1:C:450:PRO:HD2	1.84	0.59
1:A:36:PHE:HD2	1:A:40:HIS:HB3	1.66	0.59
1:A:309:MSE:HG3	1:A:310:PRO:HD2	1.84	0.59
1:D:452:LEU:HD23	1:D:452:LEU:H	1.67	0.59
1:B:108:PRO:O	1:B:109:GLN:HB2	2.01	0.59
1:C:388:GLU:O	1:C:391:SER:OG	2.17	0.59
1:C:443:LEU:HD23	1:C:460:ILE:HD11	1.83	0.59
1:D:133:ARG:NE	1:D:297:GLU:OE1	2.33	0.59
1:D:233:ARG:HG3	1:D:233:ARG:HH11	1.68	0.59
1:D:465:LYS:HD3	1:D:466:HIS:CE1	2.37	0.59
1:A:4:GLU:OE2	1:A:267:ARG:NH2	2.36	0.59
1:C:414:ILE:HG21	1:C:430:VAL:HG13	1.83	0.59
1:A:157:LEU:HD22	1:A:290:VAL:HG22	1.85	0.59
1:B:309:MSE:HG3	1:B:310:PRO:HD2	1.84	0.59
1:A:90:TYR:CE2	1:A:106:ILE:HD11	2.38	0.59
1:A:385:LEU:O	1:A:390:ARG:NH2	2.36	0.59
1:B:102:ARG:HG3	1:B:103:PRO:CD	2.33	0.59
1:A:335:MSE:HE3	1:A:364:GLU:HA	1.83	0.59
1:C:137:VAL:O	1:C:141:ILE:HG13	2.02	0.59
1:C:150:LEU:O	1:C:267:ARG:NH1	2.34	0.59
1:A:457:ASP:OD1	1:A:457:ASP:N	2.35	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:342:PRO:O	1:D:343:LYS:HB3	2.02	0.58
1:C:216:ASP:HA	1:C:219:ASN:HB2	1.85	0.58
1:A:395:PHE:HD2	1:A:415:PRO:HD3	1.68	0.58
1:B:23:PRO:HG2	1:B:53:ARG:C	2.29	0.58
1:B:111:SER:O	1:B:115:LEU:HB3	2.03	0.58
1:A:342:PRO:O	1:A:343:LYS:HB3	2.04	0.58
1:B:437:LEU:HB3	1:B:444:LEU:HD22	1.86	0.58
1:C:465:LYS:HD3	1:C:466:HIS:CE1	2.38	0.58
1:A:521:GLU:C	1:A:523:VAL:H	2.12	0.58
1:B:30:LEU:HB3	1:B:36:PHE:CD1	2.39	0.58
1:C:397:VAL:HG22	1:C:464:LEU:HB3	1.84	0.58
1:C:465:LYS:HD3	1:C:466:HIS:HE1	1.69	0.58
1:D:385:LEU:O	1:D:390:ARG:NH2	2.37	0.58
1:A:72:ILE:HG21	1:A:88:GLU:HG3	1.85	0.57
1:A:437:LEU:HB3	1:A:444:LEU:HD22	1.84	0.57
1:B:47:MSE:HE2	1:B:56:ASN:HD21	1.69	0.57
1:B:443:LEU:HD23	1:B:460:ILE:HD11	1.86	0.57
2:G:750:PHE:O	2:G:752:PHE:N	2.37	0.57
1:B:109:GLN:C	1:B:111:SER:H	2.12	0.57
1:A:414:ILE:HG21	1:A:430:VAL:HG13	1.86	0.57
1:D:24:ARG:HG3	1:D:27:LEU:HD12	1.85	0.57
2:E:750:PHE:O	2:E:752:PHE:N	2.37	0.57
1:C:395:PHE:HD2	1:C:415:PRO:HD3	1.69	0.57
1:D:436:ARG:O	1:D:440:ARG:HG3	2.04	0.57
1:A:379:GLN:O	1:A:383:GLU:HG3	2.05	0.57
1:B:128:MSE:HG3	1:B:167:VAL:HG22	1.85	0.57
1:C:527:GLU:HA	1:C:530:PRO:HG3	1.87	0.57
1:D:250:ASP:HA	1:D:271:THR:OG1	2.05	0.57
1:A:63:ARG:HE	1:A:240:ILE:CD1	2.13	0.57
1:D:335:MSE:HE3	1:D:364:GLU:HA	1.86	0.56
1:A:150:LEU:O	1:A:267:ARG:NH1	2.39	0.56
1:A:247:PHE:HE2	1:A:266:LEU:HD22	1.69	0.56
1:B:4:GLU:HG3	1:B:5:ILE:N	2.19	0.56
1:A:160:ARG:HG2	1:A:161:ALA:H	1.69	0.56
1:B:302:LEU:HD23	1:B:337:PHE:CE1	2.41	0.56
1:B:342:PRO:O	1:B:343:LYS:HB3	2.04	0.56
1:D:283:GLN:HG2	1:D:284:THR:H	1.70	0.56
1:D:354:LYS:HD3	1:D:362:GLY:O	2.04	0.56
1:D:111:SER:O	1:D:115:LEU:HB3	2.05	0.56
1:C:159:GLY:N	1:C:165:LYS:HD3	2.21	0.56
1:A:406:PRO:HA	1:A:453:THR:HG22	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:532:PHE:C	1:D:534:GLN:H	2.13	0.56
1:A:91:ILE:O	1:A:95:ILE:HG13	2.05	0.56
1:A:197:PRO:HD2	1:A:198:LYS:HE2	1.88	0.56
1:B:529:PHE:HB3	1:B:532:PHE:CZ	2.40	0.56
1:C:233:ARG:HH11	1:C:233:ARG:HG3	1.70	0.56
1:A:79:ASN:ND2	1:B:36:PHE:O	2.38	0.56
1:D:237:ASN:C	1:D:239:LEU:H	2.14	0.56
1:D:406:PRO:HA	1:D:453:THR:HG22	1.86	0.56
1:A:233:ARG:HG3	1:A:233:ARG:HH11	1.71	0.56
1:A:438:SER:O	1:A:442:ALA:HA	2.05	0.56
1:C:126:LYS:HZ3	1:D:283:GLN:HG3	1.70	0.56
1:B:452:LEU:HD23	1:B:452:LEU:H	1.71	0.55
1:C:326:SER:O	1:C:328:GLY:N	2.39	0.55
1:A:465:LYS:HD3	1:A:466:HIS:CE1	2.41	0.55
1:B:1:MSE:HA	1:B:62:ARG:O	2.06	0.55
1:D:155:LEU:HB3	1:D:269:LEU:HD23	1.88	0.55
1:A:415:PRO:O	1:A:416:VAL:HG23	2.06	0.55
1:B:274:ASP:OD2	1:B:440:ARG:NH1	2.40	0.55
2:F:750:PHE:HB3	2:F:753:LEU:HA	1.88	0.55
1:A:111:SER:C	1:A:113:GLN:H	2.14	0.55
1:A:354:LYS:HD3	1:A:362:GLY:O	2.07	0.55
1:B:410:TRP:HD1	1:B:414:ILE:HD13	1.72	0.55
1:C:273:ARG:NH2	1:C:380:ARG:HH21	2.05	0.55
1:D:255:GLU:HB3	1:D:277:ILE:HG22	1.88	0.55
1:D:457:ASP:OD1	1:D:457:ASP:N	2.39	0.55
1:A:250:ASP:HA	1:A:271:THR:OG1	2.06	0.55
1:C:532:PHE:HB2	1:C:534:GLN:HG2	1.88	0.55
1:D:160:ARG:HB3	1:D:163:SER:HB3	1.86	0.55
1:A:397:VAL:O	1:A:461:HIS:NE2	2.37	0.55
1:C:342:PRO:O	1:C:343:LYS:HB3	2.07	0.55
1:C:457:ASP:OD1	1:C:457:ASP:N	2.40	0.54
1:A:257:THR:HG22	1:A:258:ILE:N	2.22	0.54
1:B:529:PHE:N	1:B:530:PRO:HD3	2.22	0.54
1:D:397:VAL:C	1:D:399:MSE:H	2.15	0.54
1:D:487:ILE:O	1:D:531:LYS:NZ	2.40	0.54
1:D:537:GLN:O	1:D:541:ASP:N	2.38	0.54
1:A:178:GLN:O	1:A:182:ILE:HB	2.07	0.54
1:C:36:PHE:HD2	1:C:40:HIS:HB3	1.72	0.54
1:C:250:ASP:HA	1:C:271:THR:OG1	2.08	0.54
1:D:42:GLU:O	1:D:46:LYS:NZ	2.26	0.54
1:B:72:ILE:HG21	1:B:88:GLU:HG3	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:113:GLN:O	1:D:118:LYS:HG3	2.08	0.54
1:D:347:LYS:O	1:D:350:GLN:HB2	2.06	0.54
1:D:405:ILE:HG13	1:D:409:LEU:HB3	1.90	0.54
1:A:157:LEU:HD11	1:A:168:ILE:HG22	1.89	0.54
1:A:397:VAL:HG22	1:A:464:LEU:HB3	1.88	0.54
1:B:250:ASP:HA	1:B:271:THR:OG1	2.07	0.54
1:C:407:VAL:HG23	1:C:452:LEU:O	2.07	0.54
1:D:310:PRO:HA	1:D:345:PHE:HE2	1.72	0.54
1:D:326:SER:O	1:D:328:GLY:N	2.40	0.54
1:B:233:ARG:HG3	1:B:233:ARG:HH11	1.73	0.54
1:B:458:HIS:HB3	1:B:462:MSE:HE2	1.90	0.54
1:C:463:PHE:O	1:C:467:VAL:HB	2.08	0.54
1:A:44:ILE:O	1:A:53:ARG:HG2	2.08	0.53
1:A:389:ASP:HB3	1:A:437:LEU:HD21	1.89	0.53
1:B:415:PRO:O	1:B:416:VAL:HG23	2.08	0.53
1:B:457:ASP:OD1	1:B:457:ASP:N	2.40	0.53
1:B:521:GLU:C	1:B:523:VAL:H	2.16	0.53
1:C:90:TYR:CE2	1:C:106:ILE:HD11	2.40	0.53
1:B:112:ARG:HA	1:B:116:ASP:HB2	1.90	0.53
1:C:163:SER:C	1:C:330:PRO:HG2	2.34	0.53
1:C:354:LYS:O	1:C:358:ARG:HB2	2.09	0.53
1:D:136:HIS:O	1:D:140:VAL:HG23	2.08	0.53
1:A:436:ARG:O	1:A:440:ARG:HG3	2.09	0.53
1:D:397:VAL:HG22	1:D:464:LEU:HB3	1.90	0.53
1:D:529:PHE:N	1:D:530:PRO:HD3	2.24	0.53
1:A:36:PHE:CD2	1:A:40:HIS:HB3	2.44	0.53
1:C:8:ARG:NH1	1:C:185:ASP:OD1	2.42	0.53
1:C:67:GLU:HG2	1:C:69:GLY:N	2.23	0.53
1:D:135:TYR:O	1:D:139:ARG:HG3	2.08	0.53
1:D:247:PHE:HE2	1:D:266:LEU:HD22	1.74	0.53
1:B:157:LEU:HD22	1:B:290:VAL:HG22	1.91	0.53
1:D:341:GLU:CB	1:D:342:PRO:HD3	2.39	0.53
1:A:189:TRP:CE3	1:A:248:VAL:HG11	2.44	0.52
1:B:155:LEU:HB3	1:B:269:LEU:HD23	1.91	0.52
1:B:257:THR:HG22	1:B:258:ILE:N	2.24	0.52
1:B:386:SER:OG	1:B:388:GLU:OE1	2.27	0.52
1:B:35:ILE:HD12	1:B:70:PRO:HG2	1.89	0.52
1:A:93:PHE:HE2	1:A:104:VAL:HG11	1.75	0.52
1:A:443:LEU:HD23	1:A:460:ILE:HD11	1.91	0.52
1:B:364:GLU:HB2	1:B:373:SER:HB3	1.90	0.52
1:B:527:GLU:N	1:B:527:GLU:OE1	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:122:GLY:HA3	1:C:187:ILE:HG23	1.91	0.52
1:C:436:ARG:O	1:C:440:ARG:HG3	2.08	0.52
1:A:133:ARG:O	1:A:137:VAL:HG23	2.09	0.52
1:A:458:HIS:HB3	1:A:462:MSE:HE2	1.92	0.52
1:B:157:LEU:HD11	1:B:168:ILE:CG2	2.39	0.52
1:A:165:LYS:N	4:A:602:ATP:O2B	2.42	0.52
1:A:255:GLU:O	1:A:259:ARG:HG3	2.10	0.52
1:C:26:ALA:HB2	1:C:74:PHE:CE1	2.45	0.52
1:C:102:ARG:HG2	1:C:182:ILE:HG22	1.92	0.52
1:C:216:ASP:H	1:C:218:LEU:HD23	1.74	0.52
1:D:1:MSE:O	1:D:2:LEU:HD23	2.10	0.52
1:B:231:LEU:HA	1:B:234:MSE:HE3	1.91	0.52
1:B:388:GLU:HB3	1:B:433:ARG:HH21	1.74	0.52
1:B:438:SER:O	1:B:442:ALA:HA	2.10	0.52
2:H:749:MSE:O	2:H:750:PHE:HB2	2.10	0.52
1:A:67:GLU:HG2	1:A:69:GLY:N	2.24	0.51
1:A:143:LYS:O	1:A:147:MSE:HG3	2.10	0.51
1:B:255:GLU:O	1:B:259:ARG:HG3	2.10	0.51
1:D:465:LYS:HD3	1:D:466:HIS:HE1	1.73	0.51
1:A:532:PHE:HB2	1:A:534:GLN:HG2	1.91	0.51
1:B:50:ARG:HA	1:B:53:ARG:HE	1.76	0.51
1:C:18:ILE:O	1:C:50:ARG:NH2	2.42	0.51
1:C:283:GLN:HG2	1:C:284:THR:H	1.74	0.51
1:C:379:GLN:O	1:C:383:GLU:HG3	2.10	0.51
1:C:23:PRO:HG2	1:C:53:ARG:HB3	1.93	0.51
1:C:247:PHE:HE2	1:C:266:LEU:HD22	1.75	0.51
1:D:197:PRO:HD2	1:D:198:LYS:HE2	1.92	0.51
1:A:247:PHE:CE2	1:A:266:LEU:HD13	2.46	0.51
1:B:157:LEU:HD11	1:B:168:ILE:HG22	1.92	0.51
1:B:336:PHE:HE2	1:B:348:MSE:SE	2.44	0.51
1:C:72:ILE:HG13	1:C:88:GLU:HG3	1.91	0.51
1:C:313:GLU:O	1:C:316:GLU:HG3	2.09	0.51
1:D:437:LEU:HB3	1:D:444:LEU:HD22	1.93	0.51
1:B:158:HIS:CE1	1:B:289:GLU:HB2	2.45	0.51
1:B:397:VAL:HG22	1:B:464:LEU:HB3	1.93	0.51
1:C:9:ALA:HB2	1:C:90:TYR:CD1	2.46	0.51
1:D:388:GLU:HB3	1:D:433:ARG:HH21	1.74	0.51
1:B:47:MSE:HE2	1:B:56:ASN:ND2	2.26	0.51
1:B:197:PRO:HD2	1:B:198:LYS:HE2	1.92	0.51
1:D:204:PHE:HA	1:D:207:ILE:HD12	1.92	0.51
1:D:388:GLU:O	1:D:391:SER:OG	2.20	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:HIS:ND1	1:A:60:ILE:HD13	2.26	0.51
1:B:4:GLU:O	1:B:8:ARG:HB2	2.11	0.51
1:B:160:ARG:NH1	1:B:442:ALA:O	2.44	0.51
1:D:75:PHE:HB3	1:D:84:ALA:HB2	1.92	0.51
1:B:170:SER:O	1:B:174:SER:HB2	2.11	0.51
1:C:102:ARG:HG2	1:C:182:ILE:CG2	2.40	0.51
1:C:111:SER:C	1:C:113:GLN:H	2.19	0.51
1:C:527:GLU:C	1:C:530:PRO:HD3	2.36	0.51
1:D:257:THR:HG22	1:D:258:ILE:N	2.26	0.51
1:D:415:PRO:O	1:D:416:VAL:HG23	2.11	0.51
1:B:414:ILE:HG21	1:B:430:VAL:HG13	1.93	0.51
1:C:125:PRO:O	1:C:170:SER:OG	2.29	0.51
1:B:189:TRP:CE3	1:B:248:VAL:HG11	2.46	0.50
1:C:158:HIS:CE1	1:C:289:GLU:HB2	2.46	0.50
1:C:188:VAL:HB	1:C:247:PHE:CD1	2.47	0.50
1:C:347:LYS:HA	1:C:350:GLN:HB2	1.93	0.50
1:D:200:THR:HG21	1:D:256:GLU:HB3	1.93	0.50
1:A:47:MSE:HB2	1:A:53:ARG:CG	2.37	0.50
1:A:136:HIS:O	1:A:140:VAL:HG23	2.11	0.50
1:B:47:MSE:HB3	1:B:52:GLU:HB3	1.93	0.50
2:F:751:ASN:ND2	2:F:751:ASN:O	2.44	0.50
1:D:159:GLY:N	1:D:165:LYS:HD3	2.26	0.50
1:A:111:SER:O	1:A:113:GLN:N	2.44	0.50
1:A:216:ASP:H	1:A:218:LEU:HD23	1.76	0.50
1:B:32:GLY:C	1:B:34:ASN:H	2.19	0.50
1:B:160:ARG:HG2	1:B:161:ALA:H	1.76	0.50
1:B:189:TRP:CD1	1:B:189:TRP:C	2.90	0.50
1:B:463:PHE:O	1:B:467:VAL:HB	2.12	0.50
1:D:314:LYS:HA	1:D:317:ASP:HB2	1.94	0.50
1:D:355:LEU:HD21	1:D:363:VAL:HB	1.92	0.50
1:D:397:VAL:O	1:D:399:MSE:N	2.44	0.50
1:C:165:LYS:N	4:C:602:ATP:O2B	2.44	0.50
1:C:257:THR:HG22	1:C:258:ILE:N	2.26	0.50
1:A:188:VAL:HB	1:A:247:PHE:CD1	2.46	0.50
1:B:49:THR:HG22	1:B:51:LEU:H	1.77	0.50
1:B:109:GLN:C	1:B:111:SER:N	2.68	0.50
1:C:142:LYS:O	1:C:146:GLU:HG3	2.12	0.50
1:D:164:GLY:HA3	4:D:602:ATP:H2	1.75	0.50
1:A:135:TYR:O	1:A:139:ARG:HG3	2.12	0.50
1:C:189:TRP:CE3	1:C:248:VAL:HG11	2.46	0.50
1:A:410:TRP:CD1	1:A:414:ILE:HD13	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:219:ASN:O	1:C:221:PRO:HD3	2.12	0.49
1:A:465:LYS:HD3	1:A:466:HIS:HE1	1.77	0.49
1:B:136:HIS:O	1:B:140:VAL:HG23	2.12	0.49
1:C:150:LEU:HD11	1:C:284:THR:HG21	1.93	0.49
1:C:397:VAL:C	1:C:399:MSE:H	2.19	0.49
1:C:529:PHE:N	1:C:530:PRO:HD3	2.27	0.49
1:D:235:ILE:O	1:D:239:LEU:HB2	2.11	0.49
1:A:190:LEU:CD1	1:A:207:ILE:HG13	2.42	0.49
1:B:302:LEU:HD23	1:B:337:PHE:HE1	1.77	0.49
1:D:178:GLN:HB3	1:D:183:ASN:OD1	2.13	0.49
1:B:158:HIS:CD2	1:B:275:VAL:HG13	2.47	0.49
1:B:158:HIS:HE1	1:B:289:GLU:HB2	1.77	0.49
1:B:332:THR:O	1:B:336:PHE:HB2	2.12	0.49
1:C:132:ILE:HG23	1:C:137:VAL:HG21	1.93	0.49
1:B:195:THR:O	1:B:254:GLN:NE2	2.46	0.49
1:B:410:TRP:CD1	1:B:414:ILE:HD13	2.47	0.49
1:B:56:ASN:O	1:B:60:ILE:HB	2.12	0.49
1:A:299:TYR:HA	1:A:302:LEU:HD12	1.95	0.49
1:D:178:GLN:HA	1:D:182:ILE:HB	1.94	0.49
1:A:163:SER:C	1:A:330:PRO:HG2	2.38	0.49
1:B:332:THR:HG23	1:B:374:LEU:HB2	1.95	0.49
1:B:355:LEU:HD21	1:B:363:VAL:HB	1.95	0.49
1:C:524:ILE:HG13	1:C:525:ARG:N	2.27	0.49
1:B:114:MSE:SE	1:B:177:ASP:HA	2.62	0.49
1:D:157:LEU:HD11	1:D:168:ILE:HG22	1.95	0.49
1:A:166:SER:HA	1:A:271:THR:HG21	1.94	0.49
1:C:21:PHE:CD1	1:C:75:PHE:HE1	2.30	0.49
1:B:252:VAL:HG11	1:B:258:ILE:HD13	1.94	0.48
1:C:343:LYS:O	1:C:343:LYS:HG2	2.14	0.48
1:D:378:LEU:HA	1:D:381:CYS:HB2	1.94	0.48
1:D:527:GLU:N	1:D:527:GLU:OE1	2.45	0.48
1:A:72:ILE:C	1:A:74:PHE:H	2.21	0.48
1:A:248:VAL:HG22	1:A:269:LEU:HB3	1.95	0.48
1:C:40:HIS:O	1:C:44:ILE:HG13	2.13	0.48
1:D:135:TYR:CZ	1:D:139:ARG:HD2	2.48	0.48
1:B:314:LYS:HA	1:B:317:ASP:HB2	1.95	0.48
1:C:527:GLU:N	1:C:527:GLU:OE1	2.46	0.48
1:B:23:PRO:HG2	1:B:53:ARG:O	2.13	0.48
1:B:307:MSE:SE	1:B:308:PRO:HD2	2.64	0.48
1:D:327:SER:HB2	1:D:459:ILE:HD11	1.94	0.48
1:A:206:ASP:O	1:A:210:MSE:HG2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:393:LEU:HB2	1:A:437:LEU:HD11	1.94	0.48
1:C:155:LEU:HB3	1:C:269:LEU:HD23	1.96	0.48
1:C:525:ARG:HA	1:C:526:PRO:HD3	1.64	0.48
1:D:318:VAL:O	1:D:321:LYS:HB3	2.13	0.48
1:A:137:VAL:O	1:A:141:ILE:HG13	2.14	0.48
1:A:142:LYS:HE2	1:A:146:GLU:OE2	2.14	0.48
1:A:170:SER:O	1:A:174:SER:HB2	2.13	0.48
1:A:302:LEU:HD23	1:A:337:PHE:CE1	2.49	0.48
1:C:138:ASP:OD1	1:C:175:LYS:NZ	2.38	0.48
1:D:189:TRP:CE3	1:D:248:VAL:HG11	2.47	0.48
1:D:527:GLU:C	1:D:530:PRO:HD3	2.38	0.48
1:A:8:ARG:HG2	1:A:8:ARG:NH1	2.28	0.48
1:B:341:GLU:CB	1:B:342:PRO:HD3	2.44	0.48
1:D:283:GLN:C	1:D:285:CYS:H	2.22	0.48
1:D:399:MSE:HE3	1:D:410:TRP:CD1	2.49	0.48
1:A:92:ASP:C	1:A:94:ALA:H	2.22	0.48
1:C:47:MSE:HB2	1:C:53:ARG:CG	2.36	0.48
1:D:190:LEU:CD1	1:D:207:ILE:HG13	2.43	0.48
1:D:237:ASN:C	1:D:239:LEU:N	2.72	0.48
1:B:72:ILE:C	1:B:74:PHE:H	2.21	0.48
1:B:166:SER:HA	1:B:271:THR:HG21	1.96	0.48
1:B:532:PHE:C	1:B:534:GLN:H	2.22	0.48
1:C:44:ILE:HD11	1:C:60:ILE:HD12	1.96	0.48
1:C:269:LEU:HD22	1:C:270:VAL:H	1.79	0.48
1:C:334:MSE:HE3	1:C:334:MSE:HA	1.95	0.48
1:D:334:MSE:HE3	1:D:334:MSE:HA	1.96	0.48
1:A:252:VAL:CG2	1:A:257:THR:HG21	2.44	0.48
1:B:154:PHE:HE2	1:B:262:GLN:HG3	1.78	0.48
1:D:44:ILE:HG23	1:D:56:ASN:HB3	1.94	0.48
1:D:319:LEU:HD12	1:D:345:PHE:CE1	2.48	0.48
1:A:160:ARG:NH1	1:A:442:ALA:O	2.46	0.47
1:C:240:ILE:HG13	1:C:241:ASP:OD1	2.14	0.47
1:C:315:GLU:HA	1:C:318:VAL:HG23	1.96	0.47
1:A:181:GLY:N	1:A:184:TYR:O	2.31	0.47
1:B:114:MSE:HG3	1:B:180:ILE:HB	1.96	0.47
1:C:72:ILE:HG21	1:C:88:GLU:CD	2.39	0.47
1:C:206:ASP:O	1:C:210:MSE:HG2	2.14	0.47
1:D:110:PHE:HD1	1:D:110:PHE:HA	1.60	0.47
1:A:47:MSE:HE2	1:A:52:GLU:HG2	1.96	0.47
1:D:233:ARG:HH11	1:D:233:ARG:CG	2.27	0.47
1:D:329:ASN:HA	1:D:330:PRO:HD2	1.75	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:ILE:HD13	1:A:246:LEU:HB3	1.96	0.47
1:A:195:THR:O	1:A:254:GLN:NE2	2.47	0.47
1:D:233:ARG:HA	1:D:236:CYS:SG	2.55	0.47
1:D:414:ILE:HG21	1:D:430:VAL:HG13	1.96	0.47
1:A:135:TYR:CZ	1:A:139:ARG:HD2	2.49	0.47
1:B:114:MSE:O	1:B:119:LEU:HG	2.14	0.47
1:C:30:LEU:HD21	1:C:70:PRO:O	2.14	0.47
1:C:200:THR:HG21	1:C:256:GLU:HB3	1.96	0.47
1:D:154:PHE:HE2	1:D:262:GLN:HG3	1.79	0.47
1:D:259:ARG:O	1:D:263:GLU:HG3	2.15	0.47
1:A:233:ARG:HH11	1:A:233:ARG:CG	2.28	0.47
1:A:255:GLU:HB3	1:A:277:ILE:HG22	1.96	0.47
1:A:532:PHE:C	1:A:534:GLN:H	2.22	0.47
1:B:88:GLU:O	1:B:91:ILE:HG22	2.15	0.47
1:B:88:GLU:HA	1:B:91:ILE:HG22	1.96	0.47
1:B:255:GLU:HB3	1:B:277:ILE:HG22	1.96	0.47
1:B:336:PHE:CD1	1:B:363:VAL:HG21	2.50	0.47
1:C:126:LYS:HD3	1:D:283:GLN:HE21	1.79	0.47
1:C:158:HIS:HE1	1:C:289:GLU:HB2	1.80	0.47
1:C:444:LEU:HD12	1:C:444:LEU:HA	1.74	0.47
1:C:532:PHE:C	1:C:534:GLN:H	2.21	0.47
1:D:380:ARG:HA	1:D:383:GLU:HG3	1.97	0.47
1:A:44:ILE:HD11	1:A:60:ILE:HD12	1.97	0.47
1:B:133:ARG:O	1:B:137:VAL:HG23	2.15	0.47
1:C:189:TRP:CD1	1:C:189:TRP:C	2.93	0.47
1:C:407:VAL:HG21	1:C:452:LEU:HD12	1.96	0.47
1:A:7:CYS:SG	1:A:62:ARG:HD3	2.55	0.47
1:A:401:PRO:HB2	1:A:458:HIS:CE1	2.50	0.47
1:A:527:GLU:C	1:A:530:PRO:HD3	2.39	0.47
1:A:326:SER:O	1:A:328:GLY:N	2.48	0.47
1:B:163:SER:C	1:B:330:PRO:HG2	2.40	0.47
1:C:139:ARG:C	1:C:139:ARG:HD3	2.40	0.47
1:C:235:ILE:O	1:C:239:LEU:HB2	2.14	0.47
1:C:410:TRP:HD1	1:C:414:ILE:HD13	1.80	0.47
1:A:269:LEU:HD22	1:A:270:VAL:H	1.80	0.46
1:A:283:GLN:HG2	1:A:284:THR:H	1.79	0.46
1:C:88:GLU:HA	1:C:91:ILE:HG22	1.97	0.46
1:D:307:MSE:SE	1:D:308:PRO:HD2	2.65	0.46
1:A:2:LEU:HG	1:A:62:ARG:HA	1.96	0.46
1:B:237:ASN:C	1:B:239:LEU:H	2.22	0.46
1:B:257:THR:HG22	1:B:258:ILE:HD12	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:283:GLN:O	1:B:285:CYS:N	2.48	0.46
1:B:329:ASN:HA	1:B:330:PRO:HD2	1.76	0.46
1:B:354:LYS:HD3	1:B:362:GLY:O	2.15	0.46
1:D:472:THR:HG21	2:H:749:MSE:HE2	1.96	0.46
1:A:527:GLU:OE1	1:A:527:GLU:N	2.47	0.46
1:A:252:VAL:HG21	1:A:257:THR:HG21	1.98	0.46
1:C:393:LEU:HB2	1:C:437:LEU:HD11	1.97	0.46
1:D:44:ILE:HD11	1:D:60:ILE:HD12	1.96	0.46
1:D:248:VAL:HG22	1:D:269:LEU:HD12	1.98	0.46
1:A:237:ASN:C	1:A:239:LEU:H	2.23	0.46
1:C:163:SER:O	1:C:330:PRO:HG2	2.15	0.46
1:C:443:LEU:HA	1:C:460:ILE:HG12	1.96	0.46
1:D:137:VAL:O	1:D:141:ILE:HG13	2.16	0.46
1:D:337:PHE:CD1	1:D:337:PHE:C	2.94	0.46
1:A:385:LEU:HD23	1:A:440:ARG:HH21	1.81	0.46
1:B:406:PRO:CA	1:B:453:THR:HG22	2.44	0.46
1:B:481:GLU:HA	1:B:484:LEU:CD2	2.46	0.46
1:D:131:TYR:HE2	1:D:300:ASP:HB2	1.81	0.46
1:D:343:LYS:O	1:D:343:LYS:HG2	2.16	0.46
1:D:371:TYR:CD2	1:D:377:ALA:HB2	2.51	0.46
1:A:347:LYS:HA	1:A:350:GLN:HB2	1.97	0.46
1:B:242:ARG:HB3	1:B:245:THR:OG1	2.16	0.46
1:A:30:LEU:HD21	1:A:70:PRO:O	2.15	0.46
1:A:302:LEU:CD1	1:A:319:LEU:HD11	2.40	0.46
1:B:39:ASP:HA	1:B:42:GLU:HB2	1.97	0.46
1:C:130:CYS:O	1:C:304:ALA:HB1	2.15	0.46
1:C:368:PRO:HD3	1:D:279:ASN:O	2.16	0.46
1:D:247:PHE:CE2	1:D:266:LEU:HD13	2.51	0.46
1:A:404:ASP:OD1	1:A:447:LYS:HD3	2.15	0.46
1:A:471:GLN:HE21	1:D:471:GLN:HG3	1.80	0.46
1:B:252:VAL:HG22	1:B:257:THR:HG21	1.98	0.46
1:A:471:GLN:HG3	1:D:471:GLN:NE2	2.31	0.46
1:B:465:LYS:HD3	1:B:466:HIS:CE1	2.51	0.46
1:C:178:GLN:HB3	1:C:183:ASN:OD1	2.16	0.46
1:C:237:ASN:O	1:C:240:ILE:HG23	2.16	0.46
1:D:409:LEU:HD22	1:D:409:LEU:HA	1.77	0.46
1:A:347:LYS:O	1:A:347:LYS:HD3	2.15	0.45
1:B:408:LYS:HD2	1:B:408:LYS:HA	1.73	0.45
1:D:337:PHE:C	1:D:337:PHE:HD1	2.23	0.45
1:A:189:TRP:CD1	1:A:189:TRP:C	2.94	0.45
1:C:109:GLN:HE21	1:C:109:GLN:HB2	1.53	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:142:LYS:O	1:D:146:GLU:HG3	2.17	0.45
1:A:78:ASN:HA	1:B:37:THR:HG21	1.98	0.45
1:A:106:ILE:HG21	1:A:182:ILE:HA	1.99	0.45
1:B:253:VAL:HG11	1:B:380:ARG:HD3	1.98	0.45
1:C:23:PRO:O	1:C:27:LEU:HG	2.17	0.45
1:C:111:SER:O	1:C:113:GLN:N	2.50	0.45
1:C:329:ASN:HA	1:C:330:PRO:HD2	1.82	0.45
1:C:386:SER:OG	1:C:388:GLU:OE1	2.34	0.45
1:D:28:THR:O	1:D:32:GLY:N	2.49	0.45
1:D:36:PHE:HE2	1:D:44:ILE:HD12	1.82	0.45
1:D:110:PHE:C	1:D:112:ARG:H	2.23	0.45
1:D:491:ASN:OD1	1:D:491:ASN:N	2.50	0.45
1:A:395:PHE:O	1:A:398:VAL:HG22	2.17	0.45
1:B:399:MSE:HE1	1:B:409:LEU:O	2.15	0.45
1:D:354:LYS:O	1:D:358:ARG:HB2	2.17	0.45
1:A:21:PHE:CG	1:A:22:GLU:N	2.84	0.45
1:B:35:ILE:CD1	1:B:70:PRO:HG2	2.47	0.45
1:B:137:VAL:O	1:B:141:ILE:HG13	2.17	0.45
1:B:327:SER:HB2	1:B:459:ILE:HG13	1.97	0.45
1:D:61:TYR:HE2	1:D:68:LEU:HD22	1.81	0.45
1:A:37:THR:H	1:A:40:HIS:HB2	1.81	0.45
1:A:156:PHE:HB2	1:A:287:PHE:CD2	2.52	0.45
1:C:248:VAL:HG22	1:C:269:LEU:HD12	1.98	0.45
1:A:18:ILE:O	1:A:50:ARG:NH2	2.49	0.45
1:D:32:GLY:C	1:D:34:ASN:H	2.24	0.45
1:D:100:LEU:HD12	1:D:100:LEU:HA	1.77	0.45
1:D:403:VAL:HG12	1:D:405:ILE:HG22	1.99	0.45
1:A:351:LEU:HD13	1:A:363:VAL:HG23	1.99	0.45
1:B:105:VAL:HG22	1:B:110:PHE:CE2	2.52	0.45
1:C:126:LYS:HE2	1:C:126:LYS:O	2.18	0.45
1:C:233:ARG:NH1	1:C:237:ASN:OD1	2.49	0.45
1:C:237:ASN:C	1:C:239:LEU:H	2.24	0.45
1:A:405:ILE:HA	1:A:406:PRO:HD3	1.71	0.44
1:B:23:PRO:O	1:B:27:LEU:HG	2.17	0.44
1:B:218:LEU:HG	1:B:219:ASN:N	2.33	0.44
1:C:167:VAL:O	1:C:171:GLN:HG3	2.17	0.44
1:C:240:ILE:C	1:C:242:ARG:H	2.25	0.44
1:A:116:ASP:OD1	1:A:116:ASP:N	2.49	0.44
1:A:368:PRO:HB2	4:A:602:ATP:O2'	2.17	0.44
1:D:39:ASP:OD1	1:D:39:ASP:N	2.47	0.44
1:D:124:VAL:HA	1:D:125:PRO:HD3	1.85	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:158:HIS:HE1	1:D:287:PHE:HB3	1.82	0.44
1:A:341:GLU:CB	1:A:342:PRO:HD3	2.47	0.44
1:B:49:THR:HG22	1:B:50:ARG:N	2.31	0.44
1:B:477:ILE:HD11	1:B:543:LEU:HG	2.00	0.44
1:C:116:ASP:OD1	1:C:116:ASP:N	2.51	0.44
1:C:388:GLU:CD	1:C:388:GLU:H	2.26	0.44
1:D:405:ILE:HA	1:D:406:PRO:HD3	1.77	0.44
1:D:444:LEU:HB3	1:D:445:SER:H	1.54	0.44
1:C:233:ARG:HH11	1:C:233:ARG:CG	2.29	0.44
1:C:247:PHE:CE2	1:C:266:LEU:HD13	2.53	0.44
1:D:17:LEU:O	1:D:21:PHE:HB3	2.17	0.44
1:D:135:TYR:CE1	1:D:139:ARG:HD2	2.52	0.44
1:D:190:LEU:HD12	1:D:207:ILE:HG13	1.98	0.44
1:D:269:LEU:HD22	1:D:270:VAL:H	1.82	0.44
1:D:447:LYS:NZ	1:D:449:MSE:O	2.41	0.44
1:B:233:ARG:HH11	1:B:233:ARG:CG	2.30	0.44
1:B:426:LEU:HD23	1:B:426:LEU:HA	1.81	0.44
1:B:525:ARG:HA	1:B:525:ARG:HD3	1.76	0.44
1:C:466:HIS:ND1	1:C:466:HIS:N	2.65	0.44
1:D:443:LEU:HA	1:D:460:ILE:HG12	2.00	0.44
1:A:40:HIS:O	1:A:44:ILE:HG13	2.18	0.44
1:A:247:PHE:CD2	1:A:266:LEU:HD13	2.52	0.44
1:B:44:ILE:HD13	1:B:57:PHE:HA	1.99	0.44
1:C:210:MSE:HE1	1:D:233:ARG:HD3	1.99	0.44
1:D:165:LYS:NZ	4:D:602:ATP:O1G	2.51	0.44
1:A:155:LEU:HB3	1:A:269:LEU:HD23	1.99	0.44
1:B:347:LYS:HA	1:B:350:GLN:HB2	1.99	0.44
1:D:189:TRP:C	1:D:189:TRP:CD1	2.96	0.44
1:D:237:ASN:O	1:D:239:LEU:N	2.51	0.44
1:A:88:GLU:HA	1:A:91:ILE:HG22	2.00	0.44
1:A:307:MSE:SE	1:A:308:PRO:HD2	2.67	0.44
1:A:444:LEU:HD12	1:A:444:LEU:HA	1.72	0.44
1:C:4:GLU:OE2	1:C:267:ARG:NH2	2.50	0.44
1:C:131:TYR:CD1	1:C:131:TYR:C	2.95	0.44
1:C:434:LEU:HD22	1:C:444:LEU:HD21	1.98	0.44
1:D:21:PHE:CE2	1:D:23:PRO:HG3	2.52	0.44
1:D:23:PRO:HG2	1:D:53:ARG:HB3	1.99	0.44
1:A:98:PRO:O	1:A:99:ASP:HB2	2.17	0.44
1:A:126:LYS:HD3	1:B:283:GLN:OE1	2.18	0.44
1:A:163:SER:O	1:A:330:PRO:HG2	2.18	0.44
1:B:449:MSE:HE3	1:B:450:PRO:HA	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:477:ILE:HG21	1:C:544:LYS:HE3	1.99	0.44
1:A:23:PRO:HG2	1:A:53:ARG:C	2.43	0.43
1:A:255:GLU:CB	1:A:277:ILE:HG22	2.48	0.43
1:A:463:PHE:O	1:A:467:VAL:HB	2.18	0.43
1:B:2:LEU:HB3	1:B:6:GLU:HB2	2.00	0.43
1:B:69:GLY:C	1:B:71:LEU:H	2.25	0.43
1:B:319:LEU:HA	1:B:348:MSE:HE1	1.99	0.43
1:B:334:MSE:HE2	1:B:338:LYS:CD	2.48	0.43
1:C:1:MSE:HE2	1:C:2:LEU:N	2.33	0.43
1:D:75:PHE:CB	1:D:84:ALA:HB2	2.48	0.43
1:A:30:LEU:HA	1:A:30:LEU:HD23	1.82	0.43
1:A:235:ILE:O	1:A:239:LEU:HB2	2.18	0.43
1:A:243:PRO:O	1:A:244:ASN:C	2.61	0.43
1:A:343:LYS:O	1:A:343:LYS:HG2	2.18	0.43
1:B:44:ILE:O	1:B:53:ARG:HG2	2.18	0.43
1:B:187:ILE:HD13	1:B:246:LEU:HB3	2.00	0.43
1:B:481:GLU:O	1:B:484:LEU:HG	2.17	0.43
1:C:341:GLU:CB	1:C:342:PRO:HD3	2.46	0.43
1:C:426:LEU:HD23	1:C:426:LEU:HA	1.76	0.43
1:D:163:SER:C	1:D:330:PRO:HG2	2.43	0.43
1:A:395:PHE:CD2	1:A:415:PRO:HD3	2.50	0.43
1:B:92:ASP:O	1:B:96:ASN:HB2	2.18	0.43
1:B:527:GLU:C	1:B:530:PRO:HD3	2.43	0.43
1:C:485:LEU:HD12	1:C:486:GLU:H	1.83	0.43
1:D:437:LEU:HD23	1:D:437:LEU:HA	1.77	0.43
1:A:301:PHE:CE1	1:A:305:TYR:HE2	2.36	0.43
1:C:334:MSE:HG3	1:C:338:LYS:HD3	1.99	0.43
1:D:161:ALA:O	1:D:374:LEU:HD21	2.19	0.43
1:B:190:LEU:HD23	1:B:190:LEU:HA	1.83	0.43
1:D:31:GLU:OE1	1:D:38:GLU:HG2	2.18	0.43
1:D:525:ARG:HA	1:D:526:PRO:HD3	1.73	0.43
1:B:185:ASP:N	1:B:244:ASN:O	2.51	0.43
1:B:399:MSE:HE3	1:B:405:ILE:HD11	2.01	0.43
1:C:72:ILE:C	1:C:74:PHE:H	2.26	0.43
1:D:93:PHE:N	1:D:93:PHE:HD1	2.17	0.43
1:A:159:GLY:N	1:A:165:LYS:HD3	2.34	0.43
1:A:529:PHE:N	1:A:530:PRO:HD3	2.34	0.43
1:C:21:PHE:CG	1:C:22:GLU:N	2.86	0.43
1:C:383:GLU:HG2	2:G:753:LEU:HD13	1.99	0.43
1:C:460:ILE:O	1:C:464:LEU:HG	2.18	0.43
1:A:525:ARG:HA	1:A:526:PRO:HD3	1.72	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:190:LEU:HD21	1:D:229:VAL:HG21	2.00	0.43
1:D:133:ARG:O	1:D:137:VAL:HG23	2.18	0.43
1:D:231:LEU:HA	1:D:234:MSE:HE3	2.00	0.43
1:A:216:ASP:H	1:A:218:LEU:CD2	2.32	0.43
1:A:248:VAL:HG22	1:A:269:LEU:HD12	2.00	0.43
1:A:529:PHE:HB3	1:A:532:PHE:CZ	2.53	0.43
1:C:155:LEU:HD23	1:C:155:LEU:HA	1.88	0.43
1:D:6:GLU:O	1:D:9:ALA:HB3	2.19	0.43
1:D:165:LYS:H	1:D:165:LYS:HG3	1.64	0.43
1:D:303:GLU:HA	1:D:309:MSE:HE1	2.01	0.43
1:A:204:PHE:HB3	1:A:231:LEU:HD21	2.01	0.43
1:A:240:ILE:C	1:A:242:ARG:H	2.26	0.43
1:C:20:ASP:HB2	1:D:63:ARG:HB3	2.01	0.43
1:C:75:PHE:HB3	1:C:80:GLN:O	2.18	0.43
1:D:237:ASN:O	1:D:240:ILE:HG23	2.19	0.43
1:D:460:ILE:HD12	1:D:460:ILE:HA	1.75	0.43
1:A:187:ILE:HD12	1:A:246:LEU:O	2.18	0.42
1:B:326:SER:O	1:B:328:GLY:N	2.51	0.42
1:D:444:LEU:HD12	1:D:444:LEU:HA	1.76	0.42
1:A:150:LEU:HD13	1:A:153:PHE:HB3	2.01	0.42
1:B:460:ILE:HA	1:B:460:ILE:HD12	1.85	0.42
1:D:426:LEU:HD23	1:D:426:LEU:HA	1.78	0.42
2:H:752:PHE:HB3	2:H:753:LEU:HA	2.01	0.42
1:A:247:PHE:HD2	1:A:266:LEU:HB3	1.84	0.42
1:B:178:GLN:HB3	1:B:183:ASN:OD1	2.19	0.42
1:C:187:ILE:HD13	1:C:246:LEU:HB3	2.01	0.42
1:A:13:ALA:HB2	1:A:86:PHE:CE1	2.54	0.42
1:A:444:LEU:HB3	1:A:445:SER:H	1.61	0.42
1:B:412:CYS:O	1:B:483:ARG:NH1	2.51	0.42
1:C:456:ILE:HG13	1:C:460:ILE:CG2	2.49	0.42
1:C:529:PHE:HB3	1:C:532:PHE:CZ	2.53	0.42
1:D:157:LEU:O	1:D:165:LYS:HD2	2.20	0.42
1:D:313:GLU:O	1:D:316:GLU:HG3	2.20	0.42
1:D:327:SER:HB2	1:D:459:ILE:CD1	2.50	0.42
1:A:26:ALA:HA	1:A:74:PHE:CE1	2.55	0.42
1:A:111:SER:C	1:A:113:GLN:N	2.78	0.42
1:B:240:ILE:C	1:B:242:ARG:H	2.27	0.42
1:C:462:MSE:HE3	1:C:462:MSE:HB2	1.92	0.42
1:D:24:ARG:HA	1:D:27:LEU:HB2	2.01	0.42
1:D:93:PHE:N	1:D:93:PHE:CD1	2.87	0.42
1:A:397:VAL:C	1:A:399:MSE:H	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:343:LYS:O	1:B:343:LYS:HG2	2.18	0.42
1:B:474:ALA:HB2	1:B:547:ALA:HB1	2.01	0.42
1:D:35:ILE:O	1:D:64:GLN:HG3	2.19	0.42
1:D:121:LEU:HD23	1:D:121:LEU:HA	1.82	0.42
1:D:532:PHE:C	1:D:534:GLN:N	2.77	0.42
1:A:435:LYS:HD3	1:A:435:LYS:HA	1.90	0.42
1:C:124:VAL:HA	1:C:125:PRO:HD3	1.82	0.42
1:D:30:LEU:C	1:D:36:PHE:HB2	2.45	0.42
1:D:100:LEU:HB3	1:D:101:LEU:H	1.71	0.42
1:A:69:GLY:C	1:A:71:LEU:H	2.27	0.42
1:B:255:GLU:CB	1:B:277:ILE:HG22	2.50	0.42
1:C:385:LEU:HD22	1:C:389:ASP:HB3	2.02	0.42
1:C:403:VAL:HG12	1:C:405:ILE:HG22	2.02	0.42
1:A:86:PHE:HE2	1:A:109:GLN:HE22	1.67	0.42
1:A:368:PRO:HD3	1:B:279:ASN:O	2.18	0.42
1:B:200:THR:HG21	1:B:256:GLU:HB3	2.02	0.42
1:B:212:LYS:HE2	1:B:212:LYS:HA	2.02	0.42
1:B:235:ILE:O	1:B:239:LEU:HB2	2.20	0.42
1:B:405:ILE:HA	1:B:406:PRO:HD3	1.79	0.42
1:C:157:LEU:O	1:C:165:LYS:HD2	2.19	0.42
1:D:405:ILE:HG12	1:D:410:TRP:CE3	2.55	0.42
1:B:243:PRO:O	1:B:244:ASN:C	2.63	0.42
1:B:354:LYS:O	1:B:358:ARG:HB2	2.20	0.42
1:B:388:GLU:CD	1:B:388:GLU:H	2.27	0.42
1:B:456:ILE:HD13	1:B:461:HIS:HD2	1.84	0.42
1:C:23:PRO:HG2	1:C:53:ARG:CB	2.50	0.42
1:C:33:LYS:HD3	1:C:70:PRO:HG3	2.01	0.42
1:C:107:ALA:C	1:C:109:GLN:H	2.28	0.42
1:C:188:VAL:HB	1:C:247:PHE:HD1	1.84	0.42
1:C:255:GLU:HB3	1:C:277:ILE:HG22	2.02	0.42
1:C:307:MSE:SE	1:C:308:PRO:HD2	2.70	0.42
1:C:443:LEU:O	1:C:460:ILE:HG21	2.20	0.42
1:C:460:ILE:HD12	1:C:460:ILE:HA	1.76	0.42
1:D:240:ILE:HG13	1:D:241:ASP:OD1	2.19	0.42
1:D:385:LEU:HD11	1:D:437:LEU:HD21	2.01	0.42
1:D:529:PHE:HB3	1:D:532:PHE:CZ	2.55	0.42
1:A:211:LEU:HD11	1:A:239:LEU:HD23	2.02	0.41
1:A:524:ILE:HD12	1:A:525:ARG:H	1.85	0.41
1:B:131:TYR:HD1	4:B:602:ATP:HN62	1.67	0.41
1:B:401:PRO:HB2	1:B:458:HIS:CE1	2.54	0.41
1:B:481:GLU:HA	1:B:484:LEU:HG	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:237:ASN:HD22	1:C:237:ASN:HA	1.66	0.41
1:C:302:LEU:HD13	1:C:319:LEU:HD21	2.01	0.41
1:D:44:ILE:HD13	1:D:57:PHE:HD1	1.84	0.41
1:D:154:PHE:CE2	1:D:262:GLN:HG3	2.54	0.41
1:D:165:LYS:N	4:D:602:ATP:O2B	2.53	0.41
1:D:292:SER:HB3	1:D:328:GLY:HA3	2.02	0.41
1:A:397:VAL:O	1:A:399:MSE:N	2.52	0.41
1:B:240:ILE:HG13	1:B:241:ASP:N	2.35	0.41
1:B:351:LEU:HD13	1:B:363:VAL:HG23	2.02	0.41
1:C:126:LYS:HE2	1:C:126:LYS:HB3	1.87	0.41
1:C:189:TRP:CD2	1:C:248:VAL:HG11	2.55	0.41
1:C:347:LYS:HD3	1:C:347:LYS:O	2.19	0.41
1:C:349:ALA:HA	1:C:352:ASN:HD22	1.85	0.41
1:D:156:PHE:CE2	1:D:278:SER:HB3	2.56	0.41
1:A:482:GLN:C	1:A:484:LEU:H	2.27	0.41
1:C:39:ASP:HA	1:C:42:GLU:HB2	2.02	0.41
1:C:126:LYS:HZ2	1:D:283:GLN:HG3	1.82	0.41
1:D:29:TYR:CD2	1:D:74:PHE:HB2	2.55	0.41
1:A:1:MSE:HG2	1:A:63:ARG:HA	2.02	0.41
1:A:115:LEU:HD23	1:A:115:LEU:HA	1.67	0.41
1:A:449:MSE:HE2	1:A:449:MSE:HB2	1.73	0.41
1:A:456:ILE:HD13	1:A:461:HIS:HD2	1.86	0.41
1:B:336:PHE:CE2	1:B:348:MSE:SE	3.23	0.41
1:C:333:LEU:O	1:C:336:PHE:HB2	2.21	0.41
1:D:105:VAL:HG22	1:D:110:PHE:CE2	2.55	0.41
1:A:314:LYS:HA	1:A:317:ASP:HB2	2.02	0.41
1:A:395:PHE:C	1:A:397:VAL:H	2.28	0.41
1:B:8:ARG:HD3	1:B:90:TYR:OH	2.21	0.41
1:B:40:HIS:O	1:B:44:ILE:HG13	2.21	0.41
1:B:269:LEU:HD22	1:B:270:VAL:H	1.86	0.41
1:B:273:ARG:NE	4:B:602:ATP:O3G	2.46	0.41
1:B:485:LEU:HD23	1:B:485:LEU:HA	1.86	0.41
1:D:208:LEU:HD23	1:D:235:ILE:HA	2.02	0.41
1:D:343:LYS:HB3	1:D:343:LYS:HE3	1.78	0.41
1:D:438:SER:O	1:D:442:ALA:HA	2.20	0.41
1:A:157:LEU:HD11	1:A:168:ILE:CG2	2.50	0.41
1:A:274:ASP:O	1:A:277:ILE:HG12	2.20	0.41
1:A:313:GLU:O	1:A:316:GLU:HG3	2.20	0.41
1:B:247:PHE:CE2	1:B:266:LEU:HD22	2.56	0.41
1:B:313:GLU:O	1:B:316:GLU:HG3	2.21	0.41
1:C:9:ALA:HA	1:C:12:THR:HG22	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:128:MSE:HG3	1:C:167:VAL:CG2	2.45	0.41
1:D:110:PHE:O	1:D:112:ARG:N	2.54	0.41
1:D:393:LEU:HD11	1:D:443:LEU:HD13	2.03	0.41
1:D:407:VAL:HG22	1:D:454:PHE:HE1	1.84	0.41
1:A:131:TYR:CD1	1:A:131:TYR:C	2.98	0.41
1:A:437:LEU:HA	1:A:437:LEU:HD22	1.75	0.41
1:B:159:GLY:N	1:B:165:LYS:HD3	2.35	0.41
1:B:444:LEU:HD12	1:B:444:LEU:HA	1.72	0.41
1:C:39:ASP:O	1:C:43:LEU:HD13	2.20	0.41
1:C:326:SER:O	1:C:327:SER:C	2.63	0.41
1:D:26:ALA:HB2	1:D:74:PHE:CE2	2.56	0.41
1:D:283:GLN:HG2	1:D:284:THR:N	2.35	0.41
1:D:301:PHE:CE1	1:D:305:TYR:HE1	2.39	0.41
1:D:466:HIS:ND1	1:D:466:HIS:N	2.67	0.41
1:B:30:LEU:HD23	1:B:30:LEU:HA	1.71	0.41
1:C:30:LEU:HD23	1:C:30:LEU:HA	1.82	0.41
1:C:257:THR:HG22	1:C:258:ILE:HD12	2.03	0.41
1:C:408:LYS:HD2	1:C:408:LYS:HA	1.76	0.41
1:C:434:LEU:HA	1:C:437:LEU:HB2	2.02	0.41
1:C:445:SER:HB3	1:C:455:LYS:O	2.21	0.41
1:A:157:LEU:HD22	1:A:290:VAL:CG2	2.49	0.41
1:A:527:GLU:HA	1:A:530:PRO:HG3	2.02	0.41
1:B:124:VAL:HA	1:B:125:PRO:HD3	1.90	0.41
1:B:207:ILE:O	1:B:210:MSE:N	2.54	0.41
1:C:322:THR:O	1:C:326:SER:OG	2.37	0.41
1:C:395:PHE:CD2	1:C:415:PRO:HD3	2.51	0.41
1:D:104:VAL:HB	1:D:105:VAL:H	1.63	0.41
1:D:326:SER:O	1:D:327:SER:C	2.63	0.41
1:A:347:LYS:O	1:A:350:GLN:HB2	2.21	0.41
1:B:347:LYS:O	1:B:347:LYS:HD3	2.20	0.41
1:C:205:THR:HA	1:C:231:LEU:HD11	2.02	0.41
1:C:243:PRO:O	1:C:244:ASN:C	2.64	0.41
1:C:406:PRO:CA	1:C:453:THR:HG22	2.47	0.41
1:C:444:LEU:HB3	1:C:445:SER:H	1.55	0.41
1:D:131:TYR:CD1	1:D:131:TYR:C	2.98	0.41
1:D:399:MSE:HG3	1:D:410:TRP:NE1	2.36	0.41
1:A:9:ALA:HA	1:A:12:THR:HG22	2.04	0.40
1:B:154:PHE:CE2	1:B:262:GLN:HG3	2.55	0.40
1:C:292:SER:HB3	1:C:328:GLY:HA3	2.03	0.40
1:C:385:LEU:HD23	1:C:440:ARG:HH21	1.86	0.40
1:D:445:SER:HB3	1:D:455:LYS:O	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:PHE:CE2	1:A:104:VAL:HG11	2.56	0.40
1:A:253:VAL:HG11	1:A:380:ARG:HD3	2.03	0.40
1:C:69:GLY:C	1:C:71:LEU:H	2.30	0.40
1:C:415:PRO:O	1:C:416:VAL:HG13	2.21	0.40
1:D:187:ILE:HD13	1:D:246:LEU:HB3	2.04	0.40
1:A:189:TRP:CD2	1:A:248:VAL:HG11	2.55	0.40
1:A:212:LYS:HE2	1:A:212:LYS:HA	2.03	0.40
1:B:395:PHE:C	1:B:397:VAL:H	2.29	0.40
1:B:409:LEU:HA	1:B:409:LEU:HD22	1.85	0.40
1:B:430:VAL:O	1:B:434:LEU:HG	2.21	0.40
1:C:240:ILE:HG13	1:C:241:ASP:N	2.37	0.40
1:C:538:LYS:C	1:C:540:TYR:H	2.29	0.40
1:D:157:LEU:HD22	1:D:290:VAL:HG22	2.02	0.40
1:D:325:LEU:HA	1:D:325:LEU:HD23	1.83	0.40
1:A:240:ILE:HG13	1:A:241:ASP:OD1	2.22	0.40
1:A:466:HIS:ND1	1:A:466:HIS:N	2.69	0.40
1:B:395:PHE:CD2	1:B:415:PRO:HD3	2.52	0.40
1:B:409:LEU:HD21	1:B:530:PRO:HB3	2.02	0.40
1:C:7:CYS:HB3	1:C:243:PRO:HG3	2.02	0.40
1:C:307:MSE:HA	1:C:308:PRO:HD3	1.85	0.40
1:C:318:VAL:O	1:C:321:LYS:HB3	2.20	0.40
1:D:130:CYS:O	1:D:304:ALA:HB1	2.22	0.40
1:D:255:GLU:CB	1:D:277:ILE:HG22	2.51	0.40
1:D:273:ARG:HD2	1:D:381:CYS:SG	2.60	0.40
1:D:349:ALA:HA	1:D:352:ASN:HB2	2.03	0.40
1:A:98:PRO:C	1:A:100:LEU:H	2.29	0.40
1:A:115:LEU:HD21	1:A:180:ILE:O	2.21	0.40
1:A:477:ILE:HD11	1:A:543:LEU:HB2	2.02	0.40
1:B:21:PHE:CG	1:B:22:GLU:N	2.89	0.40
1:B:155:LEU:HD23	1:B:155:LEU:HA	1.89	0.40
1:C:233:ARG:HA	1:C:236:CYS:SG	2.62	0.40
2:G:752:PHE:O	2:G:753:LEU:HG	2.22	0.40
1:D:27:LEU:HD23	1:D:27:LEU:HA	1.86	0.40
1:D:487:ILE:HG22	1:D:531:LYS:HG2	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:LEU:O	1:B:113:GLN:NE2[3_545]	2.19	0.01

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	504/549 (92%)	404 (80%)	77 (15%)	23 (5%)	2	12
1	B	505/549 (92%)	406 (80%)	73 (14%)	26 (5%)	1	10
1	C	504/549 (92%)	406 (81%)	72 (14%)	26 (5%)	1	10
1	D	505/549 (92%)	405 (80%)	76 (15%)	24 (5%)	2	11
2	E	4/8 (50%)	2 (50%)	1 (25%)	1 (25%)	0	0
2	F	3/8 (38%)	1 (33%)	1 (33%)	1 (33%)	0	0
2	G	3/8 (38%)	1 (33%)	0	2 (67%)	0	0
2	H	4/8 (50%)	1 (25%)	2 (50%)	1 (25%)	0	0
All	All	2032/2228 (91%)	1626 (80%)	302 (15%)	104 (5%)	1	10

All (104) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	327	SER
1	A	341	GLU
1	A	398	VAL
1	A	445	SER
1	A	486	GLU
2	E	751	ASN
1	B	109	GLN
1	B	127	GLN
1	B	327	SER
1	B	341	GLU
1	B	398	VAL
1	B	445	SER
1	C	127	GLN
1	C	327	SER
1	C	341	GLU
1	C	398	VAL
1	C	445	SER

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Mol	Chain	Res	Type
1	D	35	ILE
1	D	127	GLN
1	D	284	THR
1	D	327	SER
1	D	341	GLU
1	D	398	VAL
1	D	445	SER
1	A	112	ARG
1	A	238	ALA
1	A	244	ASN
1	A	396	ALA
1	A	465	LYS
1	A	522	THR
1	B	244	ASN
1	B	396	ALA
1	B	465	LYS
1	B	522	THR
1	C	112	ARG
1	C	160	ARG
1	C	244	ASN
1	C	396	ALA
1	C	465	LYS
1	C	522	THR
2	G	751	ASN
1	D	104	VAL
1	D	160	ARG
1	D	208	LEU
1	D	244	ASN
1	D	465	LYS
1	D	522	THR
2	H	750	PHE
1	A	2	LEU
1	A	73	ASP
1	A	160	ARG
1	A	208	LEU
1	A	284	THR
1	B	2	LEU
1	B	104	VAL
1	B	160	ARG
1	B	208	LEU
1	B	238	ALA
1	B	284	THR

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Mol	Chain	Res	Type
2	F	751	ASN
1	C	2	LEU
1	C	73	ASP
1	C	208	LEU
1	C	238	ALA
1	C	284	THR
1	C	486	GLU
1	D	111	SER
1	D	238	ALA
1	D	396	ALA
1	A	485	LEU
1	B	73	ASP
1	B	176	SER
1	C	485	LEU
2	G	752	PHE
1	D	3	CYS
1	A	484	LEU
1	B	35	ILE
1	B	257	THR
1	B	328	GLY
1	B	354	LYS
1	D	34	ASN
1	D	102	ARG
1	D	237	ASN
1	A	237	ASN
1	A	257	THR
1	B	102	ARG
1	C	3	CYS
1	C	33	LYS
1	C	237	ASN
1	C	257	THR
1	D	64	GLN
1	B	106	ILE
1	B	311	VAL
1	C	108	PRO
1	D	311	VAL
1	A	70	PRO
1	A	108	PRO
1	A	311	VAL
1	B	70	PRO
1	C	70	PRO
1	C	311	VAL

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Mol	Chain	Res	Type
1	D	103	PRO
1	D	328	GLY
1	C	35	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	A	462/483 (96%)	379 (82%)	83 (18%)	2	7
1	B	464/483 (96%)	376 (81%)	88 (19%)	1	6
1	C	462/483 (96%)	370 (80%)	92 (20%)	1	5
1	D	464/483 (96%)	366 (79%)	98 (21%)	1	3
2	E	5/6 (83%)	4 (80%)	1 (20%)	1	4
2	F	5/6 (83%)	5 (100%)	0	100	100
2	G	5/6 (83%)	5 (100%)	0	100	100
2	H	6/6 (100%)	5 (83%)	1 (17%)	2	9
All	All	1873/1956 (96%)	1510 (81%)	363 (19%)	1	5

All (363) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	8	ARG
1	A	15	THR
1	A	17	LEU
1	A	43	LEU
1	A	46	LYS
1	A	47	MSE
1	A	51	LEU
1	A	52	GLU
1	A	67	GLU
1	A	68	LEU
1	A	77	TYR

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Mol	Chain	Res	Type
1	A	82	HIS
1	A	95	ILE
1	A	100	LEU
1	A	113	GLN
1	A	115	LEU
1	A	116	ASP
1	A	126	LYS
1	A	130	CYS
1	A	131	TYR
1	A	155	LEU
1	A	157	LEU
1	A	167	VAL
1	A	174	SER
1	A	179	LEU
1	A	180	ILE
1	A	185	ASP
1	A	198	LYS
1	A	203	LEU
1	A	205	THR
1	A	217	LEU
1	A	220	PHE
1	A	223	VAL
1	A	226	VAL
1	A	227	THR
1	A	233	ARG
1	A	239	LEU
1	A	241	ASP
1	A	253	VAL
1	A	255	GLU
1	A	257	THR
1	A	264	LEU
1	A	269	LEU
1	A	277	ILE
1	A	284	THR
1	A	286	GLU
1	A	290	VAL
1	A	296	ASP
1	A	316	GLU
1	A	332	THR
1	A	344	THR
1	A	348	MSE
1	A	355	LEU

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Mol	Chain	Res	Type
1	A	363	VAL
1	A	364	GLU
1	A	366	ILE
1	A	369	TYR
1	A	374	LEU
1	A	381	CYS
1	A	388	GLU
1	A	404	ASP
1	A	405	ILE
1	A	409	LEU
1	A	410	TRP
1	A	414	ILE
1	A	416	VAL
1	A	437	LEU
1	A	444	LEU
1	A	449	MSE
1	A	451	VAL
1	A	452	LEU
1	A	456	ILE
1	A	457	ASP
1	A	459	ILE
1	A	460	ILE
1	A	466	HIS
1	A	467	VAL
1	A	485	LEU
1	A	525	ARG
1	A	533	MSE
1	A	539	PHE
1	A	542	SER
2	E	750	PHE
1	B	1	MSE
1	B	4	GLU
1	B	8	ARG
1	B	15	THR
1	B	17	LEU
1	B	35	ILE
1	B	43	LEU
1	B	46	LYS
1	B	51	LEU
1	B	52	GLU
1	B	64	GLN
1	B	67	GLU

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Mol	Chain	Res	Type
1	B	68	LEU
1	B	77	TYR
1	B	97	GLU
1	B	99	ASP
1	B	104	VAL
1	B	106	ILE
1	B	109	GLN
1	B	112	ARG
1	B	116	ASP
1	B	117	ARG
1	B	126	LYS
1	B	130	CYS
1	B	131	TYR
1	B	148	CYS
1	B	155	LEU
1	B	157	LEU
1	B	167	VAL
1	B	174	SER
1	B	179	LEU
1	B	180	ILE
1	B	185	ASP
1	B	198	LYS
1	B	203	LEU
1	B	208	LEU
1	B	210	MSE
1	B	217	LEU
1	B	218	LEU
1	B	220	PHE
1	B	223	VAL
1	B	226	VAL
1	B	227	THR
1	B	233	ARG
1	B	239	LEU
1	B	241	ASP
1	B	253	VAL
1	B	255	GLU
1	B	257	THR
1	B	262	GLN
1	B	269	LEU
1	B	277	ILE
1	B	283	GLN
1	B	284	THR

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Mol	Chain	Res	Type
1	B	290	VAL
1	B	296	ASP
1	B	316	GLU
1	B	332	THR
1	B	344	THR
1	B	355	LEU
1	B	363	VAL
1	B	364	GLU
1	B	366	ILE
1	B	369	TYR
1	B	374	LEU
1	B	381	CYS
1	B	382	VAL
1	B	404	ASP
1	B	405	ILE
1	B	409	LEU
1	B	410	TRP
1	B	414	ILE
1	B	416	VAL
1	B	444	LEU
1	B	451	VAL
1	B	452	LEU
1	B	457	ASP
1	B	459	ILE
1	B	460	ILE
1	B	466	HIS
1	B	467	VAL
1	B	485	LEU
1	B	528	ASP
1	B	533	MSE
1	B	539	PHE
1	B	542	SER
1	B	546	PHE
1	B	548	CYS
1	C	1	MSE
1	C	8	ARG
1	C	11	SER
1	C	15	THR
1	C	17	LEU
1	C	46	LYS
1	C	47	MSE
1	C	51	LEU

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Mol	Chain	Res	Type
1	C	52	GLU
1	C	67	GLU
1	C	68	LEU
1	C	77	TYR
1	C	78	ASN
1	C	79	ASN
1	C	82	HIS
1	C	99	ASP
1	C	100	LEU
1	C	109	GLN
1	C	113	GLN
1	C	116	ASP
1	C	117	ARG
1	C	126	LYS
1	C	130	CYS
1	C	131	TYR
1	C	132	ILE
1	C	139	ARG
1	C	148	CYS
1	C	155	LEU
1	C	157	LEU
1	C	160	ARG
1	C	167	VAL
1	C	174	SER
1	C	179	LEU
1	C	180	ILE
1	C	185	ASP
1	C	198	LYS
1	C	203	LEU
1	C	217	LEU
1	C	220	PHE
1	C	223	VAL
1	C	224	GLU
1	C	226	VAL
1	C	227	THR
1	C	233	ARG
1	C	237	ASN
1	C	239	LEU
1	C	241	ASP
1	C	253	VAL
1	C	255	GLU
1	C	257	THR

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Mol	Chain	Res	Type
1	C	264	LEU
1	C	269	LEU
1	C	277	ILE
1	C	284	THR
1	C	286	GLU
1	C	290	VAL
1	C	296	ASP
1	C	307	MSE
1	C	309	MSE
1	C	316	GLU
1	C	319	LEU
1	C	326	SER
1	C	332	THR
1	C	355	LEU
1	C	360	LEU
1	C	361	VAL
1	C	363	VAL
1	C	364	GLU
1	C	366	ILE
1	C	369	TYR
1	C	374	LEU
1	C	381	CYS
1	C	388	GLU
1	C	405	ILE
1	C	409	LEU
1	C	410	TRP
1	C	414	ILE
1	C	416	VAL
1	C	437	LEU
1	C	444	LEU
1	C	451	VAL
1	C	452	LEU
1	C	456	ILE
1	C	457	ASP
1	C	459	ILE
1	C	460	ILE
1	C	466	HIS
1	C	467	VAL
1	C	485	LEU
1	C	524	ILE
1	C	525	ARG
1	C	533	MSE

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Mol	Chain	Res	Type
1	D	1	MSE
1	D	3	CYS
1	D	12	THR
1	D	15	THR
1	D	17	LEU
1	D	18	ILE
1	D	19	HIS
1	D	31	GLU
1	D	35	ILE
1	D	37	THR
1	D	38	GLU
1	D	39	ASP
1	D	43	LEU
1	D	46	LYS
1	D	51	LEU
1	D	52	GLU
1	D	64	GLN
1	D	68	LEU
1	D	71	LEU
1	D	77	TYR
1	D	82	HIS
1	D	83	LEU
1	D	93	PHE
1	D	104	VAL
1	D	110	PHE
1	D	116	ASP
1	D	117	ARG
1	D	126	LYS
1	D	130	CYS
1	D	131	TYR
1	D	148	CYS
1	D	155	LEU
1	D	157	LEU
1	D	160	ARG
1	D	167	VAL
1	D	174	SER
1	D	176	SER
1	D	179	LEU
1	D	180	ILE
1	D	185	ASP
1	D	198	LYS
1	D	203	LEU

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Mol	Chain	Res	Type
1	D	217	LEU
1	D	218	LEU
1	D	220	PHE
1	D	223	VAL
1	D	226	VAL
1	D	227	THR
1	D	233	ARG
1	D	239	LEU
1	D	241	ASP
1	D	253	VAL
1	D	255	GLU
1	D	257	THR
1	D	262	GLN
1	D	264	LEU
1	D	269	LEU
1	D	277	ILE
1	D	284	THR
1	D	286	GLU
1	D	290	VAL
1	D	296	ASP
1	D	307	MSE
1	D	316	GLU
1	D	319	LEU
1	D	337	PHE
1	D	345	PHE
1	D	355	LEU
1	D	360	LEU
1	D	361	VAL
1	D	363	VAL
1	D	364	GLU
1	D	366	ILE
1	D	369	TYR
1	D	374	LEU
1	D	381	CYS
1	D	382	VAL
1	D	388	GLU
1	D	404	ASP
1	D	405	ILE
1	D	409	LEU
1	D	410	TRP
1	D	414	ILE
1	D	416	VAL

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Mol	Chain	Res	Type
1	D	444	LEU
1	D	451	VAL
1	D	452	LEU
1	D	456	ILE
1	D	457	ASP
1	D	459	ILE
1	D	460	ILE
1	D	466	HIS
1	D	467	VAL
1	D	491	ASN
1	D	525	ARG
1	D	538	LYS
1	D	539	PHE
1	D	548	CYS
2	H	751	ASN

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

Mol	Chain	Res	Type
1	A	171	GLN
1	A	320	ASN
1	A	458	HIS
1	A	471	GLN
1	B	56	ASN
1	B	123	ASN
1	B	136	HIS
1	B	158	HIS
1	B	183	ASN
1	B	254	GLN
1	B	458	HIS
1	C	56	ASN
1	C	78	ASN
1	C	158	HIS
1	C	350	GLN
1	C	458	HIS
1	D	109	GLN
1	D	136	HIS
1	D	171	GLN
1	D	183	ASN
1	D	237	ASN
1	D	283	GLN
1	D	458	HIS

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Mol	Chain	Res	Type
1	D	461	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	ATP	C	602	3	32,33,33	1.43	5 (15%)	48,52,52	1.71	7 (14%)
4	ATP	A	602	3	32,33,33	1.39	4 (12%)	48,52,52	1.68	8 (16%)
4	ATP	D	602	3	32,33,33	1.45	6 (18%)	48,52,52	1.77	9 (18%)
4	ATP	B	602	3	32,33,33	1.40	4 (12%)	48,52,52	1.83	10 (20%)

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
4	ATP	C	602	3	-	2/22/38/38	0/3/3/3
4	ATP	A	602	3	-	1/22/38/38	0/3/3/3
4	ATP	D	602	3	-	2/22/38/38	0/3/3/3
4	ATP	B	602	3	-	3/22/38/38	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	602	ATP	C5-C4	4.83	1.47	1.39
4	C	602	ATP	C5-C4	4.82	1.47	1.39
4	B	602	ATP	C5-C4	4.79	1.47	1.39
4	A	602	ATP	C5-C4	4.78	1.47	1.39
4	B	602	ATP	C5-C6	2.81	1.48	1.41
4	D	602	ATP	C5-C6	2.75	1.48	1.41
4	C	602	ATP	C5-C6	2.73	1.48	1.41
4	A	602	ATP	C5-C6	2.65	1.48	1.41
4	D	602	ATP	PB-O3B	2.37	1.62	1.59
4	C	602	ATP	C5-N7	-2.33	1.34	1.39
4	A	602	ATP	C5-N7	-2.30	1.34	1.39
4	B	602	ATP	C5-N7	-2.27	1.34	1.39
4	D	602	ATP	C8-N7	2.24	1.36	1.31
4	D	602	ATP	C5-N7	-2.23	1.35	1.39
4	B	602	ATP	C8-N7	2.23	1.36	1.31
4	A	602	ATP	C8-N7	2.22	1.36	1.31
4	C	602	ATP	PB-O3B	2.15	1.61	1.59
4	C	602	ATP	C8-N7	2.12	1.35	1.31
4	D	602	ATP	PB-O3A	2.03	1.61	1.59

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	602	ATP	C5-C4-N3	-6.00	118.46	126.72
4	D	602	ATP	C5-C4-N3	-5.97	118.50	126.72
4	B	602	ATP	C5-C4-N3	-5.96	118.52	126.72
4	A	602	ATP	C5-C4-N3	-5.72	118.84	126.72
4	D	602	ATP	N3-C4-N9	5.07	135.79	127.17
4	C	602	ATP	N3-C4-N9	4.99	135.65	127.17
4	B	602	ATP	N3-C4-N9	4.89	135.49	127.17
4	A	602	ATP	N3-C4-N9	4.78	135.30	127.17
4	B	602	ATP	C2-N3-C4	3.79	121.09	111.83
4	D	602	ATP	C2-N3-C4	3.70	120.87	111.83
4	C	602	ATP	C2-N3-C4	3.60	120.61	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	602	ATP	C2-N3-C4	3.44	120.22	111.83
4	B	602	ATP	C4-C5-N7	-3.39	106.70	110.58
4	B	602	ATP	N3-C2-N1	-3.28	123.61	128.58
4	D	602	ATP	C4-C5-N7	-3.20	106.92	110.58
4	C	602	ATP	C4-C5-N7	-3.18	106.94	110.58
4	A	602	ATP	C4-C5-N7	-3.16	106.97	110.58
4	D	602	ATP	N3-C2-N1	-3.16	123.80	128.58
4	D	602	ATP	C4-N9-C8	3.04	108.92	105.74
4	A	602	ATP	C4-N9-C8	2.91	108.80	105.74
4	C	602	ATP	N3-C2-N1	-2.87	124.24	128.58
4	B	602	ATP	C2'-C1'-N9	-2.83	106.26	113.30
4	A	602	ATP	N3-C2-N1	-2.81	124.33	128.58
4	B	602	ATP	C4-N9-C8	2.71	108.58	105.74
4	C	602	ATP	C4-N9-C8	2.63	108.50	105.74
4	B	602	ATP	C5-N7-C8	2.59	107.52	103.45
4	D	602	ATP	C5-N7-C8	2.57	107.50	103.45
4	A	602	ATP	C5-N7-C8	2.46	107.32	103.45
4	C	602	ATP	C5-N7-C8	2.37	107.18	103.45
4	B	602	ATP	O4'-C1'-N9	2.34	112.58	108.09
4	D	602	ATP	C2'-C1'-N9	-2.19	107.86	113.30
4	D	602	ATP	N9-C8-N7	-2.11	110.94	113.94
4	B	602	ATP	C6-C5-N7	2.03	136.01	132.09
4	A	602	ATP	N9-C8-N7	-2.03	111.06	113.94

There are no chirality outliers.

All (8) torsion outliers are listed below:

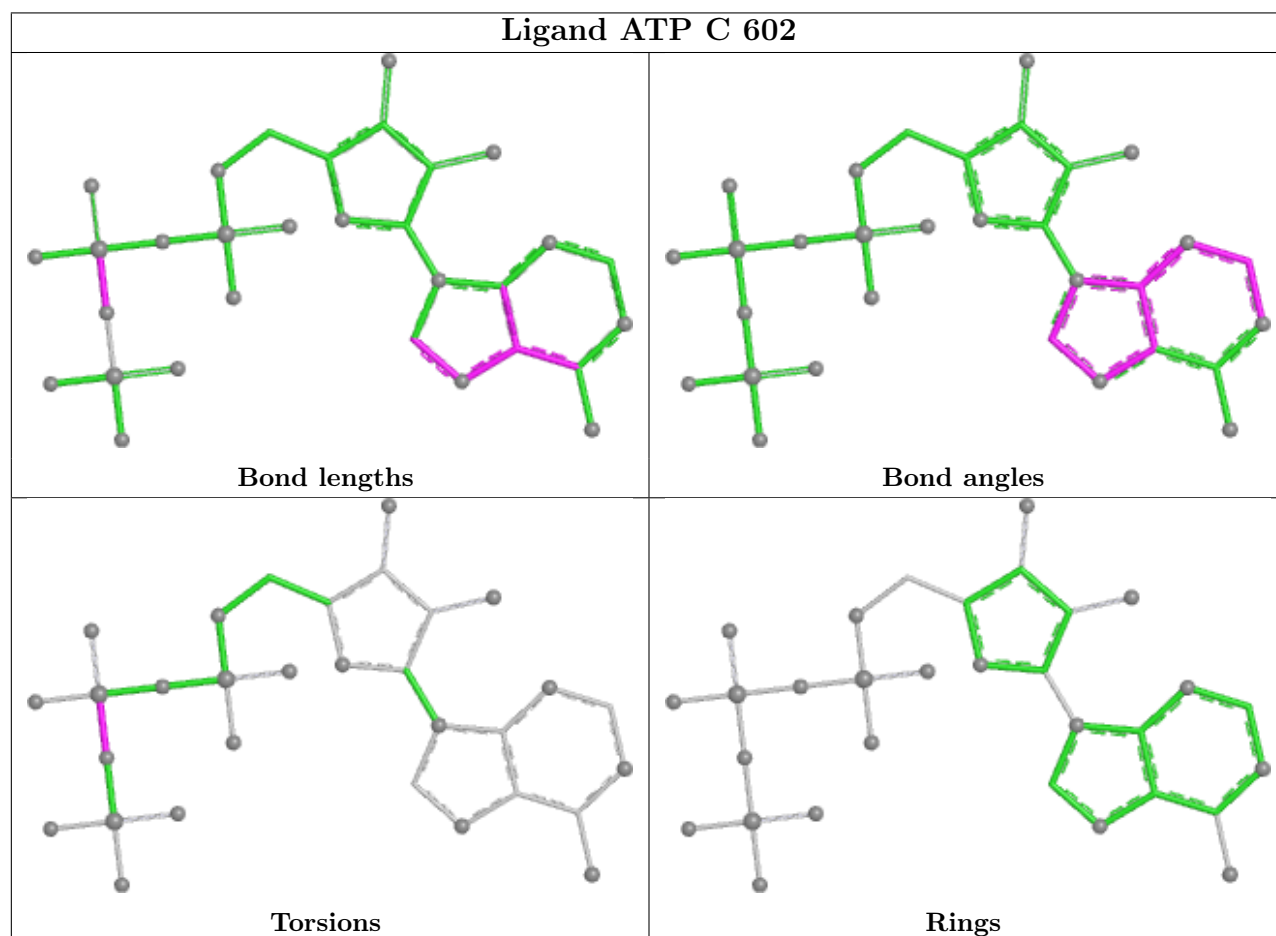
Mol	Chain	Res	Type	Atoms
4	B	602	ATP	O4'-C1'-N9-C8
4	B	602	ATP	O4'-C1'-N9-C4
4	D	602	ATP	O4'-C1'-N9-C8
4	D	602	ATP	O4'-C1'-N9-C4
4	A	602	ATP	PG-O3B-PB-O2B
4	B	602	ATP	PG-O3B-PB-O2B
4	C	602	ATP	PG-O3B-PB-O1B
4	C	602	ATP	PG-O3B-PB-O2B

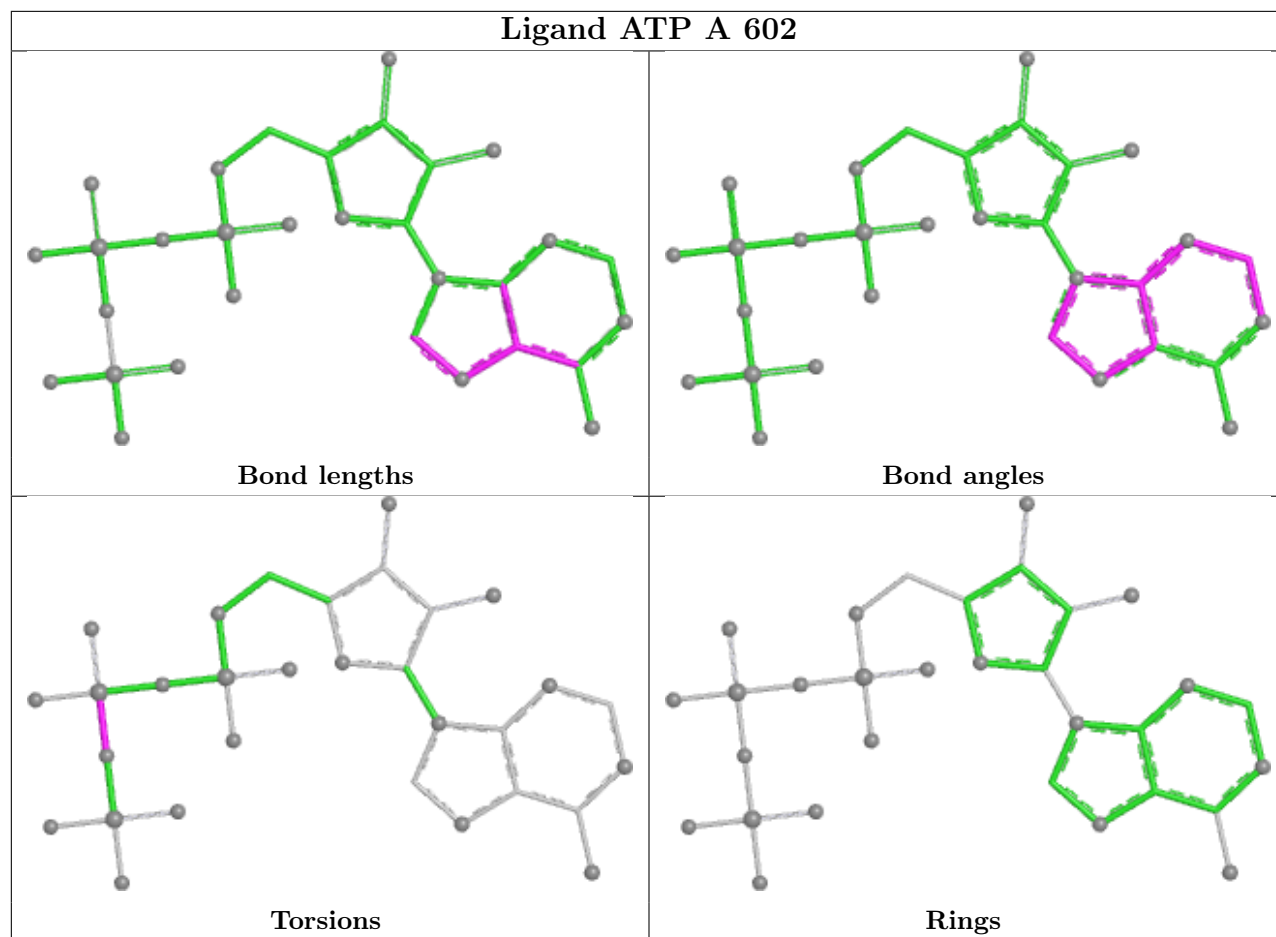
There are no ring outliers.

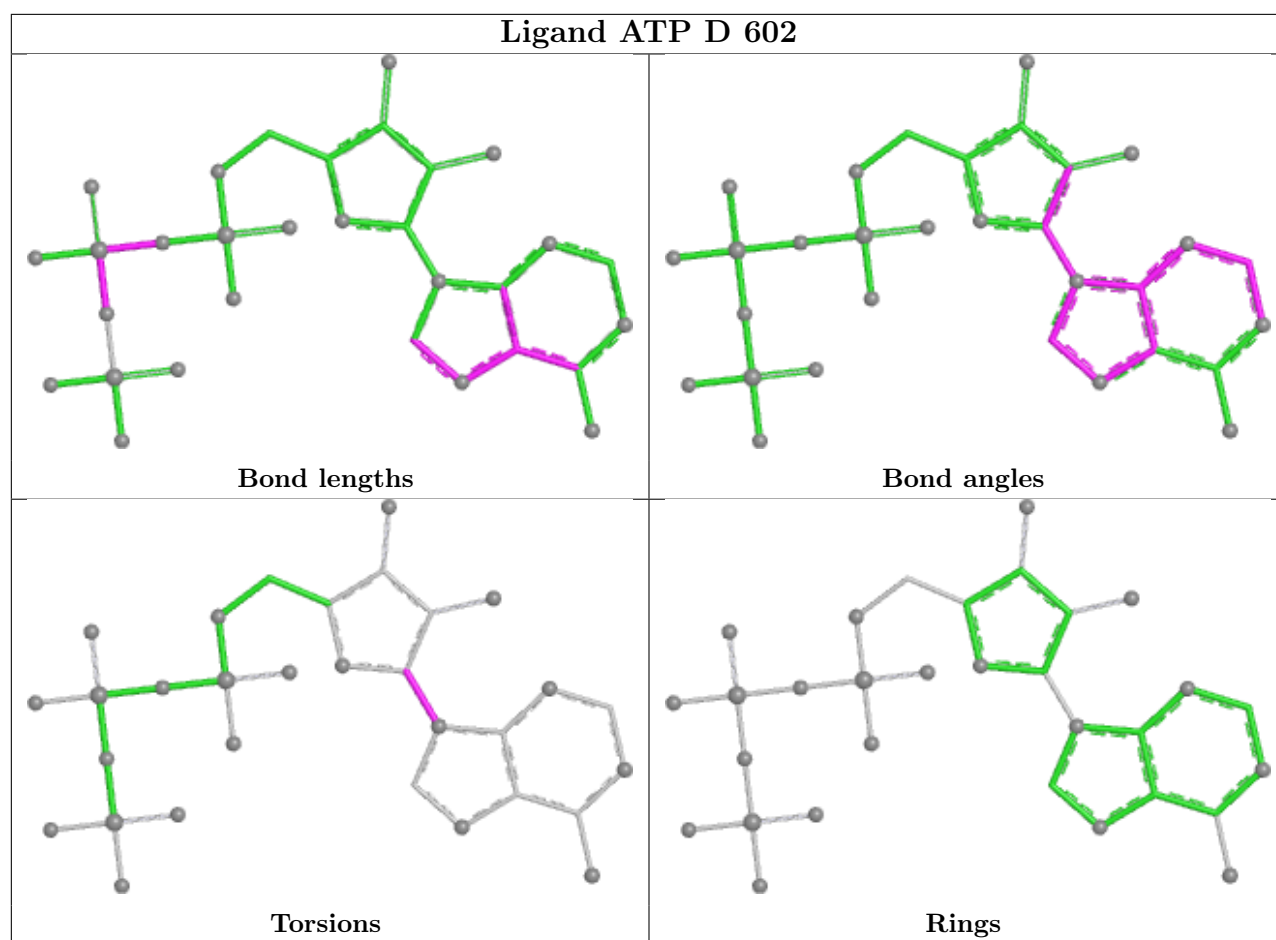
4 monomers are involved in 8 short contacts:

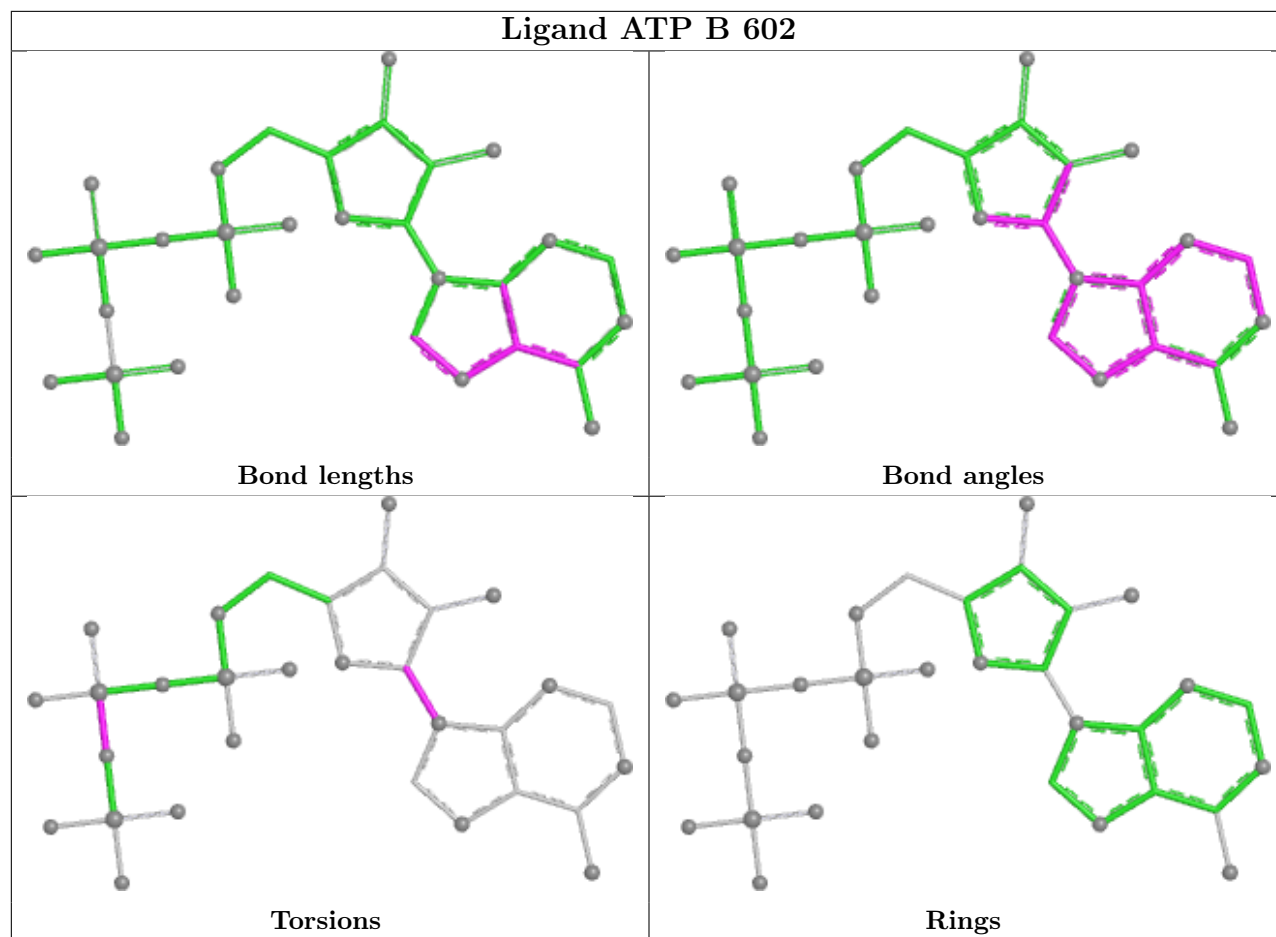
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	602	ATP	1	0
4	A	602	ATP	2	0
4	D	602	ATP	3	0
4	B	602	ATP	2	0

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 ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2			OWAB(Å ²)	Q < 0.9
1	A	493/549 (89%)	0.72	23 (4%)	36	27	55, 109, 155, 194	0
1	B	494/549 (89%)	0.75	40 (8%)	18	16	62, 108, 206, 225	0
1	C	493/549 (89%)	0.78	35 (7%)	22	18	67, 119, 159, 205	0
1	D	494/549 (89%)	0.82	36 (7%)	21	18	62, 122, 203, 247	0
2	E	5/8 (62%)	0.59	1 (20%)	3	4	90, 98, 101, 110	0
2	F	4/8 (50%)	0.93	0	100	100	90, 102, 105, 139	0
2	G	4/8 (50%)	0.20	0	100	100	94, 100, 112, 131	0
2	H	5/8 (62%)	0.94	0	100	100	95, 124, 153, 162	0
All	All	1992/2228 (89%)	0.77	135 (6%)	23	19	55, 116, 184, 247	0

All (135) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	49	THR	5.3
1	B	105	VAL	4.4
1	A	64	GLN	4.4
1	C	49	THR	4.4
1	A	337	PHE	4.2
1	B	116	ASP	3.7
1	A	29	TYR	3.7
1	B	100	LEU	3.7
1	D	410	TRP	3.7
1	C	29	TYR	3.6
1	A	159	GLY	3.5
1	B	51	LEU	3.5
1	A	76	ASN	3.5
1	D	121	LEU	3.3
1	C	487	ILE	3.3
1	A	539	PHE	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	10	LEU	3.3
1	C	470	ALA	3.3
1	D	51	LEU	3.3
1	D	110	PHE	3.2
1	C	65	ALA	3.2
1	A	409	LEU	3.2
1	C	293	LEU	3.2
1	C	238	ALA	3.2
1	A	359	GLY	3.2
1	C	77	TYR	3.2
1	B	106	ILE	3.1
1	B	427	ASP	3.0
1	A	398	VAL	3.0
1	C	417	ASP	3.0
1	B	337	PHE	3.0
1	C	394	ALA	2.9
1	B	75	PHE	2.9
1	D	109	GLN	2.9
1	C	179	LEU	2.9
1	C	66	SER	2.9
1	C	409	LEU	2.9
1	C	32	GLY	2.9
1	C	283	GLN	2.8
1	D	71	LEU	2.8
1	D	214	GLU	2.8
1	B	6	GLU	2.8
1	B	426	LEU	2.7
1	C	532	PHE	2.7
1	C	476	GLY	2.7
1	D	181	GLY	2.7
1	D	61	TYR	2.7
1	A	360	LEU	2.7
1	D	452	LEU	2.7
1	C	381	CYS	2.6
1	D	100	LEU	2.6
1	B	5	ILE	2.6
1	D	381	CYS	2.6
1	B	336	PHE	2.6
1	D	530	PRO	2.6
1	A	32	GLY	2.6
1	D	101	LEU	2.5
1	B	214	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	56	ASN	2.5
1	A	487	ILE	2.5
1	C	539	PHE	2.5
1	C	484	LEU	2.5
1	D	106	ILE	2.5
1	B	101	LEU	2.5
1	D	10	LEU	2.5
1	C	401	PRO	2.5
1	A	476	GLY	2.5
1	B	359	GLY	2.5
1	B	107	ALA	2.4
1	B	302	LEU	2.4
1	D	176	SER	2.4
1	D	414	ILE	2.4
1	B	355	LEU	2.4
1	A	77	TYR	2.4
1	D	456	ILE	2.4
2	E	753	LEU	2.4
1	B	181	GLY	2.4
1	B	115	LEU	2.4
1	A	284	THR	2.3
1	C	524	ILE	2.3
1	D	72	ILE	2.3
1	D	11	SER	2.3
1	B	522	THR	2.3
1	C	34	ASN	2.3
1	D	116	ASP	2.3
1	D	427	ASP	2.3
1	D	283	GLN	2.3
1	B	398	VAL	2.3
1	C	105	VAL	2.3
1	A	302	LEU	2.3
1	B	121	LEU	2.3
1	C	157	LEU	2.3
1	A	475	ASN	2.3
1	C	427	ASP	2.3
1	B	62	ARG	2.3
1	D	341	GLU	2.3
1	D	95	ILE	2.3
1	A	301	PHE	2.3
1	D	166	SER	2.3
1	D	463	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	76	ASN	2.2
1	D	471	GLN	2.2
1	B	217	LEU	2.2
1	B	345	PHE	2.2
1	C	272	THR	2.2
1	B	77	TYR	2.2
1	B	223	VAL	2.2
1	B	341	GLU	2.2
1	D	161	ALA	2.2
1	B	460	ILE	2.2
1	C	35	ILE	2.2
1	A	51	LEU	2.2
1	C	280	ALA	2.2
1	B	272	THR	2.2
1	B	284	THR	2.2
1	D	105	VAL	2.2
1	B	452	LEU	2.2
1	C	279	ASN	2.2
1	D	160	ARG	2.2
1	D	268	CYS	2.2
1	B	481	GLU	2.1
1	C	301	PHE	2.1
1	B	158	HIS	2.1
1	B	464	LEU	2.1
1	A	262	GLN	2.1
1	B	2	LEU	2.1
1	C	91	ILE	2.0
1	D	91	ILE	2.0
1	A	336	PHE	2.0
1	D	328	GLY	2.0
1	B	487	ILE	2.0
1	C	33	LYS	2.0
1	A	224	GLU	2.0
1	C	51	LEU	2.0
1	C	180	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

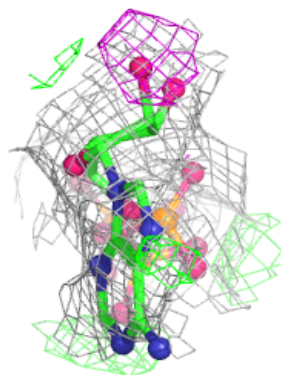
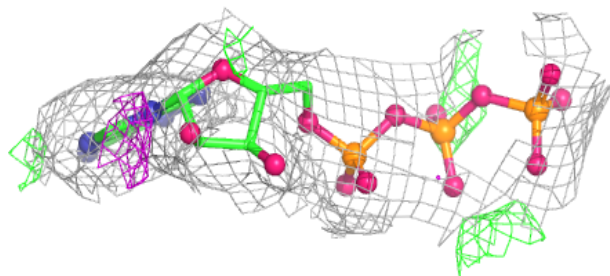
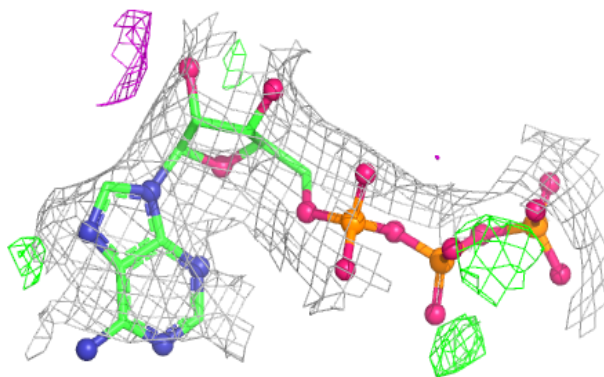
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	601	1/1	0.82	0.15	77,77,77,77	0
4	ATP	D	602	31/31	0.87	0.12	66,91,123,128	0
3	MG	D	601	1/1	0.89	0.23	75,75,75,75	0
4	ATP	A	602	31/31	0.90	0.10	62,79,97,107	0
4	ATP	C	602	31/31	0.91	0.09	68,91,117,119	0
4	ATP	B	602	31/31	0.91	0.10	60,75,88,102	0
3	MG	A	601	1/1	0.94	0.10	60,60,60,60	0
3	MG	B	601	1/1	0.95	0.15	58,58,58,58	0

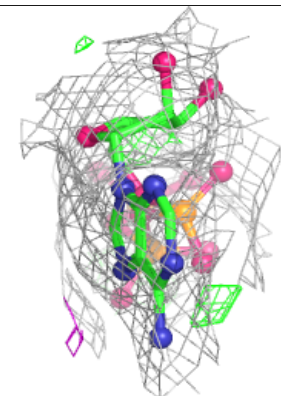
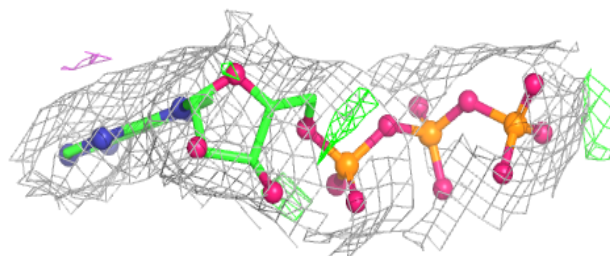
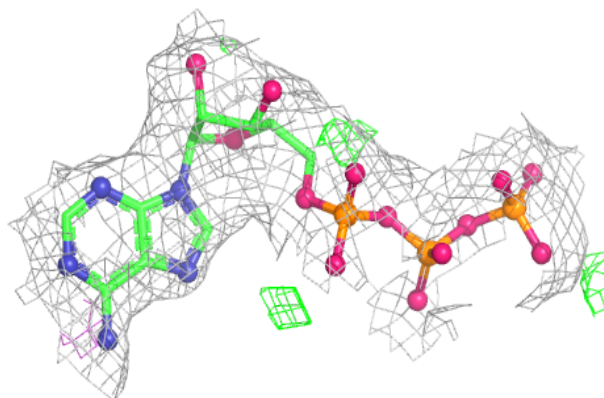
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ATP D 602:

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

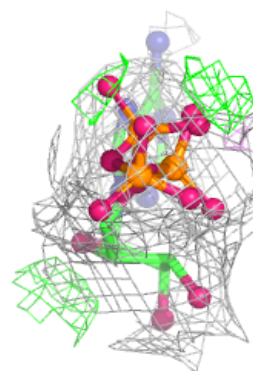
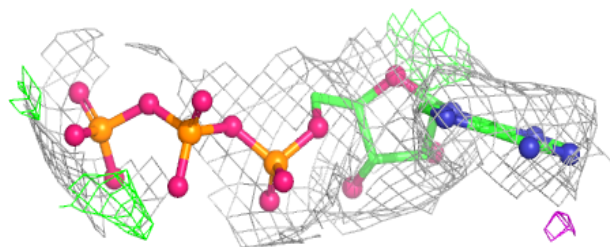
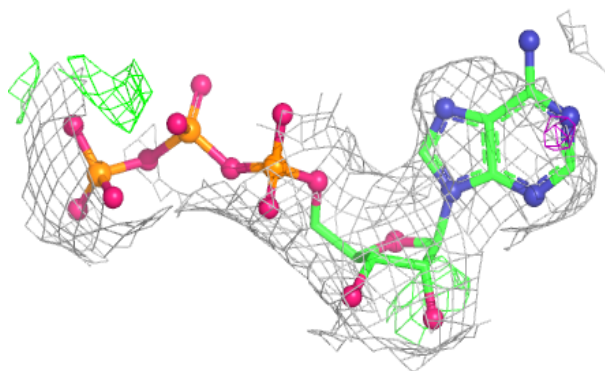
**Electron density around ATP A 602:**

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

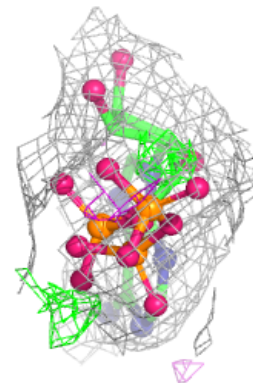
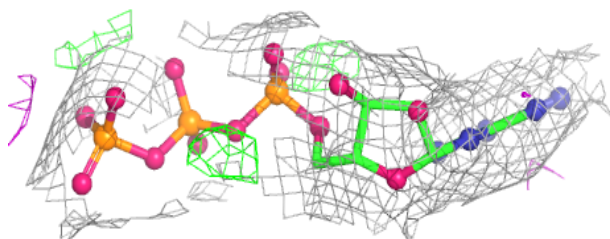
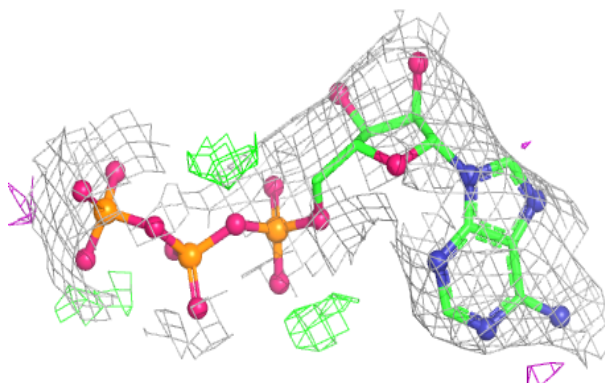


Electron density around ATP C 602:

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

**Electron density around ATP B 602:**

$2mF_o-DF_c$ (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.