



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 04:12 PM UTC

PDB ID : 6KID / pdb_00006kid
Title : Crystal structure of human leucyl-tRNA synthetase, ATP-bound form
Authors : Kim, S.; Son, J.; Kim, S.; Hwang, K.Y.
Deposited on : 2019-07-18
Resolution : 3.15 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

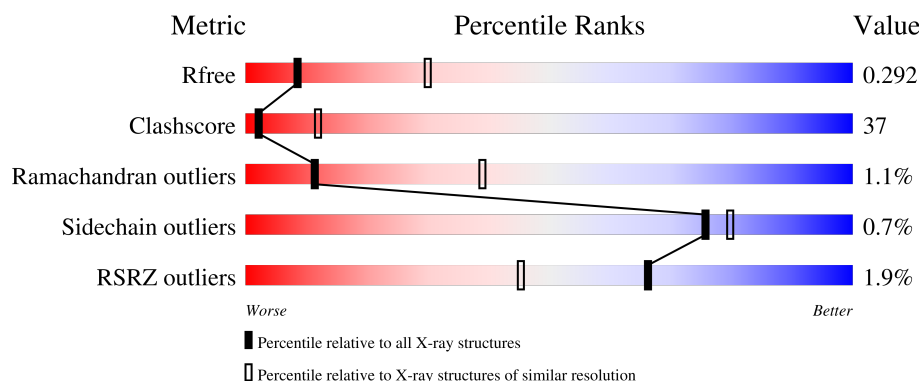
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

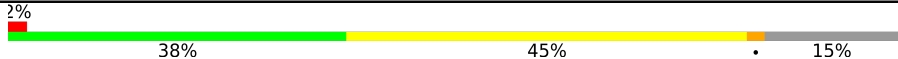
The reported resolution of this entry is 3.15 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2361 (3.20-3.12)
Clashscore	190562	2486 (3.20-3.12)
Ramachandran outliers	187476	2405 (3.20-3.12)
Sidechain outliers	187428	2404 (3.20-3.12)
RSRZ outliers	180081	2361 (3.20-3.12)

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

Mol	Chain	Length	Quality of chain
1	A	1188	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8353 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 Leucine-tRNA ligase, cytoplasmic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1004	Total	C	N	O	S	0	0	0
			8103	5213	1346	1491	53			

There are 12 discrepancies between the modelled and reference sequences:

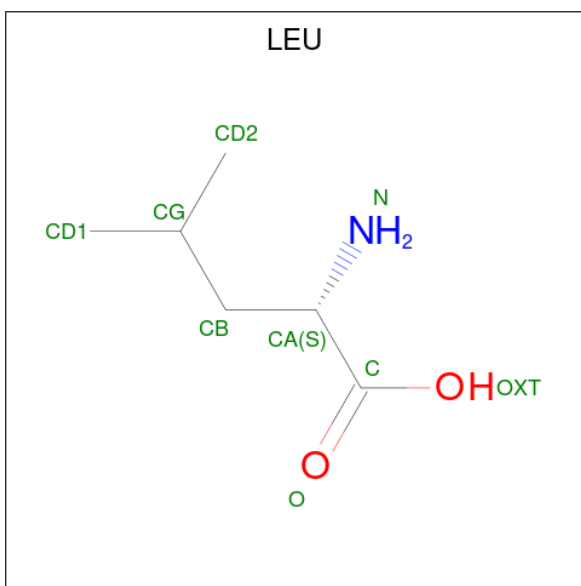
Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	initiating methionine	UNP Q9P2J5
A	-10	ARG	-	expression tag	UNP Q9P2J5
A	-9	GLY	-	expression tag	UNP Q9P2J5
A	-8	SER	-	expression tag	UNP Q9P2J5
A	-7	HIS	-	expression tag	UNP Q9P2J5
A	-6	HIS	-	expression tag	UNP Q9P2J5
A	-5	HIS	-	expression tag	UNP Q9P2J5
A	-4	HIS	-	expression tag	UNP Q9P2J5
A	-3	HIS	-	expression tag	UNP Q9P2J5
A	-2	HIS	-	expression tag	UNP Q9P2J5
A	-1	GLY	-	expression tag	UNP Q9P2J5
A	0	SER	-	expression tag	UNP Q9P2J5

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is LEUCINE (CCD ID: LEU) (formula: $\text{C}_6\text{H}_{13}\text{NO}_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			9	6	1	2		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



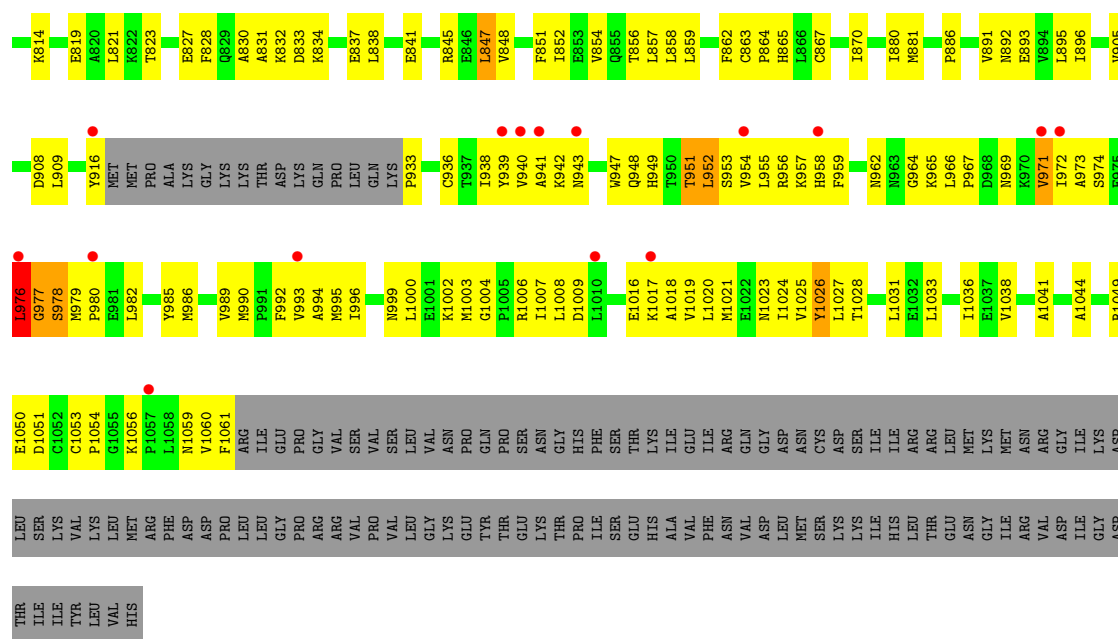
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	200	Total	O	0	0
			200	200		

- Molecule 1: Leucine-tRNA ligase, cytoplasmic





4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	137.22Å 137.22Å 433.32Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.89 – 3.15 42.89 – 3.15	Depositor EDS
% Data completeness (in resolution range)	91.0 (42.89-3.15) 91.2 (42.89-3.15)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.40 (at 2.96Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.245 , 0.293 0.247 , 0.292	Depositor DCC
R_{free} test set	1832 reflections (3.68%)	wwPDB-VP
Wilson B-factor (Å ²)	32.6	Xtriage
Anisotropy	0.345	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 59.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.31$, $\langle L^2 \rangle = 0.15$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.73	EDS
Total number of atoms	8353	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, ATP

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	0.78	2/8303 (0.0%)	1.07	17/11222 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	346	VAL	CA-C	7.73	1.59	1.53
1	A	534	TYR	CA-C	-5.95	1.42	1.52

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1026	TYR	N-CA-C	-10.57	99.78	112.89
1	A	449	LEU	N-CA-C	-9.29	101.21	114.12
1	A	261	PRO	N-CA-C	-8.67	96.90	110.95
1	A	976	LEU	N-CA-C	-7.79	103.39	113.12
1	A	249	MET	CG-SD-CE	7.49	117.38	100.90
1	A	977	GLY	N-CA-C	-7.26	103.86	112.64
1	A	249	MET	CB-CG-SD	6.91	133.44	112.70
1	A	955	LEU	N-CA-C	-6.72	103.14	111.75
1	A	971	VAL	N-CA-C	-5.72	106.82	111.91
1	A	282	LEU	CA-CB-CG	-5.70	96.36	116.30
1	A	847	LEU	CA-CB-CG	5.54	135.69	116.30
1	A	391	SER	CA-C-N	-5.46	112.04	122.13
1	A	391	SER	C-N-CA	-5.46	112.04	122.13
1	A	262	GLN	N-CA-C	5.36	117.64	108.90
1	A	575	ARG	NE-CZ-NH1	-5.24	116.27	121.50
1	A	896	ILE	N-CA-C	-5.08	106.73	111.45
1	A	249	MET	CA-CB-CG	5.00	124.10	114.10

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	8103	0	8062	598	0
2	A	31	0	11	5	0
3	A	9	0	10	0	0
4	A	10	0	0	0	0
5	A	200	0	0	22	0
All	All	8353	0	8083	600	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 37.

All (600) 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:1017:LYS:NZ	1:A:1038:VAL:CG1	1.68	1.55
1:A:1017:LYS:NZ	1:A:1038:VAL:HG12	1.03	1.36
1:A:1017:LYS:CE	1:A:1038:VAL:HG12	1.70	1.19
1:A:629:GLN:HE21	1:A:650:ILE:HD11	0.99	1.14
1:A:629:GLN:NE2	1:A:650:ILE:HD11	1.65	1.11
1:A:1017:LYS:CE	1:A:1038:VAL:CG1	2.29	1.09
1:A:1017:LYS:HZ3	1:A:1038:VAL:CG1	1.42	1.08
1:A:629:GLN:HE21	1:A:650:ILE:CD1	1.66	1.07
1:A:268:LYS:HE3	1:A:288:PHE:CE2	1.92	1.04
1:A:1003:MET:HG2	1:A:1007:ILE:HD11	1.37	1.03
1:A:954:VAL:O	1:A:957:LYS:HB2	1.63	0.96
1:A:940:VAL:HG21	1:A:1017:LYS:CD	1.96	0.95
1:A:9:LYS:HE3	1:A:765:LEU:HD22	1.46	0.95
1:A:419:ARG:H	1:A:422:MET:HE3	1.28	0.94
1:A:1017:LYS:HZ3	1:A:1038:VAL:HG13	1.32	0.94
1:A:629:GLN:NE2	1:A:650:ILE:CD1	2.30	0.91
1:A:267:LEU:HD21	1:A:498:MET:SD	2.11	0.90
1:A:224:THR:HG21	1:A:692:MET:HE3	1.51	0.90
1:A:237:TYR:CZ	1:A:526:LEU:HD22	2.07	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:CYS:SG	1:A:206:ILE:HD12	2.15	0.87
1:A:629:GLN:HG2	1:A:650:ILE:CD1	2.05	0.86
1:A:1017:LYS:HZ2	1:A:1038:VAL:CG1	1.55	0.85
1:A:751:GLU:O	5:A:1301:HOH:O	1.94	0.85
1:A:320:ASN:OD1	5:A:1302:HOH:O	1.95	0.84
1:A:972:ILE:HD12	1:A:990:MET:HG2	1.58	0.84
1:A:1017:LYS:HE3	1:A:1038:VAL:HG11	1.59	0.83
1:A:949:HIS:O	1:A:953:SER:HB2	1.79	0.83
1:A:940:VAL:HG21	1:A:1017:LYS:HD3	1.61	0.82
1:A:1017:LYS:NZ	1:A:1038:VAL:HG13	1.89	0.81
1:A:942:LYS:HG3	1:A:943:ASN:H	1.44	0.81
1:A:976:LEU:HD13	1:A:982:LEU:HD12	1.63	0.80
1:A:434:ILE:HG12	1:A:471:GLY:HA2	1.61	0.80
1:A:195:MET:HE2	1:A:197:LEU:HD21	1.61	0.80
1:A:213:TYR:HD2	1:A:611:GLN:HG2	1.46	0.80
1:A:10:VAL:HG22	1:A:764:ILE:HD13	1.62	0.80
1:A:586:ASP:HB3	1:A:589:TRP:CD1	2.17	0.80
1:A:629:GLN:CG	1:A:650:ILE:HD11	2.12	0.80
1:A:941:ALA:HB3	1:A:1054:PRO:HD3	1.63	0.80
1:A:774:VAL:HG13	1:A:852:ILE:HD13	1.64	0.79
1:A:1017:LYS:CE	1:A:1038:VAL:HG11	2.11	0.78
1:A:1017:LYS:HE3	1:A:1038:VAL:CG1	2.13	0.78
1:A:761:ASP:OD2	5:A:1303:HOH:O	2.01	0.78
1:A:954:VAL:O	1:A:957:LYS:N	2.15	0.78
1:A:1049:ARG:HH11	1:A:1050:GLU:HG3	1.50	0.77
1:A:330:ALA:O	1:A:334:MET:HG2	1.86	0.76
1:A:799:SER:HB3	1:A:892:ASN:H	1.51	0.76
1:A:940:VAL:HG12	1:A:941:ALA:H	1.51	0.76
1:A:770:TRP:HE1	1:A:870:ILE:HD11	1.49	0.76
1:A:837:GLU:HG2	1:A:1026:TYR:OH	1.86	0.76
1:A:569:GLN:HG2	1:A:570:GLU:H	1.51	0.76
1:A:31:GLU:OE1	1:A:202:ARG:NH1	2.19	0.75
1:A:605:THR:O	1:A:661:GLU:HG2	1.87	0.75
1:A:261:PRO:HG2	5:A:1360:HOH:O	1.87	0.74
1:A:977:GLY:HA2	1:A:986:MET:HG3	1.69	0.74
1:A:16:ILE:O	1:A:20:ILE:HG12	1.88	0.73
1:A:538:ASN:ND2	5:A:1307:HOH:O	2.18	0.73
1:A:24:TRP:CH2	1:A:864:PRO:HB2	2.25	0.72
1:A:586:ASP:HB3	1:A:589:TRP:HD1	1.55	0.71
1:A:268:LYS:HE3	1:A:288:PHE:CD2	2.25	0.71
1:A:282:LEU:HD11	1:A:349:VAL:HG21	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:892:ASN:HB3	1:A:895:LEU:HB2	1.73	0.71
2:A:1201:ATP:H5'1	2:A:1201:ATP:O1B	1.91	0.71
1:A:629:GLN:CD	1:A:650:ILE:HD11	2.16	0.70
1:A:990:MET:HA	1:A:993:VAL:HB	1.74	0.70
1:A:602:ALA:O	1:A:605:THR:OG1	2.09	0.69
1:A:645:PHE:HZ	1:A:652:LYS:HA	1.56	0.69
1:A:940:VAL:CG2	1:A:1017:LYS:NZ	2.56	0.69
1:A:419:ARG:N	1:A:422:MET:HE3	2.04	0.69
1:A:287:ILE:HG23	1:A:323:ILE:HB	1.73	0.69
1:A:629:GLN:CG	1:A:650:ILE:CD1	2.71	0.69
1:A:947:TRP:HZ3	1:A:989:VAL:HA	1.58	0.69
1:A:471:GLY:O	1:A:487:VAL:HG21	1.94	0.68
1:A:366:LEU:HD21	1:A:498:MET:HG3	1.75	0.68
1:A:60:HIS:HA	1:A:726:THR:HA	1.75	0.68
1:A:439:ASN:HB2	1:A:440:LEU:HG	1.77	0.67
1:A:200:ASP:OD2	1:A:202:ARG:NH2	2.27	0.67
1:A:203:ARG:HG3	1:A:608:HIS:HB3	1.76	0.67
1:A:335:SER:HA	1:A:340:THR:HG23	1.76	0.67
1:A:949:HIS:O	1:A:949:HIS:ND1	2.28	0.67
1:A:167:SER:O	1:A:171:ILE:HG13	1.94	0.67
1:A:10:VAL:HG22	1:A:764:ILE:CD1	2.25	0.66
1:A:573:CYS:HA	1:A:592:GLU:HG2	1.77	0.66
1:A:614:ASN:OD1	1:A:615:LEU:N	2.27	0.66
1:A:1003:MET:HE3	1:A:1007:ILE:HG13	1.75	0.66
1:A:1056:LYS:NZ	5:A:1312:HOH:O	2.28	0.66
1:A:259:VAL:HG13	1:A:511:GLU:HB2	1.76	0.66
1:A:423:VAL:HG23	1:A:424:LEU:HG	1.78	0.66
1:A:940:VAL:HG21	1:A:1017:LYS:HD2	1.77	0.66
1:A:530:TRP:H	1:A:574:SER:HB3	1.61	0.66
1:A:736:ALA:O	1:A:740:ARG:N	2.26	0.66
1:A:1000:LEU:HD22	1:A:1004:GLY:O	1.96	0.66
1:A:55:MET:SD	1:A:90:LEU:HD22	2.37	0.65
1:A:237:TYR:CE2	1:A:526:LEU:HD22	2.31	0.65
1:A:733:LYS:NZ	1:A:757:GLU:OE2	2.25	0.65
1:A:10:VAL:CG2	1:A:764:ILE:HD13	2.26	0.65
1:A:1049:ARG:NH1	1:A:1050:GLU:HG3	2.11	0.65
1:A:14:LYS:HA	1:A:17:GLU:HB2	1.78	0.65
1:A:267:LEU:HD13	1:A:299:MET:HE2	1.79	0.65
1:A:859:LEU:HB3	1:A:867:CYS:SG	2.38	0.64
1:A:282:LEU:CD1	1:A:349:VAL:HG21	2.28	0.64
1:A:1053:CYS:HB2	1:A:1054:PRO:HD2	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:893:GLU:OE2	5:A:1304:HOH:O	2.14	0.64
1:A:409:LYS:HB2	1:A:412:LEU:HD13	1.80	0.64
1:A:965:LYS:HD3	1:A:966:LEU:O	1.97	0.64
1:A:969:ASN:HA	1:A:972:ILE:HG23	1.78	0.64
1:A:60:HIS:O	1:A:63:HIS:HB2	1.98	0.64
1:A:977:GLY:HA2	1:A:986:MET:CG	2.28	0.63
1:A:22:GLN:O	1:A:26:THR:HG22	1.99	0.63
1:A:601:MET:O	1:A:604:TYR:HB2	1.98	0.63
1:A:280:SER:O	1:A:283:LYS:HE3	1.97	0.63
1:A:673:SER:OG	1:A:674:GLY:N	2.32	0.63
1:A:78:ARG:HD2	1:A:881:MET:HE2	1.79	0.63
1:A:12:PHE:HB3	5:A:1305:HOH:O	1.99	0.62
1:A:61:LEU:O	1:A:64:THR:N	2.29	0.62
1:A:947:TRP:CE3	1:A:989:VAL:HG12	2.34	0.62
1:A:77:GLN:O	1:A:80:LYS:N	2.33	0.62
1:A:194:ARG:O	1:A:731:ILE:HD13	1.98	0.62
1:A:940:VAL:CG2	1:A:1017:LYS:HD3	2.28	0.62
1:A:158:TRP:CH2	1:A:171:ILE:HD12	2.35	0.62
1:A:185:PRO:HG3	1:A:206:ILE:HD11	1.81	0.62
1:A:256:GLY:O	1:A:259:VAL:HG23	1.98	0.62
1:A:271:VAL:HG23	1:A:287:ILE:O	1.99	0.62
1:A:648:THR:HG21	1:A:655:LEU:HD12	1.82	0.62
1:A:413:ARG:HD3	1:A:423:VAL:HG21	1.82	0.62
1:A:267:LEU:HD11	1:A:299:MET:HG3	1.82	0.62
1:A:995:MET:SD	1:A:996:ILE:HG12	2.40	0.62
1:A:282:LEU:HD22	1:A:285:LYS:HE3	1.82	0.61
1:A:9:LYS:O	5:A:1305:HOH:O	2.16	0.61
1:A:314:ILE:HG23	1:A:348:VAL:HG13	1.82	0.61
1:A:629:GLN:HE21	1:A:650:ILE:CG1	2.12	0.61
1:A:948:GLN:HA	1:A:951:THR:OG1	2.01	0.61
1:A:93:THR:HG21	1:A:596:ASP:OD2	2.01	0.61
1:A:916:TYR:HE1	1:A:1061:PHE:HD2	1.46	0.61
1:A:954:VAL:C	1:A:957:LYS:H	2.07	0.61
1:A:650:ILE:HD12	1:A:650:ILE:N	2.16	0.61
1:A:199:VAL:HB	1:A:201:TRP:CD1	2.36	0.61
1:A:557:THR:HG22	1:A:561:PHE:HE1	1.66	0.60
1:A:939:TYR:HB3	1:A:1041:ALA:HB2	1.83	0.60
1:A:423:VAL:HG23	1:A:424:LEU:N	2.16	0.60
1:A:39:LYS:O	5:A:1306:HOH:O	2.17	0.60
1:A:940:VAL:HG23	1:A:1017:LYS:NZ	2.16	0.60
1:A:1024:ILE:HA	1:A:1027:LEU:HD12	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:569:GLN:HG2	1:A:570:GLU:N	2.15	0.60
1:A:631:THR:O	1:A:634:VAL:N	2.34	0.60
1:A:982:LEU:O	1:A:986:MET:HG2	2.01	0.60
1:A:213:TYR:CD2	1:A:611:GLN:HG2	2.33	0.60
1:A:267:LEU:CD1	1:A:299:MET:HG3	2.31	0.60
1:A:281:GLY:O	1:A:283:LYS:N	2.31	0.60
1:A:316:PHE:HB2	1:A:347:PRO:HD2	1.83	0.60
1:A:190:GLN:O	1:A:193:LYS:HB3	2.01	0.60
1:A:564:THR:HG21	1:A:679:PRO:HA	1.84	0.60
1:A:737:ASP:OD1	1:A:865:HIS:ND1	2.26	0.60
1:A:67:LEU:HD11	1:A:195:MET:HE1	1.84	0.59
1:A:268:LYS:CE	1:A:288:PHE:CE2	2.80	0.59
1:A:558:ARG:O	1:A:562:GLU:HB2	2.03	0.59
1:A:583:LEU:HD12	1:A:584:PRO:HD2	1.84	0.59
1:A:263:GLU:HA	1:A:508:MET:HG2	1.84	0.59
1:A:299:MET:HE1	1:A:390:THR:HB	1.84	0.59
1:A:559:ARG:NH1	5:A:1318:HOH:O	2.34	0.59
1:A:626:ARG:NH2	5:A:1321:HOH:O	2.35	0.59
1:A:972:ILE:HG13	1:A:973:ALA:H	1.67	0.59
1:A:770:TRP:HZ3	1:A:851:PHE:CE2	2.21	0.59
1:A:214:TYR:O	1:A:218:VAL:HG23	2.03	0.58
1:A:266:LEU:O	1:A:504:ALA:HB1	2.03	0.58
1:A:543:THR:HG21	1:A:687:TYR:CD1	2.38	0.58
1:A:731:ILE:HG23	1:A:736:ALA:HB2	1.85	0.58
1:A:606:VAL:HG11	1:A:658:LEU:HB3	1.85	0.58
1:A:1023:ASN:O	1:A:1027:LEU:HG	2.03	0.58
1:A:64:THR:HG21	1:A:727:LEU:HD13	1.85	0.58
1:A:266:LEU:CG	1:A:290:VAL:HG13	2.33	0.58
1:A:539:TRP:O	1:A:543:THR:HG23	2.03	0.58
1:A:24:TRP:HH2	1:A:864:PRO:HB2	1.66	0.58
1:A:1006:ARG:O	1:A:1008:LEU:N	2.36	0.58
1:A:9:LYS:HB3	1:A:764:ILE:HD12	1.86	0.58
1:A:942:LYS:HG3	1:A:943:ASN:N	2.18	0.58
1:A:262:GLN:HG3	1:A:511:GLU:HG2	1.84	0.58
1:A:278:LYS:O	1:A:279:LEU:HD23	2.03	0.58
1:A:776:GLU:HG2	1:A:777:MET:HE2	1.86	0.58
1:A:959:PHE:HA	1:A:962:ASN:HB2	1.86	0.58
1:A:543:THR:HG21	1:A:687:TYR:HD1	1.70	0.57
1:A:458:GLU:HG3	1:A:459:LYS:HG3	1.85	0.57
1:A:21:GLN:OE1	1:A:194:ARG:HA	2.04	0.57
1:A:571:HIS:NE2	1:A:592:GLU:OE2	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:821:LEU:HD12	1:A:862:PHE:CD2	2.39	0.57
1:A:21:GLN:CD	1:A:194:ARG:HA	2.29	0.57
1:A:10:VAL:HG11	1:A:734:PHE:HA	1.86	0.57
1:A:85:LEU:HG	1:A:87:PRO:HD3	1.85	0.56
1:A:675:LYS:HD2	1:A:708:GLY:HA2	1.87	0.56
1:A:189:ILE:HD11	1:A:201:TRP:CD1	2.40	0.56
1:A:266:LEU:HG	1:A:290:VAL:HG13	1.86	0.56
1:A:976:LEU:HD13	1:A:982:LEU:CD1	2.35	0.56
1:A:1017:LYS:HZ2	1:A:1038:VAL:HG12	0.73	0.56
1:A:78:ARG:HD2	1:A:881:MET:CE	2.36	0.56
1:A:637:TYR:HA	1:A:643:ALA:HB3	1.86	0.56
1:A:798:ALA:HB2	1:A:847:LEU:HD11	1.88	0.56
1:A:972:ILE:C	1:A:974:SER:H	2.14	0.56
1:A:13:LEU:O	1:A:16:ILE:HG12	2.06	0.56
1:A:93:THR:HG23	1:A:207:THR:HB	1.86	0.56
1:A:308:ARG:HD3	1:A:310:ASP:OD1	2.06	0.56
1:A:397:PRO:HA	1:A:400:ILE:CG2	2.36	0.56
1:A:971:VAL:HA	1:A:974:SER:HB3	1.86	0.56
1:A:821:LEU:HB2	1:A:862:PHE:CE2	2.41	0.56
1:A:954:VAL:O	1:A:957:LYS:CB	2.46	0.56
1:A:242:PRO:HG2	1:A:332:ARG:NH2	2.20	0.56
1:A:418:ILE:HG22	1:A:422:MET:HB2	1.87	0.56
1:A:437:PHE:HE2	1:A:470:LYS:HD3	1.69	0.56
1:A:881:MET:HA	1:A:881:MET:HE3	1.87	0.56
1:A:269:LEU:HD12	1:A:289:LEU:O	2.07	0.55
1:A:905:VAL:O	1:A:908:ASP:HB3	2.06	0.55
1:A:449:LEU:CD1	1:A:463:ALA:HB2	2.37	0.55
1:A:947:TRP:CZ3	1:A:989:VAL:HG12	2.40	0.55
1:A:1025:VAL:HA	1:A:1028:THR:OG1	2.07	0.55
1:A:1017:LYS:O	1:A:1020:LEU:N	2.31	0.55
1:A:58:ARG:NH1	1:A:728:THR:HG21	2.23	0.55
1:A:192:LEU:HD22	1:A:197:LEU:HD12	1.90	0.54
1:A:669:ASP:HA	1:A:702:THR:OG1	2.07	0.54
1:A:103:LYS:HD3	1:A:157:GLN:HB2	1.88	0.54
1:A:266:LEU:HD21	1:A:290:VAL:CG1	2.38	0.54
1:A:397:PRO:HA	1:A:400:ILE:HG22	1.90	0.54
1:A:586:ASP:O	1:A:588:GLN:N	2.41	0.54
1:A:50:PHE:CD1	1:A:51:PRO:HD2	2.43	0.54
1:A:71:GLU:HG3	1:A:740:ARG:NH2	2.23	0.54
1:A:482:PHE:O	1:A:483:LYS:C	2.49	0.54
1:A:854:VAL:O	1:A:858:LEU:HG	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:793:ASN:ND2	1:A:841:GLU:OE2	2.40	0.53
1:A:845:ARG:O	1:A:848:VAL:N	2.32	0.53
1:A:947:TRP:NE1	1:A:1051:ASP:OD1	2.41	0.53
1:A:597:SER:O	1:A:681:HIS:HE1	1.91	0.53
1:A:770:TRP:HE1	1:A:870:ILE:CD1	2.18	0.53
1:A:250:ASP:OD2	1:A:680:ASN:HB3	2.08	0.53
1:A:383:ASP:OD1	1:A:384:LYS:HG2	2.08	0.53
1:A:773:TRP:CD2	1:A:832:LYS:HE2	2.43	0.53
1:A:804:GLY:O	1:A:805:ILE:C	2.51	0.53
1:A:940:VAL:HG21	1:A:1017:LYS:NZ	2.23	0.53
1:A:52:TYR:HE1	1:A:91:HIS:CE1	2.26	0.53
1:A:583:LEU:HD11	1:A:585:TRP:CZ2	2.44	0.53
1:A:761:ASP:HA	1:A:764:ILE:HG12	1.91	0.53
1:A:940:VAL:HG11	1:A:1016:GLU:HG2	1.90	0.53
1:A:432:ILE:HD11	1:A:477:MET:SD	2.49	0.53
1:A:678:VAL:HB	1:A:679:PRO:HD3	1.91	0.53
1:A:600:TYR:HE1	1:A:604:TYR:CE2	2.27	0.52
1:A:717:MET:HA	1:A:723:ASN:O	2.09	0.52
1:A:333:ASN:O	1:A:334:MET:C	2.52	0.52
1:A:770:TRP:HZ3	1:A:851:PHE:HE2	1.56	0.52
1:A:305:CYS:HB3	1:A:374:VAL:HA	1.92	0.52
1:A:354:GLY:HA2	1:A:357:ILE:HD12	1.90	0.52
1:A:594:LEU:HD13	1:A:684:TYR:HE2	1.75	0.52
1:A:785:ARG:HB3	1:A:785:ARG:CZ	2.39	0.52
1:A:412:LEU:C	1:A:414:ALA:H	2.16	0.52
1:A:545:GLN:NE2	5:A:1325:HOH:O	2.43	0.52
1:A:65:PHE:HB2	1:A:710:LEU:HD22	1.91	0.52
1:A:221:GLN:HG3	1:A:692:MET:SD	2.49	0.52
1:A:601:MET:CE	1:A:685:TYR:HE1	2.23	0.52
1:A:713:ASN:HD21	1:A:757:GLU:CD	2.18	0.52
1:A:956:ARG:O	1:A:959:PHE:N	2.41	0.51
1:A:967:PRO:HB2	1:A:971:VAL:HG11	1.92	0.51
1:A:985:TYR:O	1:A:989:VAL:HG13	2.09	0.51
1:A:480:ASP:OD1	1:A:483:LYS:HE2	2.10	0.51
1:A:335:SER:O	1:A:343:ASN:ND2	2.43	0.51
1:A:933:PRO:HG3	1:A:1061:PHE:CD1	2.46	0.51
1:A:793:ASN:HA	1:A:796:VAL:HB	1.91	0.51
1:A:263:GLU:HG3	1:A:264:TYR:N	2.24	0.51
1:A:70:CYS:O	1:A:74:VAL:HG22	2.10	0.51
1:A:92:CYS:O	1:A:97:ILE:HD13	2.11	0.51
1:A:978:SER:O	1:A:978:SER:OG	2.16	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:990:MET:O	1:A:994:ALA:N	2.43	0.51
1:A:856:THR:OG1	1:A:870:ILE:HG21	2.10	0.51
1:A:938:ILE:HD11	1:A:1036:ILE:HG21	1.93	0.51
1:A:966:LEU:HD12	1:A:966:LEU:H	1.76	0.51
1:A:52:TYR:CE1	1:A:91:HIS:CG	2.99	0.51
1:A:266:LEU:HD11	1:A:290:VAL:CG1	2.41	0.51
1:A:979:MET:HG3	1:A:980:PRO:HD2	1.91	0.51
1:A:282:LEU:HA	5:A:1346:HOH:O	2.10	0.50
1:A:396:SER:HB2	1:A:468:TYR:HE1	1.76	0.50
1:A:48:VAL:HG21	1:A:70:CYS:HB2	1.94	0.50
1:A:332:ARG:O	1:A:335:SER:OG	2.30	0.50
1:A:756:VAL:O	1:A:759:MET:N	2.42	0.50
1:A:629:GLN:CD	1:A:650:ILE:CD1	2.82	0.50
1:A:947:TRP:CH2	1:A:992:PHE:HD2	2.30	0.50
1:A:1025:VAL:HG22	1:A:1025:VAL:O	2.11	0.50
1:A:106:ARG:NH2	5:A:1327:HOH:O	2.45	0.50
1:A:288:PHE:N	1:A:323:ILE:O	2.30	0.50
1:A:111:TYR:HB3	1:A:116:ASP:H	1.76	0.49
1:A:201:TRP:NE1	5:A:1322:HOH:O	2.36	0.49
1:A:253:ARG:HD3	1:A:514:VAL:HG11	1.94	0.49
1:A:334:MET:HB3	1:A:339:PHE:CD2	2.47	0.49
1:A:763:GLY:O	1:A:765:LEU:N	2.44	0.49
1:A:266:LEU:HD21	1:A:290:VAL:HG11	1.93	0.49
1:A:831:ALA:O	1:A:832:LYS:C	2.55	0.49
1:A:199:VAL:CG1	1:A:201:TRP:HD1	2.24	0.49
1:A:568:LEU:HD21	1:A:684:TYR:HE1	1.78	0.49
1:A:293:THR:HB	1:A:390:THR:HG21	1.94	0.49
1:A:336:TYR:HA	1:A:343:ASN:HD21	1.77	0.49
1:A:568:LEU:HD21	1:A:684:TYR:CE1	2.47	0.49
1:A:65:PHE:CB	1:A:710:LEU:HD22	2.43	0.49
1:A:674:GLY:N	2:A:1201:ATP:O2'	2.42	0.49
1:A:275:TYR:CE2	1:A:287:ILE:HD12	2.48	0.49
1:A:24:TRP:HD1	1:A:29:VAL:HG21	1.77	0.49
1:A:569:GLN:O	1:A:570:GLU:C	2.55	0.49
1:A:819:GLU:O	1:A:823:THR:HG23	2.12	0.49
1:A:448:GLU:O	1:A:448:GLU:HG3	2.12	0.49
1:A:804:GLY:HA3	1:A:828:PHE:HE1	1.78	0.49
1:A:977:GLY:C	1:A:979:MET:H	2.21	0.49
1:A:857:LEU:HD13	1:A:886:PRO:HD3	1.94	0.48
1:A:1027:LEU:O	1:A:1031:LEU:HB2	2.13	0.48
1:A:999:ASN:O	1:A:1003:MET:HB3	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1017:LYS:HZ3	1:A:1038:VAL:HG11	1.62	0.48
1:A:53:PRO:HA	2:A:1201:ATP:O5'	2.13	0.48
1:A:103:LYS:HE2	1:A:107:GLU:OE2	2.13	0.48
1:A:569:GLN:HG2	1:A:570:GLU:HG3	1.93	0.48
1:A:967:PRO:HB2	1:A:971:VAL:CG1	2.43	0.48
2:A:1201:ATP:H3'	2:A:1201:ATP:O3A	2.14	0.48
1:A:265:THR:HG21	1:A:296:PRO:HB3	1.96	0.48
1:A:421:ASP:O	1:A:425:PRO:HD2	2.14	0.48
1:A:867:CYS:HB3	1:A:880:ILE:HD13	1.95	0.48
1:A:486:LYS:O	1:A:490:VAL:HG22	2.12	0.48
1:A:999:ASN:HA	1:A:1002:LYS:HB2	1.95	0.48
1:A:75:GLY:HA2	1:A:198:LYS:HE2	1.95	0.48
1:A:209:ASP:N	1:A:580:GLY:O	2.47	0.48
1:A:495:GLN:O	1:A:499:ILE:HG13	2.13	0.48
1:A:158:TRP:CZ3	1:A:171:ILE:HB	2.49	0.48
1:A:329:LYS:NZ	5:A:1330:HOH:O	2.46	0.48
1:A:608:HIS:HD2	5:A:1326:HOH:O	1.95	0.48
1:A:726:THR:HG22	1:A:729:GLN:HB2	1.95	0.48
1:A:763:GLY:C	1:A:765:LEU:N	2.71	0.48
1:A:267:LEU:CD1	1:A:299:MET:HE2	2.43	0.48
1:A:279:LEU:HD23	1:A:279:LEU:HA	1.42	0.48
1:A:821:LEU:HB2	1:A:862:PHE:HE2	1.77	0.48
1:A:211:ASN:O	1:A:215:ASP:HB2	2.14	0.48
1:A:366:LEU:HD21	1:A:498:MET:CG	2.43	0.48
1:A:635:TRP:O	1:A:638:VAL:N	2.47	0.48
1:A:311:MET:HB3	1:A:313:TYR:CE1	2.49	0.47
1:A:969:ASN:C	1:A:972:ILE:HG12	2.39	0.47
1:A:92:CYS:N	1:A:207:THR:OG1	2.42	0.47
1:A:693:TRP:HB3	1:A:696:GLN:HG3	1.95	0.47
1:A:312:LYS:HE2	1:A:353:MET:HE2	1.97	0.47
1:A:951:THR:HA	1:A:976:LEU:HD11	1.97	0.47
1:A:61:LEU:HD21	1:A:755:PHE:CD2	2.50	0.47
1:A:433:GLU:HB3	1:A:478:LEU:HD11	1.96	0.47
1:A:78:ARG:HG2	1:A:78:ARG:HH11	1.80	0.47
1:A:107:GLU:OE2	1:A:156:TYR:HD1	1.97	0.47
1:A:192:LEU:HD22	1:A:197:LEU:CD1	2.44	0.47
1:A:324:PHE:O	1:A:325:ILE:HG12	2.15	0.47
1:A:674:GLY:H	2:A:1201:ATP:HO2'	1.59	0.47
1:A:940:VAL:CG1	1:A:1016:GLU:HG2	2.45	0.47
1:A:543:THR:HA	1:A:690:VAL:HG21	1.97	0.47
1:A:675:LYS:HE2	5:A:1442:HOH:O	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:892:ASN:O	1:A:895:LEU:N	2.47	0.47
1:A:18:LYS:HE3	1:A:194:ARG:HH11	1.80	0.47
1:A:284:GLY:O	1:A:285:LYS:HG3	2.14	0.47
1:A:976:LEU:O	1:A:979:MET:N	2.47	0.47
1:A:1021:MET:SD	1:A:1038:VAL:HG21	2.54	0.47
1:A:195:MET:CE	1:A:197:LEU:HD21	2.39	0.47
1:A:250:ASP:O	1:A:253:ARG:HG3	2.15	0.47
1:A:299:MET:HE1	1:A:302:GLN:OE1	2.15	0.47
1:A:610:LEU:CD2	1:A:630:MET:HE1	2.45	0.47
1:A:50:PHE:HD1	1:A:51:PRO:HD2	1.79	0.46
1:A:916:TYR:HE1	1:A:1061:PHE:CD2	2.31	0.46
1:A:158:TRP:CZ2	1:A:171:ILE:HD12	2.50	0.46
1:A:353:MET:O	1:A:355:GLU:N	2.48	0.46
1:A:977:GLY:C	1:A:979:MET:N	2.74	0.46
1:A:272:LEU:HD23	1:A:272:LEU:HA	1.34	0.46
1:A:405:ASP:HA	1:A:408:LYS:HB2	1.98	0.46
1:A:541:LYS:O	1:A:541:LYS:HG3	2.15	0.46
1:A:1024:ILE:HD12	1:A:1025:VAL:N	2.30	0.46
1:A:253:ARG:HD3	1:A:514:VAL:CG1	2.46	0.46
1:A:629:GLN:HG2	1:A:650:ILE:HD12	1.91	0.46
1:A:940:VAL:HG12	1:A:941:ALA:N	2.24	0.46
1:A:58:ARG:HD3	1:A:187:LEU:HD13	1.98	0.46
1:A:362:LEU:HD22	1:A:374:VAL:HG22	1.96	0.46
1:A:449:LEU:HD12	1:A:463:ALA:HB2	1.96	0.46
1:A:304:ASN:ND2	1:A:391:SER:OG	2.43	0.46
1:A:1041:ALA:HA	1:A:1044:ALA:HB2	1.98	0.46
1:A:225:LEU:HB2	1:A:231:ILE:HD11	1.98	0.46
1:A:300:PHE:N	1:A:300:PHE:CD1	2.83	0.46
1:A:318:THR:HG21	1:A:320:ASN:OD1	2.16	0.46
1:A:379:THR:O	1:A:379:THR:OG1	2.34	0.46
1:A:749:THR:HB	1:A:751:GLU:H	1.80	0.46
1:A:947:TRP:O	1:A:951:THR:HG23	2.16	0.46
1:A:24:TRP:HZ3	1:A:865:HIS:NE2	2.14	0.46
1:A:476:ILE:HG13	1:A:476:ILE:O	2.15	0.46
1:A:606:VAL:HG12	1:A:658:LEU:HD22	1.98	0.46
1:A:730:ALA:O	1:A:733:LYS:N	2.47	0.46
1:A:773:TRP:CE2	1:A:832:LYS:HE2	2.50	0.46
1:A:510:PRO:HD3	1:A:524:VAL:HG23	1.97	0.46
1:A:601:MET:HE3	1:A:685:TYR:HE1	1.80	0.46
1:A:951:THR:OG1	1:A:952:LEU:N	2.49	0.46
1:A:278:LYS:HE2	1:A:350:LYS:NZ	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:916:TYR:CE1	1:A:1061:PHE:HD2	2.30	0.45
1:A:45:LYS:HD3	1:A:664:PHE:O	2.16	0.45
1:A:262:GLN:C	1:A:508:MET:HE3	2.40	0.45
1:A:476:ILE:HA	1:A:486:LYS:HA	1.98	0.45
1:A:630:MET:HB2	1:A:630:MET:HE2	1.62	0.45
1:A:678:VAL:O	1:A:683:SER:OG	2.35	0.45
1:A:221:GLN:OE1	1:A:599:ILE:HG12	2.16	0.45
1:A:403:LEU:O	1:A:403:LEU:HD12	2.17	0.45
1:A:774:VAL:HG13	1:A:852:ILE:CD1	2.41	0.45
1:A:972:ILE:HG13	1:A:973:ALA:N	2.30	0.45
1:A:71:GLU:HG2	1:A:197:LEU:HD22	1.99	0.45
1:A:579:LEU:HA	1:A:579:LEU:HD23	1.40	0.45
1:A:599:ILE:CG2	1:A:688:ASN:HB3	2.46	0.45
1:A:1007:ILE:N	1:A:1007:ILE:HD12	2.32	0.45
1:A:55:MET:HE2	1:A:57:GLY:O	2.17	0.45
1:A:67:LEU:HD11	1:A:195:MET:CE	2.47	0.45
1:A:289:LEU:HD23	1:A:289:LEU:HA	1.52	0.45
1:A:947:TRP:CZ3	1:A:989:VAL:HA	2.44	0.45
1:A:230:LYS:HG3	1:A:691:ALA:HB1	1.99	0.45
1:A:506:ILE:HD13	1:A:506:ILE:HA	1.68	0.45
1:A:675:LYS:HB3	1:A:707:ASN:O	2.16	0.45
1:A:698:ASP:OD1	1:A:698:ASP:N	2.46	0.45
1:A:772:GLU:OE2	1:A:775:LYS:HD2	2.16	0.45
1:A:1019:VAL:HG13	1:A:1020:LEU:CD2	2.46	0.45
1:A:71:GLU:OE2	1:A:198:LYS:HG3	2.16	0.45
1:A:833:ASP:O	5:A:1308:HOH:O	2.21	0.45
1:A:1008:LEU:HD12	1:A:1008:LEU:HA	1.58	0.45
1:A:8:ALA:HB1	5:A:1422:HOH:O	2.17	0.45
1:A:95:MET:HE3	1:A:95:MET:HB3	1.65	0.45
1:A:456:ASP:O	1:A:460:LEU:HB2	2.17	0.45
1:A:726:THR:HG22	1:A:729:GLN:CG	2.47	0.45
1:A:795:ARG:NH2	1:A:895:LEU:HD11	2.32	0.45
1:A:892:ASN:HD22	1:A:895:LEU:HD12	1.82	0.45
1:A:601:MET:SD	1:A:685:TYR:HE1	2.40	0.45
1:A:610:LEU:HD21	1:A:630:MET:HE1	1.99	0.45
1:A:490:VAL:O	1:A:491:LYS:C	2.58	0.44
1:A:539:TRP:CE2	1:A:691:ALA:HB2	2.51	0.44
1:A:550:LEU:CD2	1:A:552:THR:HG22	2.47	0.44
1:A:55:MET:HE1	1:A:188:ALA:HB2	1.99	0.44
1:A:273:GLU:HA	1:A:274:PRO:C	2.41	0.44
1:A:316:PHE:O	1:A:323:ILE:HA	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:487:VAL:H	1:A:487:VAL:HG22	1.56	0.44
1:A:492:LYS:HE3	1:A:492:LYS:HB3	1.56	0.44
1:A:543:THR:C	1:A:546:CYS:H	2.25	0.44
1:A:956:ARG:O	1:A:958:HIS:N	2.50	0.44
1:A:543:THR:O	1:A:546:CYS:N	2.48	0.44
1:A:703:ALA:O	1:A:704:VAL:HG23	2.18	0.44
1:A:105:LYS:CE	1:A:175:SER:HB3	2.47	0.44
1:A:246:GLN:OE1	1:A:575:ARG:HB3	2.17	0.44
1:A:373:TYR:CD2	1:A:422:MET:HA	2.53	0.44
1:A:440:LEU:O	1:A:444:THR:OG1	2.26	0.44
1:A:736:ALA:O	1:A:740:ARG:HB2	2.18	0.44
1:A:230:LYS:HD2	1:A:539:TRP:CE3	2.52	0.44
1:A:692:MET:HE2	1:A:692:MET:HB3	1.70	0.44
1:A:726:THR:HG22	1:A:729:GLN:CB	2.48	0.44
1:A:947:TRP:HB2	1:A:982:LEU:HD22	1.99	0.44
1:A:482:PHE:CD1	1:A:490:VAL:HG11	2.53	0.44
1:A:909:LEU:HD13	1:A:1059:ASN:CG	2.43	0.44
1:A:30:PHE:CE1	1:A:881:MET:HG3	2.53	0.44
1:A:214:TYR:HA	1:A:603:PHE:CE2	2.52	0.44
1:A:601:MET:HE2	1:A:601:MET:HB2	1.66	0.44
1:A:1000:LEU:HD13	1:A:1000:LEU:O	2.17	0.44
1:A:1006:ARG:O	1:A:1009:ASP:N	2.50	0.44
1:A:1006:ARG:C	1:A:1008:LEU:N	2.75	0.44
1:A:217:PHE:O	1:A:218:VAL:C	2.60	0.44
1:A:601:MET:HE3	1:A:685:TYR:CE1	2.52	0.44
1:A:682:LEU:O	1:A:685:TYR:HB3	2.18	0.44
1:A:936:CYS:HB2	1:A:1033:LEU:HD13	2.00	0.44
1:A:185:PRO:CG	1:A:206:ILE:HD11	2.47	0.44
1:A:197:LEU:O	1:A:199:VAL:N	2.51	0.44
1:A:206:ILE:HD13	1:A:206:ILE:HA	1.87	0.44
1:A:301:GLY:O	1:A:393:PRO:HD2	2.17	0.44
1:A:410:GLN:HA	1:A:413:ARG:HH21	1.83	0.44
1:A:493:THR:O	1:A:496:LYS:HB3	2.17	0.44
1:A:546:CYS:HG	1:A:700:TRP:HE3	1.66	0.44
1:A:60:HIS:CD2	1:A:63:HIS:CE1	3.06	0.43
1:A:316:PHE:HZ	1:A:334:MET:HE3	1.83	0.43
1:A:539:TRP:NE1	1:A:691:ALA:HB2	2.33	0.43
1:A:316:PHE:CE1	1:A:324:PHE:HB2	2.53	0.43
1:A:892:ASN:O	1:A:895:LEU:HB2	2.17	0.43
1:A:224:THR:HG22	1:A:225:LEU:HD23	2.00	0.43
1:A:202:ARG:HE	1:A:202:ARG:HB2	1.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:464:LYS:HE3	1:A:468:TYR:CZ	2.53	0.43
1:A:648:THR:HG21	1:A:655:LEU:CD1	2.48	0.43
1:A:958:HIS:O	1:A:966:LEU:HG	2.18	0.43
1:A:977:GLY:HA2	1:A:986:MET:SD	2.59	0.43
1:A:1060:VAL:O	1:A:1061:PHE:CD1	2.71	0.43
1:A:266:LEU:HD22	1:A:507:TYR:HB2	2.00	0.43
1:A:498:MET:HE3	1:A:498:MET:HB2	1.84	0.43
1:A:585:TRP:HB3	5:A:1352:HOH:O	2.19	0.43
1:A:626:ARG:HE	1:A:628:GLN:HG3	1.83	0.43
1:A:24:TRP:CZ3	1:A:865:HIS:CD2	3.07	0.43
1:A:76:TYR:CE2	1:A:77:GLN:NE2	2.85	0.43
1:A:106:ARG:HG2	1:A:110:LEU:HD12	1.99	0.43
1:A:434:ILE:HD12	1:A:441:SER:OG	2.19	0.43
1:A:225:LEU:HD21	1:A:692:MET:SD	2.58	0.43
1:A:298:THR:OG1	1:A:392:VAL:HG21	2.19	0.43
1:A:333:ASN:O	1:A:335:SER:N	2.52	0.43
1:A:545:GLN:O	1:A:549:ASN:HB2	2.19	0.43
1:A:669:ASP:O	1:A:703:ALA:HB3	2.18	0.43
1:A:734:PHE:O	1:A:735:SER:C	2.59	0.43
1:A:105:LYS:HE2	1:A:175:SER:HB3	2.00	0.43
1:A:263:GLU:HG3	1:A:264:TYR:H	1.82	0.43
1:A:601:MET:CE	1:A:685:TYR:CE1	3.01	0.43
1:A:67:LEU:O	1:A:68:SER:C	2.60	0.43
1:A:400:ILE:O	1:A:404:ARG:HG2	2.18	0.43
1:A:689:HIS:HD1	1:A:701:PRO:HD3	1.83	0.43
1:A:342:ASP:O	1:A:345:VAL:HG23	2.19	0.42
1:A:391:SER:HA	1:A:399:ASP:OD1	2.18	0.42
1:A:667:PRO:HB3	1:A:699:LYS:O	2.19	0.42
1:A:731:ILE:HG12	1:A:739:MET:SD	2.58	0.42
1:A:795:ARG:HB3	1:A:895:LEU:HD11	2.01	0.42
1:A:1017:LYS:O	1:A:1018:ALA:C	2.62	0.42
1:A:577:TYR:HD1	1:A:578:GLY:N	2.17	0.42
1:A:614:ASN:O	1:A:615:LEU:HD23	2.18	0.42
1:A:791:THR:OG1	1:A:794:ASP:OD1	2.33	0.42
1:A:947:TRP:HH2	1:A:992:PHE:HD2	1.67	0.42
1:A:956:ARG:O	1:A:957:LYS:C	2.62	0.42
1:A:480:ASP:C	1:A:482:PHE:H	2.27	0.42
1:A:413:ARG:HH11	1:A:423:VAL:HG21	1.83	0.42
1:A:600:TYR:O	1:A:601:MET:C	2.61	0.42
1:A:620:GLU:HG3	1:A:621:SER:N	2.34	0.42
1:A:933:PRO:HG3	1:A:1061:PHE:CG	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:ASN:OD1	1:A:523:VAL:HB	2.20	0.42
1:A:686:LEU:CD2	1:A:701:PRO:HG2	2.49	0.42
1:A:834:LYS:HD3	1:A:838:LEU:HD22	2.01	0.42
1:A:105:LYS:HB2	1:A:105:LYS:HE3	1.72	0.42
1:A:154:SER:HB3	1:A:159:GLY:O	2.20	0.42
1:A:305:CYS:HB3	1:A:374:VAL:HG13	2.01	0.42
1:A:87:PRO:HA	1:A:199:VAL:HG13	2.02	0.42
1:A:304:ASN:CA	1:A:372:ILE:HD12	2.50	0.42
1:A:571:HIS:CE1	1:A:592:GLU:OE2	2.72	0.42
1:A:803:ALA:HB2	1:A:891:VAL:CG1	2.49	0.42
1:A:272:LEU:HB2	1:A:361:SER:C	2.45	0.42
1:A:381:LYS:HD2	1:A:381:LYS:HA	1.82	0.42
1:A:773:TRP:CE3	1:A:832:LYS:HE2	2.54	0.42
1:A:353:MET:C	1:A:355:GLU:H	2.27	0.42
1:A:48:VAL:HG22	1:A:86:PHE:HB2	2.01	0.41
1:A:199:VAL:CG1	1:A:201:TRP:CD1	3.03	0.41
1:A:212:PRO:HB2	1:A:627:PRO:HB3	2.00	0.41
1:A:213:TYR:O	1:A:214:TYR:C	2.63	0.41
1:A:412:LEU:C	1:A:414:ALA:N	2.78	0.41
1:A:434:ILE:HA	1:A:435:PRO:HD3	1.94	0.41
1:A:1017:LYS:O	1:A:1019:VAL:N	2.53	0.41
1:A:1021:MET:HE3	1:A:1021:MET:HB3	1.90	0.41
1:A:262:GLN:OE1	1:A:264:TYR:HE1	2.03	0.41
1:A:286:ASN:OD1	1:A:286:ASN:N	2.54	0.41
1:A:313:TYR:HB2	1:A:352:LEU:HD21	2.02	0.41
1:A:499:ILE:HA	1:A:504:ALA:O	2.20	0.41
1:A:696:GLN:O	1:A:697:SER:C	2.62	0.41
1:A:1002:LYS:HA	1:A:1002:LYS:HD2	1.89	0.41
1:A:598:THR:OG1	1:A:688:ASN:OD1	2.32	0.41
1:A:637:TYR:HA	1:A:643:ALA:CB	2.50	0.41
1:A:268:LYS:HE2	1:A:268:LYS:HB3	1.78	0.41
1:A:420:ASP:HB3	1:A:424:LEU:HD12	2.02	0.41
1:A:670:LEU:HD12	1:A:670:LEU:HA	1.86	0.41
1:A:424:LEU:HB2	1:A:425:PRO:HD3	2.01	0.41
1:A:532:LEU:O	1:A:534:TYR:N	2.52	0.41
1:A:568:LEU:HD12	1:A:568:LEU:HA	1.78	0.41
1:A:774:VAL:O	1:A:778:VAL:HG23	2.20	0.41
1:A:964:GLY:O	1:A:965:LYS:HB2	2.20	0.41
1:A:976:LEU:HD22	1:A:976:LEU:HA	1.78	0.41
1:A:266:LEU:HD11	1:A:290:VAL:HG12	2.02	0.41
1:A:284:GLY:C	1:A:285:LYS:HG3	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:314:ILE:O	1:A:325:ILE:HA	2.20	0.41
1:A:413:ARG:HB3	1:A:418:ILE:O	2.19	0.41
1:A:614:ASN:C	1:A:615:LEU:HD23	2.45	0.41
1:A:52:TYR:CD1	1:A:91:HIS:CG	3.09	0.41
1:A:52:TYR:HB3	1:A:89:GLY:O	2.20	0.41
1:A:171:ILE:HG13	1:A:171:ILE:H	1.52	0.41
1:A:633:GLU:HB3	1:A:646:PRO:HB3	2.03	0.41
1:A:770:TRP:CZ3	1:A:851:PHE:CE2	3.07	0.41
1:A:1007:ILE:N	1:A:1007:ILE:CD1	2.83	0.41
1:A:75:GLY:CA	1:A:198:LYS:HE2	2.51	0.41
1:A:335:SER:HB3	1:A:340:THR:HG21	2.02	0.41
1:A:401:ALA:O	1:A:404:ARG:HB2	2.21	0.41
1:A:581:THR:HB	1:A:591:ILE:HD12	2.02	0.41
1:A:773:TRP:O	1:A:776:GLU:HB3	2.21	0.41
1:A:799:SER:CB	1:A:892:ASN:H	2.27	0.41
1:A:48:VAL:HA	1:A:670:LEU:HB3	2.03	0.41
1:A:61:LEU:O	1:A:63:HIS:N	2.54	0.41
1:A:311:MET:HB3	1:A:313:TYR:CZ	2.56	0.41
1:A:311:MET:O	1:A:353:MET:HA	2.21	0.41
1:A:630:MET:HG3	1:A:634:VAL:CG1	2.51	0.41
1:A:714:SER:HG	1:A:754:ASN:HD21	1.68	0.41
1:A:725:LEU:HD13	1:A:730:ALA:HB2	2.03	0.41
1:A:827:GLU:O	1:A:830:ALA:HB3	2.21	0.41
1:A:92:CYS:HA	1:A:97:ILE:HD11	2.03	0.41
1:A:755:PHE:CD1	1:A:755:PHE:C	2.99	0.41
1:A:810:GLN:O	1:A:814:LYS:HG3	2.21	0.41
1:A:199:VAL:HG11	1:A:201:TRP:HD1	1.86	0.40
1:A:242:PRO:HB3	1:A:336:TYR:CZ	2.56	0.40
1:A:803:ALA:HB2	1:A:891:VAL:HG13	2.03	0.40
1:A:369:TYR:O	1:A:370:LYS:C	2.64	0.40
1:A:413:ARG:H	1:A:413:ARG:HG3	1.67	0.40
1:A:1027:LEU:HB3	1:A:1031:LEU:HD12	2.02	0.40
1:A:269:LEU:O	1:A:288:PHE:HD1	2.02	0.40
1:A:354:GLY:HA2	1:A:357:ILE:CD1	2.52	0.40
1:A:434:ILE:CG1	1:A:471:GLY:HA2	2.40	0.40
1:A:493:THR:HA	1:A:496:LYS:CB	2.51	0.40
1:A:737:ASP:OD2	1:A:863:CYS:HB3	2.22	0.40
1:A:996:ILE:HG23	1:A:1007:ILE:HG21	2.03	0.40
1:A:1007:ILE:O	1:A:1007:ILE:HG22	2.22	0.40
1:A:62:GLY:H	1:A:717:MET:HE1	1.87	0.40
1:A:117:PHE:CE2	1:A:156:TYR:CD1	3.09	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:ARG:HG2	1:A:584:PRO:HG3	2.02	0.40
1:A:479:VAL:HG22	1:A:480:ASP:N	2.37	0.40
1:A:645:PHE:HA	1:A:646:PRO:HD3	1.83	0.40
1:A:705:ARG:HH22	1:A:748:ASP:HB2	1.86	0.40
1:A:240:TYR:HB3	1:A:523:VAL:O	2.22	0.40
1:A:423:VAL:CG2	1:A:424:LEU:N	2.83	0.40
1:A:1060:VAL:HG12	1:A:1061:PHE:H	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	998/1188 (84%)	810 (81%)	177 (18%)	11 (1%)	11 39

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	587	GLU
1	A	978	SER
1	A	62	GLY
1	A	164	LEU
1	A	354	GLY
1	A	198	LYS
1	A	334	MET
1	A	390	THR
1	A	720	SER
1	A	951	THR
1	A	376	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	885/1047 (84%)	879 (99%)	6 (1%)	76	80

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	163	SER
1	A	448	GLU
1	A	449	LEU
1	A	765	LEU
1	A	952	LEU
1	A	976	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	HIS
1	A	304	ASN
1	A	343	ASN
1	A	455	ASN
1	A	629	GLN
1	A	649	GLN
1	A	729	GLN
1	A	754	ASN
1	A	793	ASN
1	A	869	HIS
1	A	892	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 ⓘ

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
4	PO4	A	1204	-	4,4,4	0.76	0	6,6,6	0.70	0
2	ATP	A	1201	-	32,33,33	2.66	6 (18%)	48,52,52	2.28	17 (35%)
3	LEU	A	1202	-	6,8,8	1.17	1 (16%)	5,10,10	1.52	2 (40%)
4	PO4	A	1203	-	4,4,4	1.03	0	6,6,6	0.88	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	A	1201	-	-	6/22/38/38	0/3/3/3
3	LEU	A	1202	-	-	2/8/8/8	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1201	ATP	PB-O3B	10.77	1.71	1.59
2	A	1201	ATP	PA-O3A	5.25	1.65	1.59
2	A	1201	ATP	C5-C4	4.94	1.47	1.39
2	A	1201	ATP	PB-O3A	4.48	1.64	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1201	ATP	C8-N7	2.74	1.37	1.31
2	A	1201	ATP	C5-N7	-2.74	1.34	1.39
3	A	1202	LEU	OXT-C	-2.39	1.23	1.30

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1201	ATP	C5-C4-N3	-6.56	117.68	126.72
2	A	1201	ATP	N3-C4-N9	6.23	137.77	127.17
2	A	1201	ATP	N6-C6-N1	4.09	127.49	118.38
2	A	1201	ATP	C2-N3-C4	3.70	120.86	111.83
2	A	1201	ATP	N3-C2-N1	-3.38	123.47	128.58
2	A	1201	ATP	C5-C6-N6	-3.19	115.39	123.29
2	A	1201	ATP	C2'-C3'-C4'	3.15	108.69	102.61
2	A	1201	ATP	O3'-C3'-C2'	-3.09	101.91	111.82
2	A	1201	ATP	C4-N9-C8	2.88	108.76	105.74
2	A	1201	ATP	O2A-PA-O3A	2.83	114.91	107.27
2	A	1201	ATP	O2'-C2'-C3'	-2.75	103.00	111.82
2	A	1201	ATP	C1'-N9-C8	-2.68	121.15	127.09
2	A	1201	ATP	O2B-PB-O3A	-2.62	100.19	107.27
2	A	1201	ATP	C2'-C1'-N9	2.61	119.80	113.30
3	A	1202	LEU	OXT-C-O	-2.47	118.48	124.08
2	A	1201	ATP	O3G-PG-O3B	2.29	112.31	104.64
2	A	1201	ATP	O4'-C1'-N9	-2.20	103.86	108.09
2	A	1201	ATP	N9-C8-N7	-2.12	110.92	113.94
3	A	1202	LEU	CB-CA-C	-2.09	103.99	110.34

There are no chirality outliers.

All (8) torsion outliers are listed below:

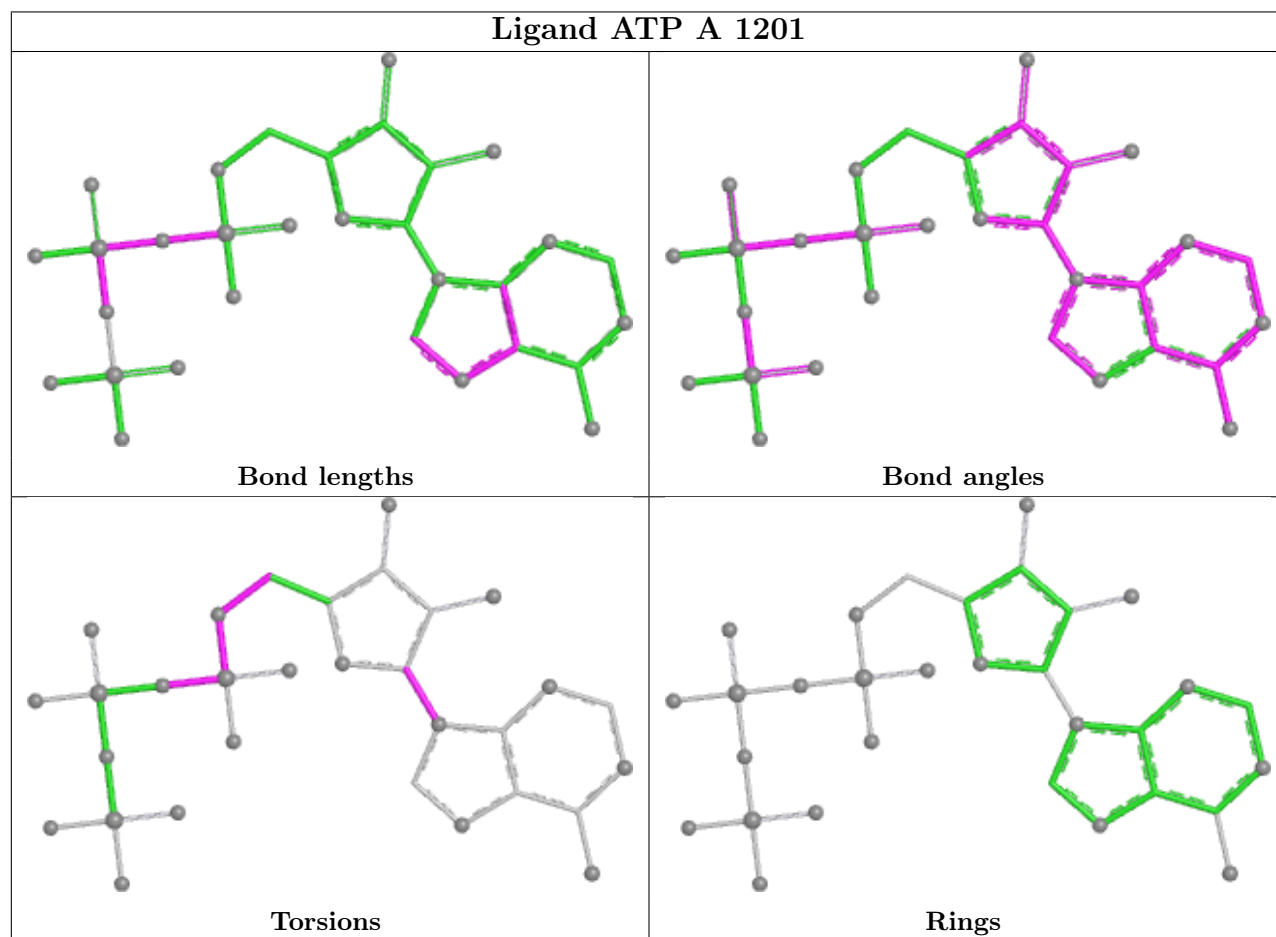
Mol	Chain	Res	Type	Atoms
2	A	1201	ATP	C5'-O5'-PA-O1A
2	A	1201	ATP	C5'-O5'-PA-O2A
3	A	1202	LEU	CA-CB-CG-CD1
3	A	1202	LEU	CA-CB-CG-CD2
2	A	1201	ATP	C5'-O5'-PA-O3A
2	A	1201	ATP	C4'-C5'-O5'-PA
2	A	1201	ATP	PB-O3A-PA-O2A
2	A	1201	ATP	O4'-C1'-N9-C8

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1201	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	1004/1188 (84%)	0.10	19 (1%)	66 45	14, 40, 107, 152	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1017	LYS	4.3
1	A	916	TYR	3.5
1	A	37	LEU	3.1
1	A	939	TYR	3.0
1	A	958	HIS	2.8
1	A	261	PRO	2.7
1	A	954	VAL	2.7
1	A	567	TRP	2.6
1	A	941	ALA	2.5
1	A	980	PRO	2.5
1	A	976	LEU	2.4
1	A	943	ASN	2.4
1	A	1010	LEU	2.4
1	A	1057	PRO	2.3
1	A	940	VAL	2.3
1	A	993	VAL	2.3
1	A	972	ILE	2.3
1	A	249	MET	2.2
1	A	971	VAL	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

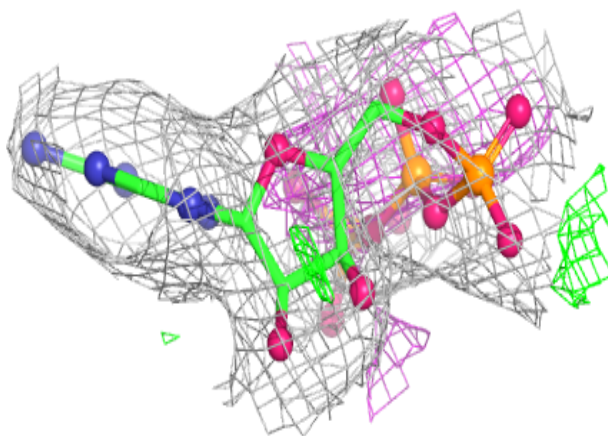
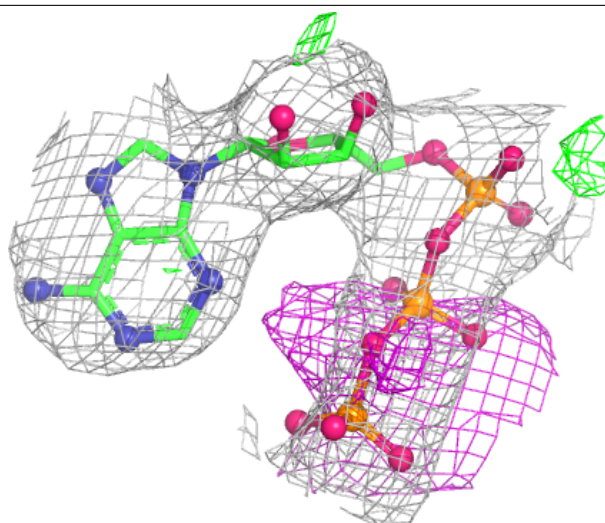
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	PO4	A	1203	5/5	0.77	0.24	59,69,71,95	0
3	LEU	A	1202	9/9	0.90	0.15	17,34,44,47	0
2	ATP	A	1201	31/31	0.90	0.12	14,38,87,99	0
4	PO4	A	1204	5/5	0.93	0.13	23,44,61,65	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 A 1201:

$2mF_o-DF_c$ (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.