



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:49 PM UTC

PDB ID : 1E94 / pdb_00001e94
Title : HslV-HslU from E.coli
Authors : Song, H.K.; Hartmann, C.; Ravishankar, R.; Bochtler, M.
Deposited on : 2000-10-07
Resolution : 2.80 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

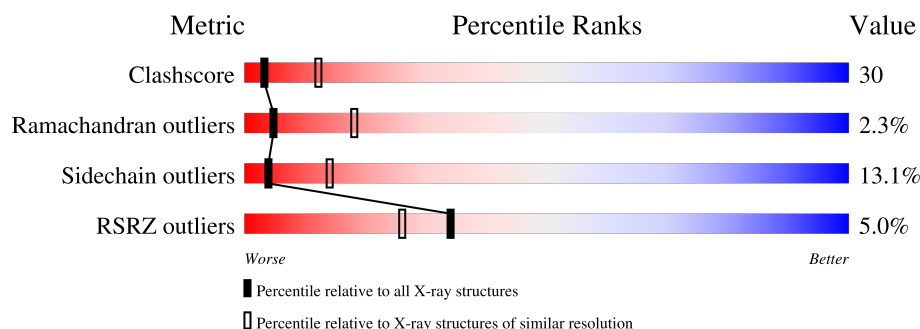
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	175	3% 62% 32% 5%
1	B	175	2% 46% 45% 8%
1	C	175	2% 55% 37% 7%
1	D	175	6% 45% 49% 6%
2	E	449	5% 43% 37% 9% 9%
2	F	449	6% 47% 33% 10% 9%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ANP	F	501	-	X	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12028 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.

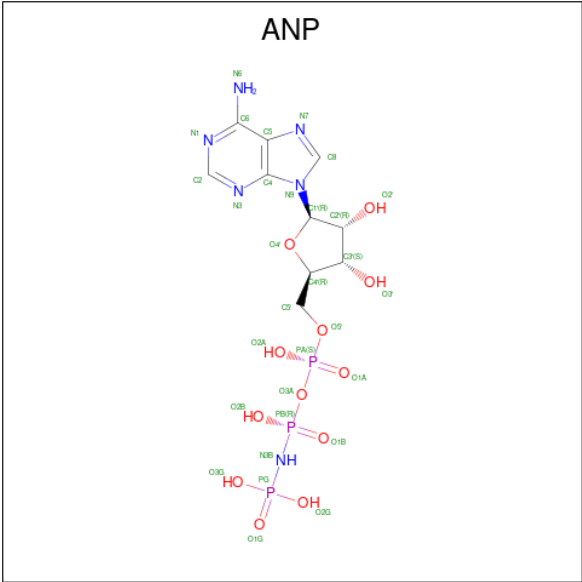
- Molecule 1 is a protein called HEAT SHOCK PROTEIN HSLV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	174	Total	C	N	O	S	25	0	0
			1328	834	237	253	4			
1	B	174	Total	C	N	O	S	24	0	0
			1328	834	237	253	4			
1	C	174	Total	C	N	O	S	18	0	0
			1328	834	237	253	4			
1	D	174	Total	C	N	O	S	24	0	0
			1328	834	237	253	4			

- Molecule 2 is a protein called HEAT SHOCK PROTEIN HSLU.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	408	Total	C	N	O	S	220	1	0
			3184	1989	570	615	10			
2	F	409	Total	C	N	O	S	443	0	1
			3184	1989	571	614	10			

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



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

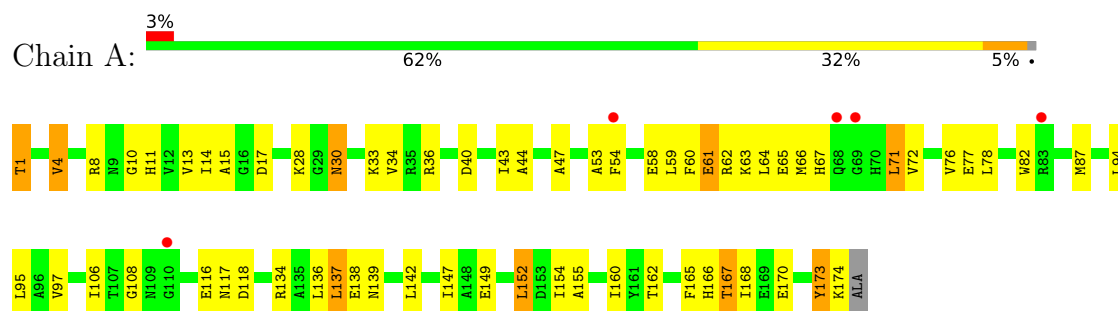
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	31	Total	O	0	0
			31	31		
4	B	29	Total	O	0	0
			29	29		
4	C	33	Total	O	0	0
			33	33		
4	D	31	Total	O	0	0
			31	31		
4	E	94	Total	O	0	0
			94	94		
4	F	68	Total	O	0	0
			68	68		

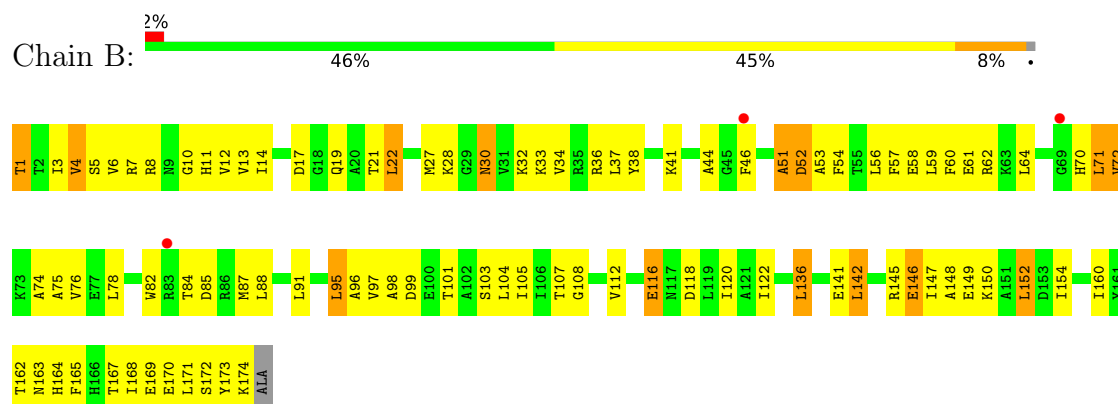
3 Residue-property plots [i](#)

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.

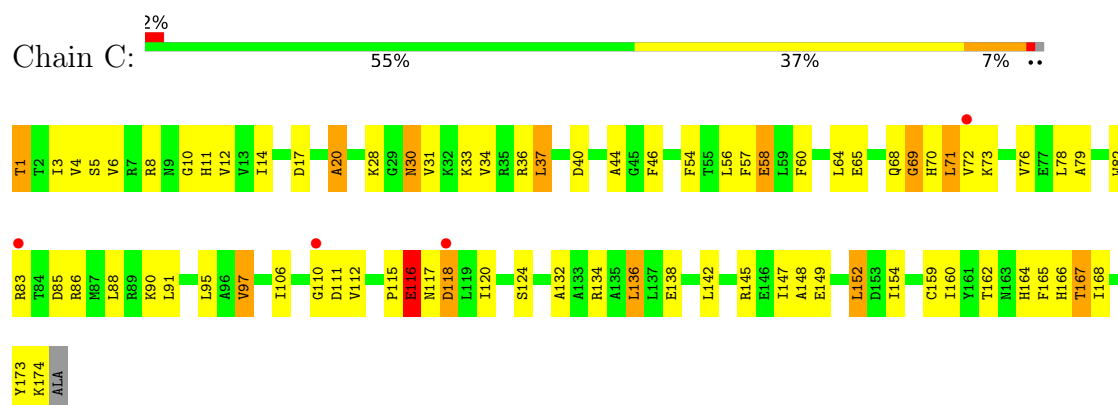
• Molecule 1: HEAT SHOCK PROTEIN HSLV



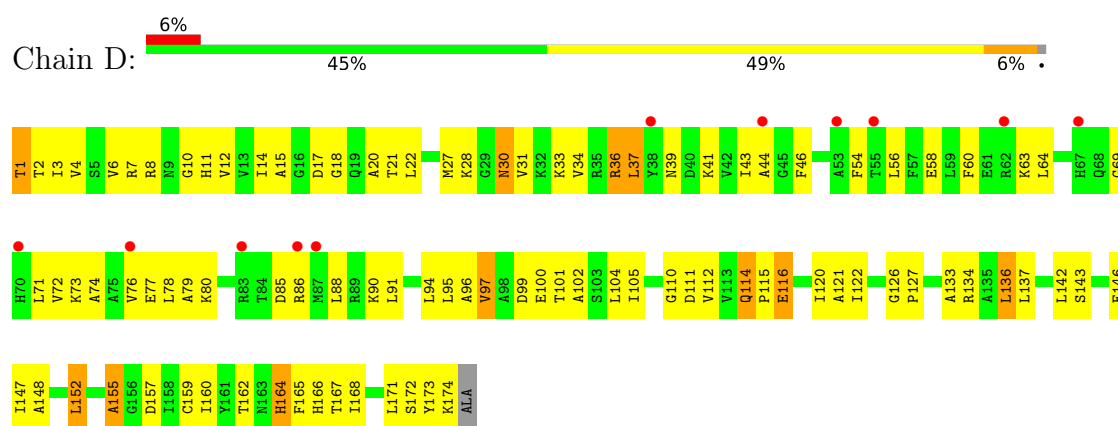
• Molecule 1: HEAT SHOCK PROTEIN HSLV



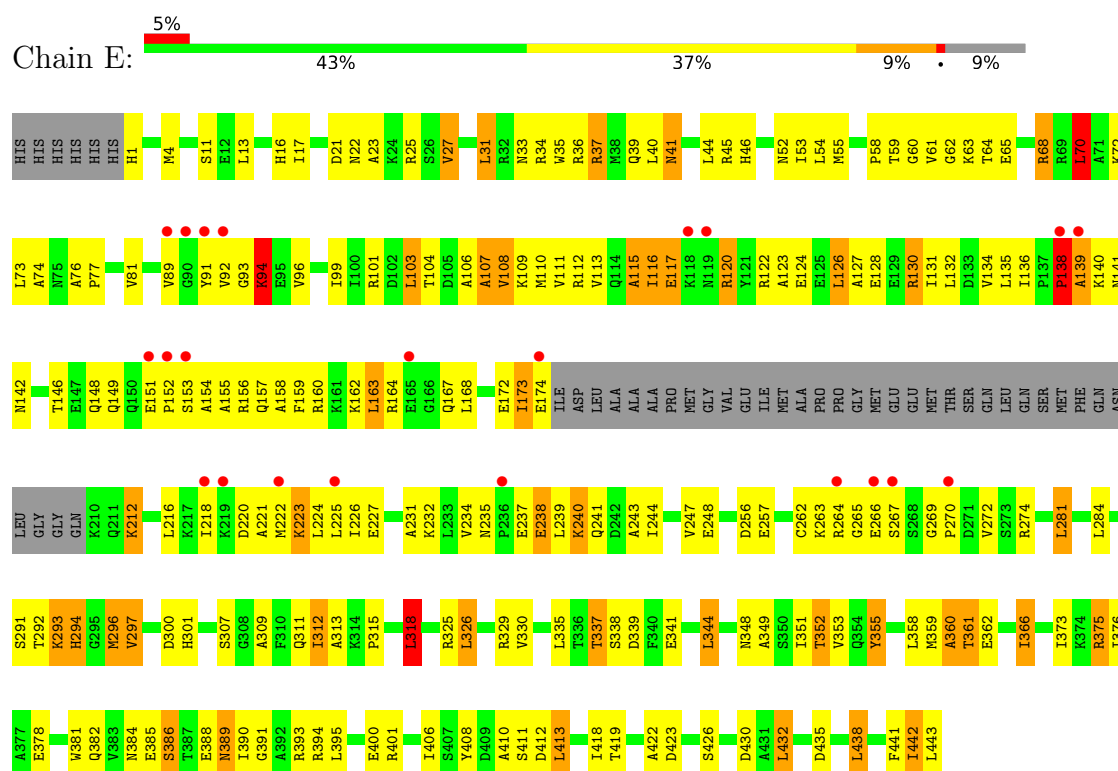
• Molecule 1: HEAT SHOCK PROTEIN HSLV



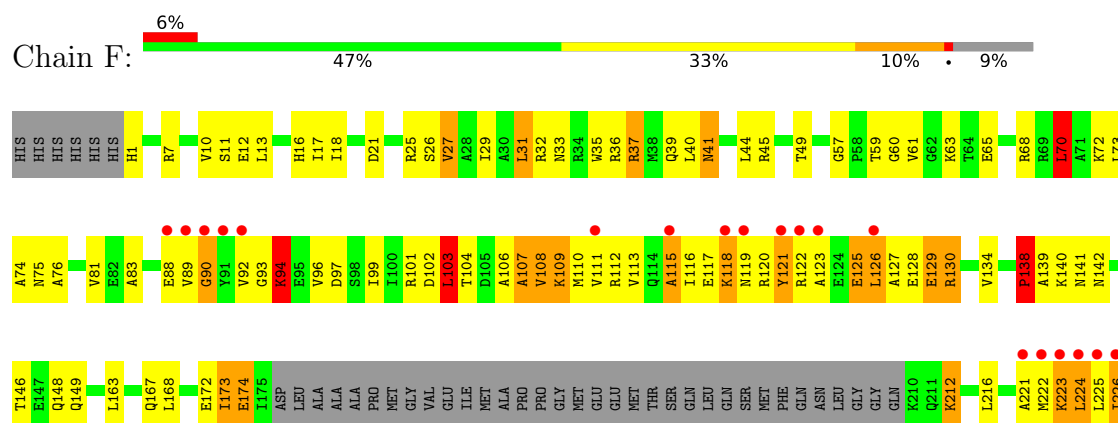
• Molecule 1: HEAT SHOCK PROTEIN HSLV

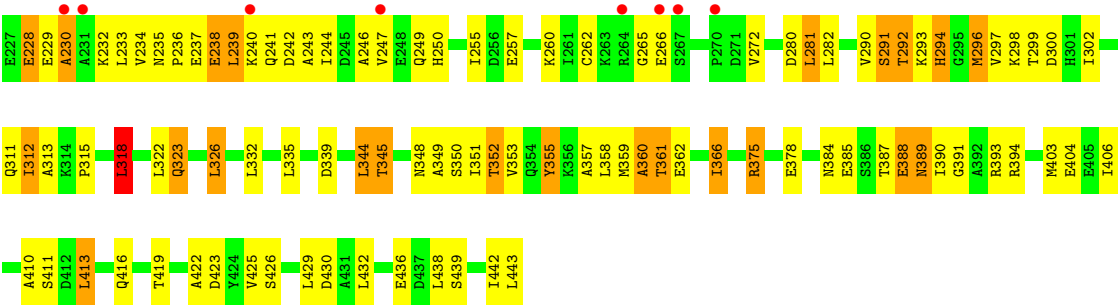


• Molecule 2: HEAT SHOCK PROTEIN HSLU



• Molecule 2: HEAT SHOCK PROTEIN HSLU





4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	172.02Å 172.02Å 276.57Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 2.80 15.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	92.3 (15.00-2.80) 91.6 (15.00-2.80)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.48 (at 2.80Å)	Xtrriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.254 , 0.304 0.259 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	68.0	Xtrriage
Anisotropy	0.372	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 83.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12028	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 61.06 % 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 1.3856e-05. 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.77	0/1345	1.12	6/1817 (0.3%)
1	B	0.73	1/1345 (0.1%)	1.14	8/1817 (0.4%)
1	C	0.63	0/1345	1.04	7/1817 (0.4%)
1	D	0.66	0/1345	1.03	3/1817 (0.2%)
2	E	0.74	1/3228 (0.0%)	1.21	26/4352 (0.6%)
2	F	0.85	4/3221 (0.1%)	1.43	32/4345 (0.7%)
All	All	0.75	6/11829 (0.1%)	1.22	82/15965 (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
2	E	0	1
2	F	0	2
All	All	0	3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	224	LEU	C-N	19.55	1.62	1.33
2	F	129	GLU	C-N	-13.78	1.14	1.33
2	E	138	PRO	C-N	-10.86	1.18	1.33
2	F	138	PRO	C-N	-10.86	1.18	1.33
1	B	27	MET	SD-CE	5.82	1.94	1.79
2	F	174	GLU	C-N	-5.71	1.25	1.33

All (82) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	129	GLU	O-C-N	-31.07	81.27	122.59
2	F	224	LEU	O-C-N	29.87	156.26	122.20
2	E	138	PRO	O-C-N	-21.43	93.71	122.64
2	F	138	PRO	O-C-N	-21.40	93.75	122.64
2	F	224	LEU	CA-C-N	-20.15	82.28	121.58
2	F	224	LEU	C-N-CA	-20.15	82.28	121.58
2	E	120	ARG	N-CA-C	-9.35	103.96	114.62
2	E	293	LYS	N-CA-C	-8.79	103.65	112.97
2	E	107	ALA	N-CA-C	-8.12	102.82	112.89
2	E	126	LEU	N-CA-C	-7.92	103.83	113.97
2	F	134	VAL	N-CA-C	-7.84	105.37	112.90
2	E	134	VAL	N-CA-C	-7.82	105.40	112.90
2	F	293	LYS	N-CA-C	-7.76	104.75	112.97
2	E	318	LEU	N-CA-C	-7.63	100.45	110.53
2	F	118	LYS	N-CA-C	-7.59	103.13	113.30
2	F	107	ALA	N-CA-C	-7.58	104.00	113.18
1	B	112	VAL	N-CA-C	-7.54	96.73	107.75
2	E	239	LEU	N-CA-C	-7.04	102.69	112.12
1	B	96	ALA	N-CA-C	-6.99	97.02	108.41
2	E	272	VAL	N-CA-C	-6.91	102.45	111.09
1	B	41	LYS	N-CA-C	-6.90	105.02	113.50
2	F	126	LEU	N-CA-C	-6.75	105.81	112.97
2	F	238	GLU	N-CA-C	-6.72	102.92	112.13
1	A	61	GLU	N-CA-C	-6.58	104.10	111.28
2	F	115	ALA	N-CA-C	-6.58	105.83	114.31
1	A	116	GLU	N-CA-C	6.48	124.60	110.80
2	E	294	HIS	N-CA-C	-6.42	104.65	114.16
1	C	165	PHE	N-CA-C	-6.38	99.61	109.76
1	D	164	HIS	N-CA-C	6.34	119.95	112.72
1	C	20	ALA	N-CA-C	-6.34	98.57	108.90
2	F	318	LEU	N-CA-C	-6.28	101.65	110.50
1	C	167	THR	N-CA-C	-6.27	97.19	108.48
2	E	376	ILE	N-CA-C	-6.27	104.64	110.53
2	F	360	ALA	N-CA-C	-6.19	104.45	111.07
2	F	94	LYS	N-CA-C	-6.17	101.35	110.48
2	E	70	LEU	N-CA-C	-5.97	104.12	111.40
2	F	294	HIS	N-CA-C	-5.96	104.85	114.09
1	A	173	TYR	N-CA-C	-5.95	105.77	114.39
2	E	62	GLY	N-CA-C	5.95	123.33	114.95
2	E	94	LYS	N-CA-C	-5.90	102.18	110.50
2	F	326	LEU	CA-C-N	5.89	125.37	119.24
2	F	326	LEU	C-N-CA	5.89	125.37	119.24
2	F	70	LEU	N-CA-C	-5.89	104.77	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	4	VAL	CB-CA-C	-5.83	101.93	110.62
2	E	360	ALA	N-CA-C	-5.83	104.12	111.11
2	F	429	LEU	N-CA-C	5.81	120.50	113.41
1	C	164	HIS	N-CA-C	5.79	120.04	112.92
1	D	155	ALA	N-CA-C	-5.79	104.66	110.97
2	F	323	GLN	N-CA-C	-5.76	105.01	111.28
2	F	163	LEU	N-CA-C	-5.73	106.19	112.72
2	E	163	LEU	N-CA-C	-5.73	106.19	112.72
1	B	52	ASP	N-CA-C	-5.61	104.38	111.11
2	E	338	SER	N-CA-C	-5.59	105.26	111.36
2	F	272	VAL	N-CA-C	-5.59	104.10	111.09
1	A	167	THR	N-CA-C	-5.57	98.46	108.48
2	E	326	LEU	CA-C-N	5.53	125.62	119.32
2	E	326	LEU	C-N-CA	5.53	125.62	119.32
1	B	104	LEU	N-CA-C	5.53	116.94	108.42
1	B	4	VAL	CB-CA-C	-5.47	103.00	110.77
2	E	46	HIS	N-CA-C	-5.46	104.87	112.45
1	D	20	ALA	N-CA-C	-5.45	100.01	108.90
2	E	173	ILE	N-CA-C	5.41	115.42	108.82
2	F	173	ILE	N-CA-C	5.41	115.42	108.82
2	F	90	GLY	N-CA-C	-5.39	107.98	114.66
2	F	83	ALA	N-CA-C	5.36	117.54	111.11
2	F	57	GLY	CA-C-N	5.35	125.34	120.21
2	F	57	GLY	C-N-CA	5.35	125.34	120.21
2	E	297	VAL	N-CA-C	5.30	117.50	109.12
2	F	436	GLU	N-CA-C	5.29	117.12	111.36
2	E	309	ALA	N-CA-C	-5.26	105.14	112.45
1	B	164	HIS	N-CA-C	5.25	118.71	112.72
2	F	241	GLN	N-CA-C	-5.25	106.93	113.38
1	C	65	GLU	N-CA-C	-5.23	106.41	112.89
1	A	142	LEU	N-CA-C	5.20	116.99	110.24
2	E	232	LYS	N-CA-C	-5.17	106.48	112.89
2	E	115	ALA	N-CA-C	-5.15	105.29	112.45
1	C	118	ASP	N-CA-C	-5.14	106.96	113.23
2	F	103	LEU	N-CA-C	-5.09	105.01	111.11
2	F	345	THR	N-CA-C	5.09	121.64	110.80
1	C	145	ARG	N-CA-C	5.08	116.82	111.28
1	B	51	ALA	N-CA-C	-5.07	105.46	111.69
2	E	442	ILE	N-CA-C	5.03	119.96	113.07

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	E	138	PRO	Mainchain
2	F	129	GLU	Mainchain
2	F	138	PRO	Mainchain

5.2 Too-close contacts [i](#)

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	1328	0	1348	53	0
1	B	1328	0	1347	71	0
1	C	1328	0	1348	60	0
1	D	1328	0	1348	80	0
2	E	3184	0	3206	232	0
2	F	3184	0	3208	183	0
3	E	31	0	12	7	0
3	F	31	0	12	9	0
4	A	31	0	0	3	0
4	B	29	0	0	1	0
4	C	33	0	0	5	0
4	D	31	0	0	5	0
4	E	94	0	0	11	0
4	F	68	0	0	10	0
All	All	12028	0	11829	672	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 30.

All (672) 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)
2:E:153:SER:CA	2:E:156:ARG:HB3	1.47	1.44
2:E:153:SER:HB3	2:E:157:GLN:CG	1.52	1.37
2:E:153:SER:C	2:E:154:ALA:C	1.93	1.35
2:E:131:ILE:HD11	2:E:218:ILE:CD1	1.57	1.34
2:E:151:GLU:O	2:E:154:ALA:N	1.57	1.32
2:F:224:LEU:O	2:F:228:GLU:HB2	1.14	1.29
2:E:152:PRO:O	2:E:156:ARG:N	1.65	1.28

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:153:SER:CA	2:E:157:GLN:H	1.48	1.26
2:E:153:SER:HA	2:E:156:ARG:CB	1.68	1.22
2:E:153:SER:O	2:E:154:ALA:O	1.58	1.18
2:E:220:ASP:C	2:E:224:LEU:HD23	1.69	1.16
2:F:221:ALA:O	2:F:225:LEU:HD23	1.46	1.15
2:E:131:ILE:CD1	2:E:218:ILE:HD11	1.75	1.15
2:F:128:GLU:CB	2:F:222:MET:HE1	1.75	1.15
2:F:223:LYS:HA	2:F:226:ILE:HG13	1.26	1.11
2:E:153:SER:HA	2:E:157:GLN:N	1.65	1.10
2:E:153:SER:HA	2:E:157:GLN:H	1.06	1.10
2:F:128:GLU:HB2	2:F:222:MET:CE	1.83	1.07
2:F:223:LYS:HA	2:F:226:ILE:CG1	1.84	1.06
2:E:153:SER:N	2:E:156:ARG:HB3	1.69	1.05
2:F:223:LYS:CA	2:F:226:ILE:HG13	1.87	1.04
2:E:153:SER:HB3	2:E:157:GLN:HG3	1.07	1.03
2:E:220:ASP:O	2:E:224:LEU:HD23	1.58	1.02
2:E:153:SER:CB	2:E:157:GLN:HG3	1.89	1.01
2:E:132:LEU:HD11	2:E:160:ARG:HG3	1.45	0.99
2:F:224:LEU:O	2:F:228:GLU:CB	2.10	0.98
2:F:223:LYS:HA	2:F:226:ILE:CD1	1.94	0.96
2:F:103:LEU:HD13	2:F:247:VAL:HG13	1.47	0.96
2:E:153:SER:CA	2:E:156:ARG:CB	2.36	0.95
2:E:220:ASP:O	2:E:224:LEU:CD2	2.13	0.94
2:E:153:SER:C	2:E:154:ALA:CA	2.39	0.94
2:E:291:SER:HA	2:E:296:MET:HE2	1.46	0.94
2:E:160:ARG:HH12	2:E:164:ARG:HH22	1.16	0.93
2:F:119:ASN:HD21	2:F:233:LEU:HD23	1.31	0.93
2:E:131:ILE:HD11	2:E:218:ILE:HD11	0.93	0.93
2:E:153:SER:HA	2:E:156:ARG:HB3	0.97	0.92
2:E:153:SER:CB	2:E:157:GLN:CG	2.45	0.91
1:C:30:ASN:HD22	1:C:30:ASN:H	1.16	0.91
2:E:131:ILE:HD11	2:E:218:ILE:CG1	2.01	0.90
2:F:221:ALA:O	2:F:225:LEU:CD2	2.19	0.90
2:F:291:SER:HA	2:F:296:MET:HE2	1.55	0.89
2:E:344:LEU:O	2:E:352:THR:HB	1.72	0.89
2:E:152:PRO:C	2:E:156:ARG:H	1.80	0.88
2:E:218:ILE:O	2:E:221:ALA:N	2.05	0.88
1:A:134:ARG:HD2	4:A:2019:HOH:O	1.72	0.88
2:E:389:ASN:ND2	2:E:391:GLY:H	1.71	0.88
2:F:221:ALA:C	2:F:225:LEU:HD23	1.99	0.87
2:F:223:LYS:O	2:F:226:ILE:N	2.07	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:220:ASP:C	2:E:224:LEU:CD2	2.47	0.86
2:E:103:LEU:HD13	2:E:247:VAL:HG13	1.58	0.86
2:E:153:SER:CB	2:E:157:GLN:NE2	2.39	0.86
1:A:1:THR:HB	1:A:33:LYS:NZ	1.92	0.85
2:E:153:SER:N	2:E:156:ARG:CB	2.37	0.85
2:F:33:ASN:HD22	2:F:36:ARG:HD2	1.41	0.84
2:E:153:SER:CA	2:E:157:GLN:N	2.30	0.84
2:F:65:GLU:HG3	3:F:501:ANP:H2'	1.58	0.84
2:F:128:GLU:HB2	2:F:222:MET:HE1	0.88	0.83
2:F:389:ASN:ND2	2:F:391:GLY:H	1.76	0.83
2:E:153:SER:HB3	2:E:157:GLN:CB	2.08	0.83
2:E:153:SER:HA	2:E:156:ARG:CA	2.08	0.82
1:A:1:THR:HB	1:A:33:LYS:HZ3	1.44	0.81
2:F:344:LEU:O	2:F:352:THR:HB	1.81	0.81
2:E:153:SER:HB3	2:E:157:GLN:CD	2.07	0.80
1:A:10:GLY:HA2	1:A:173:TYR:CE1	2.16	0.80
2:F:235:ASN:OD1	2:F:238:GLU:HB2	1.81	0.80
2:E:96:VAL:HG11	2:E:281:LEU:HD12	1.63	0.79
2:F:221:ALA:O	2:F:225:LEU:N	2.15	0.79
2:F:223:LYS:C	2:F:225:LEU:N	2.37	0.79
2:F:41:ASN:HD21	2:F:44:LEU:H	1.31	0.78
1:D:152:LEU:HD13	1:D:166:HIS:ND1	1.98	0.78
2:F:229:GLU:HA	2:F:232:LYS:HD3	1.65	0.78
2:E:153:SER:O	2:E:157:GLN:N	2.17	0.78
1:B:72:VAL:O	1:B:76:VAL:HG23	1.84	0.77
2:E:101:ARG:HG2	4:E:2057:HOH:O	1.83	0.77
2:E:65:GLU:HG3	3:E:500:ANP:H2'	1.65	0.77
2:E:131:ILE:CD1	2:E:218:ILE:CD1	2.48	0.77
1:A:10:GLY:HA3	1:A:174:LYS:HA	1.66	0.77
1:B:44:ALA:HB2	1:B:97:VAL:HG23	1.66	0.77
2:E:220:ASP:O	2:E:223:LYS:HB2	1.85	0.77
1:A:87:MET:HE1	1:B:84:THR:HG23	1.67	0.77
2:E:312:ILE:HD12	2:E:312:ILE:H	1.51	0.76
2:E:131:ILE:CD1	2:E:218:ILE:CG1	2.62	0.76
1:B:28:LYS:NZ	1:B:30:ASN:HD21	1.84	0.76
2:F:344:LEU:HD12	2:F:351:ILE:HD11	1.68	0.76
1:A:28:LYS:NZ	1:A:30:ASN:HD21	1.84	0.76
2:E:389:ASN:C	2:E:389:ASN:HD22	1.94	0.76
1:D:36:ARG:C	1:D:37:LEU:HD23	2.11	0.75
2:F:33:ASN:ND2	2:F:36:ARG:HD2	2.02	0.75
1:D:30:ASN:HD22	1:D:30:ASN:H	1.34	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:122:ARG:NH1	2:E:126:LEU:HD21	2.02	0.75
1:B:28:LYS:HZ2	1:B:30:ASN:HD21	1.32	0.75
2:F:389:ASN:HD22	2:F:390:ILE:N	1.85	0.75
2:E:384:ASN:ND2	2:E:394:ARG:HE	1.84	0.75
2:F:27:VAL:HG13	2:F:70:LEU:HG	1.68	0.75
2:E:94:LYS:HE2	2:E:94:LYS:HA	1.67	0.74
2:E:153:SER:CB	2:E:157:GLN:CD	2.60	0.74
2:E:37:ARG:HH11	2:E:37:ARG:HB3	1.53	0.74
2:F:389:ASN:HD22	2:F:389:ASN:C	1.96	0.74
2:E:153:SER:HB2	2:E:157:GLN:NE2	2.02	0.74
1:B:88:LEU:HA	1:B:91:LEU:HD12	1.69	0.74
2:E:153:SER:C	2:E:157:GLN:H	1.95	0.73
2:F:223:LYS:N	2:F:223:LYS:HD2	2.03	0.73
2:E:223:LYS:HD2	2:E:223:LYS:N	2.03	0.73
2:E:359:MET:HE1	2:F:36:ARG:NH1	2.02	0.73
2:E:432:LEU:HD12	2:E:432:LEU:H	1.54	0.73
2:F:122:ARG:HH11	2:F:126:LEU:HD23	1.54	0.73
1:B:28:LYS:NZ	1:B:30:ASN:ND2	2.37	0.73
1:C:6:VAL:HG21	1:C:147:ILE:HG22	1.70	0.73
1:B:149:GLU:OE1	1:B:168:ILE:HD11	1.89	0.72
2:F:312:ILE:HD12	2:F:313:ALA:H	1.53	0.72
2:F:312:ILE:HD12	2:F:313:ALA:N	2.05	0.72
2:E:160:ARG:NH1	2:E:164:ARG:HH22	1.87	0.72
1:A:10:GLY:HA2	1:A:173:TYR:CZ	2.26	0.71
2:E:153:SER:HA	2:E:156:ARG:C	2.15	0.71
1:C:30:ASN:H	1:C:30:ASN:ND2	1.87	0.71
2:E:91:TYR:HB2	2:F:90:GLY:O	1.90	0.70
2:F:104:THR:HG21	2:F:292:THR:HG21	1.73	0.70
2:E:152:PRO:C	2:E:156:ARG:CB	2.64	0.70
2:F:18:ILE:H	3:F:501:ANP:HN61	1.39	0.70
2:F:223:LYS:C	2:F:226:ILE:H	1.99	0.70
2:E:31:LEU:HD11	2:E:74:ALA:HB2	1.73	0.70
2:E:152:PRO:C	2:E:156:ARG:HB2	2.16	0.69
1:D:37:LEU:HD23	1:D:37:LEU:N	2.07	0.69
2:E:153:SER:O	2:E:154:ALA:C	2.12	0.69
1:D:64:LEU:HD23	1:D:74:ALA:CB	2.21	0.69
2:F:103:LEU:CD1	2:F:247:VAL:HG13	2.20	0.69
2:E:153:SER:CB	2:E:157:GLN:HE21	2.05	0.68
1:A:137:LEU:HD12	1:A:147:ILE:CD1	2.23	0.68
2:E:153:SER:CB	2:E:157:GLN:H	2.07	0.68
1:D:28:LYS:HZ2	1:D:30:ASN:ND2	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:247:VAL:HG22	4:F:2025:HOH:O	1.92	0.68
1:A:36:ARG:NH2	1:A:170:GLU:O	2.27	0.68
1:C:73:LYS:HA	1:C:76:VAL:HG12	1.75	0.68
2:F:223:LYS:C	2:F:226:ILE:HG13	2.19	0.68
2:F:94:LYS:HE2	2:F:94:LYS:HA	1.76	0.67
1:C:10:GLY:HA2	1:C:173:TYR:CE1	2.30	0.67
2:F:390:ILE:O	2:F:393:ARG:HB2	1.95	0.67
1:D:115:PRO:HG3	1:D:120:ILE:HG12	1.76	0.67
2:F:96:VAL:HG21	2:F:280:ASP:HB3	1.77	0.67
2:E:220:ASP:O	2:E:224:LEU:HD22	1.95	0.66
1:D:60:PHE:HB2	1:D:78:LEU:HD22	1.77	0.66
1:D:12:VAL:HG23	4:D:2018:HOH:O	1.95	0.66
2:F:312:ILE:HG13	4:F:2032:HOH:O	1.95	0.66
2:E:358:LEU:O	2:E:361:THR:HB	1.96	0.65
2:F:108:VAL:HG12	2:F:109:LYS:N	2.11	0.65
2:E:263:LYS:HE3	4:E:2064:HOH:O	1.97	0.65
2:E:432:LEU:H	2:E:432:LEU:CD1	2.09	0.65
1:B:170:GLU:HG2	1:B:171:LEU:H	1.61	0.65
2:E:389:ASN:HD22	2:E:390:ILE:N	1.93	0.65
1:B:85:ASP:CG	1:B:88:LEU:HD13	2.22	0.65
2:F:384:ASN:ND2	2:F:394:ARG:HE	1.95	0.65
1:C:79:ALA:HB1	1:C:110:GLY:HA2	1.79	0.64
1:D:28:LYS:NZ	1:D:30:ASN:ND2	2.45	0.64
2:E:389:ASN:HD22	2:E:391:GLY:H	1.44	0.64
2:E:136:ILE:HG22	2:E:136:ILE:O	1.97	0.64
1:B:36:ARG:O	1:B:37:LEU:HD23	1.96	0.64
2:F:41:ASN:ND2	2:F:44:LEU:H	1.95	0.64
1:D:6:VAL:HG21	1:D:147:ILE:HG22	1.80	0.64
2:E:153:SER:C	2:E:157:GLN:N	2.54	0.64
2:E:89:VAL:HA	2:E:93:GLY:N	2.12	0.64
2:E:122:ARG:CZ	2:E:126:LEU:HD21	2.28	0.64
1:D:1:THR:HB	1:D:33:LYS:NZ	2.14	0.63
2:E:218:ILE:C	2:E:221:ALA:N	2.56	0.63
2:E:41:ASN:ND2	2:E:44:LEU:H	1.97	0.63
2:F:223:LYS:HA	2:F:226:ILE:HD11	1.81	0.63
1:C:76:VAL:HA	1:C:112:VAL:HG21	1.80	0.63
2:E:131:ILE:CG1	2:E:218:ILE:HD11	2.27	0.63
2:F:41:ASN:C	2:F:41:ASN:HD22	2.05	0.63
2:F:116:ILE:O	2:F:120:ARG:HB2	1.98	0.63
1:B:17:ASP:O	1:B:33:LYS:HG3	1.99	0.62
2:E:107:ALA:O	2:E:111:VAL:HG22	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:104:THR:HG21	2:E:292:THR:HG21	1.81	0.62
2:E:432:LEU:HD12	2:E:432:LEU:N	2.13	0.62
1:B:8:ARG:NH1	1:B:142:LEU:O	2.31	0.62
1:D:105:ILE:HD11	1:D:120:ILE:HG23	1.82	0.62
2:E:41:ASN:HD21	2:E:44:LEU:H	1.45	0.62
1:D:30:ASN:HD22	1:D:30:ASN:N	1.97	0.62
1:A:28:LYS:HZ2	1:A:30:ASN:HD21	1.45	0.62
1:D:14:ILE:HG13	1:D:43:ILE:HD12	1.81	0.62
2:E:153:SER:C	2:E:155:ALA:N	2.55	0.62
2:F:27:VAL:CG1	2:F:70:LEU:HG	2.30	0.62
2:E:115:ALA:HA	4:E:2032:HOH:O	1.99	0.61
1:A:15:ALA:HB1	1:A:152:LEU:HD12	1.82	0.61
1:C:60:PHE:HB2	1:C:78:LEU:HD22	1.81	0.61
2:F:81:VAL:HG11	2:F:99:ILE:HG12	1.82	0.61
2:F:238:GLU:O	2:F:242:ASP:CG	2.44	0.61
2:E:33:ASN:HD22	2:E:36:ARG:HD2	1.66	0.61
1:D:34:VAL:HB	1:D:167:THR:HG22	1.81	0.61
2:F:128:GLU:CB	2:F:222:MET:CE	2.61	0.61
1:D:15:ALA:HB2	1:D:168:ILE:HG23	1.83	0.60
2:E:37:ARG:HH11	2:E:37:ARG:CB	2.12	0.60
2:F:355:TYR:CD1	2:F:403:MET:HE2	2.36	0.60
1:B:10:GLY:HA2	1:B:173:TYR:CZ	2.37	0.60
2:E:160:ARG:HH12	2:E:164:ARG:NH2	1.94	0.60
2:E:401:ARG:CZ	2:E:442:ILE:HG23	2.32	0.60
2:F:224:LEU:C	2:F:228:GLU:HB2	2.17	0.60
2:E:220:ASP:C	2:E:221:ALA:N	2.60	0.59
2:E:244:ILE:HD11	2:E:294:HIS:O	2.02	0.59
1:B:95:LEU:HD12	1:B:95:LEU:N	2.17	0.59
2:E:130:ARG:NH2	2:E:130:ARG:HB2	2.17	0.59
2:F:108:VAL:C	2:F:110:MET:H	2.10	0.59
2:E:122:ARG:O	2:E:126:LEU:HD23	2.02	0.59
2:E:58:PRO:HG2	2:E:61:VAL:HG11	1.85	0.59
2:E:123:ALA:O	2:E:127:ALA:HB2	2.03	0.59
2:E:235:ASN:CG	2:E:238:GLU:HB2	2.27	0.59
2:F:21:ASP:OD2	2:F:25:ARG:HD2	2.01	0.59
1:A:30:ASN:C	1:A:30:ASN:HD22	2.11	0.59
2:F:315:PRO:O	2:F:318:LEU:HB2	2.03	0.58
1:A:13:VAL:HG12	1:A:170:GLU:HG3	1.85	0.58
1:D:71:LEU:HD11	1:D:97:VAL:HG12	1.85	0.58
2:E:115:ALA:C	2:E:117:GLU:H	2.11	0.58
2:F:76:ALA:HB1	2:F:250:HIS:O	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:106:ALA:O	2:F:110:MET:HB2	2.04	0.58
1:B:38:TYR:CE1	1:B:64:LEU:HB3	2.37	0.58
1:A:166:HIS:HE1	4:A:2028:HOH:O	1.86	0.58
2:F:249:GLN:HG3	4:F:2029:HOH:O	2.03	0.58
1:B:11:HIS:CE1	1:B:172:SER:OG	2.57	0.58
2:E:96:VAL:HG12	2:E:284:LEU:HD11	1.85	0.58
2:F:123:ALA:HB1	2:F:226:ILE:O	2.04	0.58
1:B:146:GLU:O	1:B:150:LYS:HG3	2.03	0.58
2:E:426:SER:HB3	2:E:430:ASP:OD1	2.04	0.58
1:B:17:ASP:HA	1:B:165:PHE:O	2.03	0.57
2:E:40:LEU:O	2:E:45:ARG:NH1	2.37	0.57
2:F:31:LEU:HD11	2:F:74:ALA:HB2	1.85	0.57
2:F:229:GLU:OE2	2:F:232:LYS:HD3	2.04	0.57
2:E:355:TYR:HE2	2:E:400:GLU:OE2	1.87	0.57
1:A:17:ASP:O	1:A:33:LYS:HD2	2.04	0.57
1:C:83:ARG:HA	4:C:2014:HOH:O	2.04	0.57
2:E:223:LYS:HD2	2:E:223:LYS:H	1.67	0.57
1:A:63:LYS:HD2	1:A:77:GLU:HB3	1.87	0.57
1:C:95:LEU:HB2	1:C:106:ILE:HB	1.87	0.57
2:E:131:ILE:CD1	2:E:218:ILE:HG12	2.35	0.57
1:D:73:LYS:HA	1:D:76:VAL:HG12	1.87	0.57
1:D:143:SER:OG	1:D:146:GLU:HG3	2.05	0.57
1:C:30:ASN:HD22	1:C:30:ASN:N	1.85	0.56
1:D:8:ARG:NH1	1:D:142:LEU:O	2.38	0.56
2:E:33:ASN:ND2	2:E:36:ARG:HD2	2.19	0.56
2:F:72:LYS:HD2	4:F:2017:HOH:O	2.05	0.56
2:E:390:ILE:O	2:E:393:ARG:HB2	2.05	0.56
2:F:362:GLU:HG3	2:F:411:SER:HA	1.87	0.56
1:A:44:ALA:HB2	1:A:97:VAL:HG23	1.87	0.56
1:B:85:ASP:HB3	1:B:88:LEU:HB2	1.86	0.56
1:C:37:LEU:HD13	1:C:57:PHE:HB3	1.86	0.56
2:E:393:ARG:HD3	4:E:2082:HOH:O	2.06	0.56
2:F:223:LYS:O	2:F:226:ILE:HG13	2.04	0.56
2:F:257:GLU:O	2:F:257:GLU:HG3	2.05	0.56
2:F:389:ASN:HD22	2:F:391:GLY:H	1.50	0.56
2:E:408:TYR:HB2	2:F:29:ILE:HD11	1.87	0.56
1:C:1:THR:HB	1:C:33:LYS:NZ	2.21	0.56
2:F:344:LEU:CD1	2:F:351:ILE:HD11	2.34	0.56
1:D:39:ASN:ND2	4:D:2010:HOH:O	2.38	0.56
2:F:223:LYS:HD2	2:F:223:LYS:H	1.67	0.56
2:E:111:VAL:O	2:E:115:ALA:HB2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:64:LEU:HB3	1:C:69:GLY:HA2	1.88	0.56
2:E:256:ASP:O	2:E:257:GLU:HG2	2.05	0.56
2:F:96:VAL:HG11	2:F:281:LEU:HD12	1.88	0.56
1:B:105:ILE:HD11	1:B:120:ILE:HG23	1.89	0.55
1:D:76:VAL:HA	1:D:112:VAL:HG21	1.87	0.55
2:E:96:VAL:CG1	2:E:281:LEU:HD12	2.36	0.55
2:E:116:ILE:HG22	2:E:116:ILE:O	2.06	0.55
1:D:79:ALA:HB1	1:D:110:GLY:HA2	1.87	0.55
2:E:108:VAL:HG12	2:E:109:LYS:N	2.21	0.55
2:E:151:GLU:N	2:E:152:PRO:HD2	2.21	0.55
2:E:152:PRO:O	2:E:155:ALA:N	2.40	0.55
1:B:28:LYS:HZ1	1:B:30:ASN:ND2	2.03	0.55
1:C:71:LEU:HD11	1:C:97:VAL:HG12	1.89	0.55
2:F:255:ILE:HD13	2:F:281:LEU:HD21	1.87	0.55
1:A:47:ALA:HB3	1:A:94:LEU:HB2	1.89	0.55
1:B:98:ALA:HB2	1:B:103:SER:HA	1.88	0.54
1:C:10:GLY:HA2	1:C:173:TYR:CZ	2.42	0.54
2:E:375:ARG:CZ	2:E:422:ALA:HB1	2.37	0.54
2:F:389:ASN:HD21	2:F:391:GLY:H	1.55	0.54
1:A:17:ASP:CG	1:A:162:THR:HG23	2.32	0.54
2:F:35:TRP:O	2:F:39:GLN:HG2	2.07	0.54
1:B:36:ARG:NE	1:B:169:GLU:OE2	2.40	0.54
1:C:5:SER:HB2	1:C:14:ILE:HG12	1.89	0.54
2:E:366:ILE:HD13	2:E:418:ILE:HB	1.89	0.54
2:E:153:SER:C	2:E:156:ARG:N	2.65	0.54
2:E:257:GLU:HG3	2:E:257:GLU:O	2.06	0.54
2:E:106:ALA:O	2:E:110:MET:HB2	2.07	0.54
2:F:40:LEU:O	2:F:45:ARG:NH1	2.41	0.54
1:A:137:LEU:HD12	1:A:147:ILE:HD12	1.89	0.54
1:B:59:LEU:O	1:B:62:ARG:HB3	2.08	0.54
1:D:86:ARG:O	1:D:90:LYS:HG2	2.07	0.54
1:D:63:LYS:HE2	1:D:77:GLU:HB3	1.90	0.53
1:D:152:LEU:HD13	1:D:166:HIS:CE1	2.43	0.53
2:E:222:MET:O	2:E:226:ILE:HG13	2.08	0.53
1:A:54:PHE:HE1	1:A:58:GLU:OE1	1.91	0.53
1:C:6:VAL:HG21	1:C:147:ILE:CG2	2.38	0.53
1:D:71:LEU:HB2	1:D:99:ASP:OD2	2.07	0.53
2:F:413:LEU:O	2:F:416:GLN:HG3	2.08	0.53
1:D:2:THR:HG21	1:D:155:ALA:CB	2.38	0.53
1:A:134:ARG:CD	4:A:2019:HOH:O	2.41	0.53
1:C:37:LEU:HD21	1:C:44:ALA:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:262:CYS:SG	2:E:318:LEU:HD13	2.49	0.53
2:F:17:ILE:N	2:F:17:ILE:HD12	2.23	0.53
2:F:89:VAL:HA	2:F:93:GLY:N	2.23	0.53
1:B:60:PHE:CE2	1:B:97:VAL:HG21	2.44	0.53
1:C:20:ALA:HB2	1:C:31:VAL:HG21	1.90	0.53
2:F:122:ARG:HG2	2:F:122:ARG:HH21	1.73	0.53
2:F:358:LEU:O	2:F:361:THR:HB	2.09	0.53
1:B:170:GLU:HG2	1:B:171:LEU:N	2.22	0.53
1:B:3:ILE:HB	1:B:122:ILE:CG1	2.39	0.53
1:B:34:VAL:HB	1:B:167:THR:HG22	1.90	0.53
1:C:152:LEU:HD22	1:C:166:HIS:CE1	2.44	0.53
1:D:71:LEU:HD21	1:D:97:VAL:CG1	2.39	0.53
1:D:174:LYS:HE2	1:D:174:LYS:HA	1.89	0.53
2:E:153:SER:C	2:E:154:ALA:HA	2.33	0.53
1:B:46:PHE:HB3	1:B:57:PHE:CZ	2.44	0.52
2:F:262:CYS:SG	2:F:318:LEU:HD13	2.48	0.52
1:C:12:VAL:O	1:C:12:VAL:HG13	2.08	0.52
1:C:117:ASN:O	1:C:118:ASP:HB2	2.08	0.52
1:D:56:LEU:HD13	1:D:95:LEU:HD11	1.91	0.52
1:C:34:VAL:HB	1:C:167:THR:HG22	1.90	0.52
1:D:41:LYS:O	1:D:171:LEU:HD21	2.10	0.52
1:A:28:LYS:NZ	1:A:30:ASN:ND2	2.54	0.52
2:F:65:GLU:CG	3:F:501:ANP:H2'	2.33	0.52
2:F:221:ALA:HB1	2:F:225:LEU:CD2	2.40	0.52
1:C:54:PHE:O	1:C:58:GLU:HB2	2.09	0.52
1:C:78:LEU:O	1:C:82:TRP:HB2	2.10	0.52
2:E:152:PRO:O	2:E:155:ALA:HB3	2.10	0.52
2:E:237:GLU:HG3	2:E:238:GLU:OE1	2.10	0.52
2:F:61:VAL:C	3:F:501:ANP:H2	2.34	0.52
2:E:153:SER:C	2:E:154:ALA:N	2.67	0.52
2:F:108:VAL:O	2:F:110:MET:N	2.42	0.52
2:F:247:VAL:CG2	4:F:2025:HOH:O	2.53	0.51
2:E:60:GLY:N	3:E:500:ANP:O2B	2.43	0.51
2:F:229:GLU:HA	2:F:232:LYS:CD	2.38	0.51
2:E:153:SER:O	2:E:157:GLN:HB2	2.11	0.51
2:E:220:ASP:C	2:E:221:ALA:CA	2.83	0.51
2:E:244:ILE:HG23	2:E:297:VAL:HG22	1.92	0.51
2:F:222:MET:O	2:F:226:ILE:CG1	2.59	0.51
2:F:432:LEU:N	2:F:432:LEU:HD12	2.26	0.51
1:B:98:ALA:CB	1:B:103:SER:HA	2.40	0.51
2:F:222:MET:O	2:F:226:ILE:HG13	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:124:SER:OG	1:C:159:CYS:SG	2.69	0.51
1:D:54:PHE:O	1:D:58:GLU:HB2	2.10	0.51
2:E:112:ARG:O	2:E:116:ILE:HG12	2.11	0.51
2:E:408:TYR:OH	2:F:10:VAL:HG21	2.11	0.51
2:F:297:VAL:HG12	2:F:298:LYS:N	2.26	0.51
1:C:30:ASN:ND2	1:C:30:ASN:N	2.50	0.51
2:E:220:ASP:C	2:E:221:ALA:HA	2.36	0.51
1:C:28:LYS:HD3	1:C:31:VAL:HG22	1.93	0.51
1:D:22:LEU:HD12	1:D:27:MET:HE2	1.91	0.51
2:F:104:THR:HA	4:F:2025:HOH:O	2.11	0.51
2:E:153:SER:CA	2:E:156:ARG:CA	2.83	0.50
2:E:312:ILE:HD12	2:E:313:ALA:H	1.76	0.50
2:F:107:ALA:O	2:F:111:VAL:HG22	2.11	0.50
1:D:64:LEU:HD23	1:D:74:ALA:HB3	1.92	0.50
2:E:65:GLU:CG	3:E:500:ANP:H2'	2.37	0.50
2:E:401:ARG:NH2	2:E:442:ILE:HG23	2.26	0.50
2:F:18:ILE:N	3:F:501:ANP:HN61	2.07	0.50
2:F:394:ARG:NH1	2:F:443:LEU:O	2.43	0.50
1:C:36:ARG:NH1	1:C:40:ASP:HB3	2.27	0.50
2:E:138:PRO:C	2:E:139:ALA:O	2.52	0.50
2:E:240:LYS:HE3	2:E:294:HIS:O	2.10	0.50
2:E:337:THR:O	2:E:341:GLU:HG3	2.11	0.50
2:E:315:PRO:O	2:E:318:LEU:HB2	2.12	0.50
2:E:442:ILE:O	2:E:442:ILE:HG22	2.11	0.50
2:F:31:LEU:HD22	4:F:2007:HOH:O	2.11	0.50
1:A:34:VAL:HB	1:A:167:THR:HG22	1.93	0.50
1:C:5:SER:HB3	1:C:120:ILE:HB	1.92	0.50
1:D:18:GLY:O	1:D:31:VAL:HG23	2.10	0.50
1:D:85:ASP:HB3	1:D:88:LEU:HB2	1.94	0.50
1:A:117:ASN:O	1:A:118:ASP:HB2	2.11	0.50
2:E:64:THR:HB	3:E:500:ANP:O2A	2.12	0.50
2:E:64:THR:HG21	2:E:68:ARG:NH1	2.27	0.50
2:E:131:ILE:HD11	2:E:218:ILE:HG12	1.90	0.49
2:F:312:ILE:CD1	2:F:313:ALA:N	2.75	0.49
1:A:17:ASP:HA	1:A:165:PHE:O	2.12	0.49
1:B:7:ARG:HG3	1:B:12:VAL:HB	1.93	0.49
1:D:46:PHE:HA	1:D:94:LEU:O	2.12	0.49
1:D:133:ALA:O	1:D:137:LEU:HB2	2.12	0.49
1:B:36:ARG:NH2	1:B:170:GLU:O	2.45	0.49
2:E:35:TRP:O	2:E:39:GLN:HG2	2.12	0.49
2:F:123:ALA:C	2:F:127:ALA:HB3	2.38	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:ILE:HG13	4:B:2025:HOH:O	2.12	0.49
2:E:131:ILE:HD12	2:E:218:ILE:CG1	2.43	0.49
2:F:359:MET:HG3	2:F:366:ILE:HG13	1.95	0.49
2:F:384:ASN:HD21	2:F:390:ILE:HG12	1.77	0.49
2:E:335:LEU:O	2:E:381:TRP:HZ3	1.96	0.49
2:F:118:LYS:O	2:F:119:ASN:ND2	2.46	0.49
1:C:37:LEU:HA	4:C:2008:HOH:O	2.12	0.49
2:F:108:VAL:C	2:F:110:MET:N	2.66	0.49
2:F:109:LYS:O	2:F:113:VAL:HG23	2.13	0.49
2:F:223:LYS:O	2:F:226:ILE:CB	2.61	0.49
2:E:17:ILE:HD12	2:E:17:ILE:N	2.28	0.49
2:F:299:THR:O	2:F:302:ILE:HG13	2.13	0.49
1:A:95:LEU:HB2	1:A:106:ILE:HB	1.95	0.49
1:D:95:LEU:HD12	1:D:95:LEU:N	2.28	0.48
2:E:220:ASP:HB3	2:E:224:LEU:HD21	1.95	0.48
2:E:411:SER:C	2:E:413:LEU:H	2.21	0.48
2:F:299:THR:HA	2:F:302:ILE:HG13	1.94	0.48
1:D:121:ALA:HB1	1:D:126:GLY:O	2.13	0.48
2:E:122:ARG:HH11	2:E:126:LEU:CD2	2.26	0.48
1:B:88:LEU:N	1:B:88:LEU:HD12	2.28	0.48
1:C:86:ARG:O	1:C:90:LYS:HG2	2.13	0.48
1:C:152:LEU:O	1:C:152:LEU:HD23	2.14	0.48
1:D:159:CYS:HB3	1:D:162:THR:HB	1.94	0.48
2:E:16:HIS:HB2	2:E:17:ILE:HD12	1.95	0.48
2:F:37:ARG:HH11	2:F:37:ARG:HB3	1.78	0.48
1:B:1:THR:HB	1:B:33:LYS:NZ	2.29	0.48
1:D:3:ILE:HB	1:D:122:ILE:CG1	2.43	0.48
2:E:122:ARG:NH1	2:E:126:LEU:CD2	2.75	0.48
2:E:435:ASP:OD1	2:E:438:LEU:HD12	2.13	0.48
2:F:123:ALA:HB2	2:F:230:ALA:HB2	1.95	0.48
2:E:109:LYS:O	2:E:113:VAL:HG23	2.14	0.48
1:D:7:ARG:NH1	1:D:12:VAL:HB	2.29	0.48
2:F:12:GLU:HG2	2:F:73:LEU:HD13	1.95	0.48
2:E:248:GLU:HG2	2:E:297:VAL:HG13	1.95	0.48
2:E:344:LEU:HD23	2:E:373:ILE:HG23	1.96	0.48
2:F:117:GLU:OE1	2:F:120:ARG:CZ	2.62	0.48
1:A:1:THR:HB	1:A:33:LYS:HZ1	1.78	0.48
1:D:28:LYS:HG2	1:D:30:ASN:HD22	1.79	0.48
2:E:231:ALA:O	2:E:234:VAL:HG12	2.13	0.47
2:F:234:VAL:CG1	2:F:239:LEU:HD13	2.44	0.47
3:E:500:ANP:O4'	3:E:500:ANP:N3	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:117:GLU:OE1	2:F:120:ARG:NH1	2.47	0.47
2:F:221:ALA:HB1	2:F:225:LEU:HD23	1.96	0.47
2:E:23:ALA:HA	2:E:330:VAL:HG21	1.95	0.47
2:E:68:ARG:HG2	4:E:2020:HOH:O	2.13	0.47
2:F:127:ALA:HB2	2:F:229:GLU:HG3	1.95	0.47
1:D:30:ASN:ND2	1:D:30:ASN:N	2.61	0.47
2:E:72:LYS:HD2	4:E:2021:HOH:O	2.14	0.47
1:B:141:GLU:HA	1:B:141:GLU:OE2	2.14	0.47
2:F:63:LYS:HG2	2:F:332:LEU:HD13	1.95	0.47
2:E:160:ARG:O	2:E:163:LEU:HB3	2.13	0.47
1:D:101:THR:OG1	1:D:102:ALA:N	2.48	0.47
2:E:126:LEU:O	2:E:130:ARG:NH2	2.47	0.47
2:E:136:ILE:HG22	4:E:2036:HOH:O	2.14	0.47
2:F:61:VAL:C	3:F:501:ANP:C2	2.88	0.47
2:F:115:ALA:HA	4:F:2026:HOH:O	2.13	0.47
1:D:17:ASP:HA	1:D:165:PHE:O	2.15	0.47
2:E:130:ARG:HB2	2:E:130:ARG:CZ	2.45	0.47
2:E:389:ASN:HD21	2:E:391:GLY:H	1.54	0.47
1:C:149:GLU:CD	1:C:168:ILE:HD11	2.40	0.47
2:E:128:GLU:HG2	2:E:222:MET:HE1	1.97	0.47
2:E:152:PRO:O	2:E:155:ALA:CA	2.63	0.47
1:C:36:ARG:NH1	1:C:40:ASP:O	2.39	0.47
1:D:28:LYS:HD3	1:D:31:VAL:HG22	1.96	0.47
2:F:244:ILE:HD11	2:F:294:HIS:O	2.14	0.46
2:E:4:MET:HE1	2:E:73:LEU:HD11	1.97	0.46
2:F:112:ARG:O	2:F:115:ALA:HB3	2.14	0.46
1:B:17:ASP:CG	1:B:162:THR:HG23	2.40	0.46
2:F:41:ASN:ND2	2:F:41:ASN:C	2.70	0.46
2:F:108:VAL:CG1	2:F:109:LYS:N	2.78	0.46
2:F:127:ALA:HB1	2:F:229:GLU:HB3	1.97	0.46
1:A:138:GLU:O	1:A:139:ASN:CG	2.58	0.46
1:D:60:PHE:CZ	1:D:97:VAL:HG11	2.50	0.46
1:D:127:PRO:HG3	4:D:2022:HOH:O	2.15	0.46
1:C:17:ASP:CG	1:C:162:THR:HG23	2.41	0.46
1:C:116:GLU:HB3	1:C:117:ASN:ND2	2.30	0.46
2:E:52:ASN:HB2	2:E:325:ARG:O	2.16	0.46
2:E:152:PRO:O	2:E:155:ALA:C	2.50	0.46
2:E:267:SER:HB3	2:E:270:PRO:CG	2.46	0.46
2:F:375:ARG:CZ	2:F:422:ALA:HB1	2.45	0.46
3:F:501:ANP:N3	3:F:501:ANP:H5'1	2.29	0.46
1:A:34:VAL:HB	1:A:167:THR:CG2	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:PHE:CE2	1:A:97:VAL:HG21	2.51	0.46
1:B:60:PHE:CD1	1:B:78:LEU:HD22	2.50	0.46
2:E:406:ILE:O	2:E:410:ALA:N	2.47	0.46
1:A:59:LEU:HD23	1:A:78:LEU:CD1	2.46	0.46
2:E:130:ARG:HD2	2:E:225:LEU:HD11	1.98	0.46
2:E:362:GLU:HG2	2:E:410:ALA:CB	2.46	0.46
1:A:82:TRP:CD2	1:A:108:GLY:HA2	2.51	0.46
1:A:59:LEU:O	1:A:62:ARG:HB3	2.16	0.46
1:B:19:GLN:HB2	1:B:163:ASN:ND2	2.31	0.46
1:D:64:LEU:HD23	1:D:74:ALA:HB2	1.96	0.46
1:D:114:GLN:HE21	1:D:114:GLN:HB2	1.49	0.46
2:E:293:LYS:HG3	2:E:293:LYS:O	2.16	0.46
2:E:361:THR:HG21	2:F:36:ARG:HA	1.98	0.46
2:E:382:GLN:O	2:E:386:SER:HB3	2.16	0.46
2:E:432:LEU:CD1	2:E:432:LEU:N	2.76	0.46
2:E:360:ALA:O	2:E:361:THR:C	2.58	0.45
1:A:87:MET:CE	1:B:84:THR:HG23	2.41	0.45
1:D:1:THR:HB	1:D:33:LYS:HZ2	1.81	0.45
2:E:312:ILE:CD1	2:E:313:ALA:H	2.30	0.45
1:C:136:LEU:C	1:C:138:GLU:H	2.25	0.45
2:E:138:PRO:O	2:E:139:ALA:C	2.53	0.45
2:F:60:GLY:H	2:F:393:ARG:NH2	2.14	0.45
1:B:71:LEU:HG	1:B:99:ASP:OD1	2.16	0.45
1:D:17:ASP:O	1:D:33:LYS:HD2	2.17	0.45
2:E:53:ILE:HG22	2:E:54:LEU:N	2.32	0.45
1:B:148:ALA:O	1:B:152:LEU:HB2	2.16	0.45
1:D:1:THR:HB	1:D:33:LYS:HZ3	1.81	0.45
2:E:37:ARG:HH11	2:E:37:ARG:CG	2.29	0.45
1:B:136:LEU:HB3	1:B:147:ILE:HG12	1.99	0.45
2:E:153:SER:HB2	2:E:157:GLN:CD	2.37	0.45
2:E:240:LYS:O	2:E:240:LYS:HG2	2.17	0.45
2:F:323:GLN:HB3	4:F:2045:HOH:O	2.17	0.45
2:E:128:GLU:O	2:E:131:ILE:HG22	2.16	0.45
2:E:344:LEU:HD12	2:E:351:ILE:HD11	1.99	0.45
1:A:149:GLU:HG2	1:A:168:ILE:HD11	1.98	0.45
2:F:16:HIS:C	2:F:17:ILE:HD12	2.42	0.45
1:A:72:VAL:O	1:A:76:VAL:HG23	2.16	0.45
2:E:362:GLU:HG2	2:E:410:ALA:HB1	1.98	0.45
2:E:366:ILE:HD13	2:E:366:ILE:HA	1.87	0.45
2:E:442:ILE:O	2:E:442:ILE:CG2	2.65	0.45
1:B:51:ALA:O	1:B:52:ASP:C	2.59	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:85:ASP:HB3	1:C:88:LEU:HB2	1.99	0.45
2:F:102:ASP:O	2:F:103:LEU:C	2.60	0.45
2:F:122:ARG:HH11	2:F:126:LEU:CD2	2.28	0.45
2:F:350:SER:OG	2:F:353:VAL:HG23	2.17	0.45
1:A:137:LEU:HD12	1:A:147:ILE:HD13	1.98	0.44
1:D:71:LEU:HD21	1:D:97:VAL:HG13	1.98	0.44
1:B:145:ARG:O	1:B:146:GLU:C	2.60	0.44
2:E:335:LEU:HD22	2:E:339:ASP:HB3	1.98	0.44
2:F:123:ALA:O	2:F:127:ALA:HB3	2.17	0.44
1:B:70:HIS:O	1:B:71:LEU:C	2.60	0.44
1:D:28:LYS:HG2	1:D:30:ASN:ND2	2.32	0.44
2:E:441:PHE:HA	2:F:315:PRO:HG2	1.99	0.44
1:C:8:ARG:NH1	1:C:142:LEU:O	2.50	0.44
2:E:270:PRO:O	2:E:274:ARG:HD2	2.18	0.44
2:E:292:THR:HA	4:E:2057:HOH:O	2.18	0.44
2:F:243:ALA:O	2:F:244:ILE:C	2.60	0.44
1:A:65:GLU:HA	1:A:65:GLU:OE1	2.18	0.44
2:E:21:ASP:O	2:E:25:ARG:HG3	2.18	0.44
2:E:135:LEU:HD13	2:E:159:PHE:CD2	2.53	0.44
2:F:61:VAL:CA	3:F:501:ANP:H2	2.47	0.44
2:F:406:ILE:O	2:F:410:ALA:N	2.48	0.44
1:C:17:ASP:OD2	1:C:162:THR:HA	2.17	0.44
2:E:160:ARG:NH1	2:E:164:ARG:NH2	2.61	0.44
2:F:232:LYS:HD2	2:F:232:LYS:N	2.33	0.44
2:F:360:ALA:O	2:F:361:THR:C	2.61	0.44
2:F:426:SER:HB3	2:F:430:ASP:OD1	2.18	0.44
1:B:34:VAL:HB	1:B:167:THR:CG2	2.47	0.44
1:B:82:TRP:CD2	1:B:108:GLY:HA2	2.53	0.44
1:B:170:GLU:CG	1:B:171:LEU:H	2.28	0.44
1:D:7:ARG:HB2	4:D:2018:HOH:O	2.17	0.44
2:F:223:LYS:O	2:F:226:ILE:CG1	2.66	0.44
2:F:362:GLU:HG2	2:F:410:ALA:CB	2.48	0.44
1:A:66:MET:C	1:A:67:HIS:ND1	2.75	0.44
1:B:64:LEU:HD23	1:B:74:ALA:CB	2.47	0.44
2:E:312:ILE:HD12	2:E:312:ILE:N	2.26	0.44
2:E:384:ASN:HD21	2:E:390:ILE:HG12	1.83	0.44
1:C:117:ASN:ND2	4:C:2017:HOH:O	2.51	0.43
1:D:36:ARG:HH11	1:D:36:ARG:HB3	1.82	0.43
1:D:104:LEU:HD22	1:D:112:VAL:CG1	2.47	0.43
2:F:345:THR:HG22	2:F:352:THR:HG21	1.99	0.43
1:A:15:ALA:CB	1:A:152:LEU:HD12	2.45	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:136:LEU:HB3	1:D:147:ILE:HG12	2.00	0.43
2:E:243:ALA:O	2:E:247:VAL:HG23	2.17	0.43
1:B:46:PHE:CD2	1:B:53:ALA:HB2	2.53	0.43
1:C:71:LEU:HD21	1:C:97:VAL:CG1	2.49	0.43
2:E:359:MET:HE1	2:F:36:ARG:HH12	1.81	0.43
2:E:412:ASP:HB2	2:F:7:ARG:HH21	1.83	0.43
1:B:14:ILE:HD12	1:B:44:ALA:HA	2.01	0.43
1:C:148:ALA:O	1:C:152:LEU:HB2	2.18	0.43
2:F:223:LYS:C	2:F:225:LEU:H	2.13	0.43
1:B:46:PHE:HB3	1:B:57:PHE:HZ	1.82	0.43
2:E:264:ARG:HA	4:E:2052:HOH:O	2.18	0.43
2:F:432:LEU:H	2:F:432:LEU:CD1	2.31	0.43
1:D:99:ASP:HA	1:D:171:LEU:CD2	2.48	0.43
2:E:292:THR:C	2:E:294:HIS:H	2.24	0.43
1:A:154:ILE:O	1:A:155:ALA:C	2.62	0.43
1:C:1:THR:HB	1:C:33:LYS:HZ3	1.82	0.43
2:F:384:ASN:ND2	2:F:390:ILE:H	2.16	0.43
1:A:36:ARG:NH1	1:A:40:ASP:O	2.51	0.43
1:B:85:ASP:OD1	1:B:87:MET:HB2	2.19	0.43
2:E:348:ASN:O	2:E:349:ALA:C	2.61	0.43
1:C:11:HIS:CE1	1:C:174:LYS:HE2	2.54	0.43
2:E:122:ARG:HH11	2:E:126:LEU:HD21	1.79	0.43
2:E:274:ARG:HA	4:E:2026:HOH:O	2.19	0.43
2:F:290:VAL:HG12	2:F:291:SER:N	2.34	0.43
1:D:10:GLY:O	1:D:172:SER:HA	2.19	0.43
2:E:22:ASN:HD22	2:E:22:ASN:HA	1.64	0.43
2:F:127:ALA:CB	2:F:229:GLU:HB3	2.49	0.43
2:F:322:LEU:HD12	2:F:322:LEU:HA	1.81	0.43
1:A:30:ASN:C	1:A:30:ASN:ND2	2.74	0.42
1:B:54:PHE:HE1	1:B:58:GLU:OE1	2.02	0.42
1:B:21:THR:HG22	1:B:22:LEU:N	2.34	0.42
1:B:60:PHE:O	1:B:61:GLU:C	2.62	0.42
1:D:15:ALA:HA	1:D:167:THR:O	2.18	0.42
1:C:136:LEU:C	1:C:138:GLU:N	2.76	0.42
2:E:151:GLU:C	2:E:153:SER:N	2.76	0.42
2:F:32:ARG:O	2:F:36:ARG:HG3	2.19	0.42
2:F:257:GLU:OE1	2:F:260:LYS:HD2	2.19	0.42
2:F:366:ILE:HA	2:F:366:ILE:HD13	1.81	0.42
1:B:32:LYS:HB2	1:B:32:LYS:HE3	1.69	0.42
2:E:115:ALA:C	2:E:117:GLU:N	2.77	0.42
2:F:88:GLU:HB2	4:F:2021:HOH:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:122:ARG:HG2	2:F:122:ARG:NH2	2.33	0.42
2:F:221:ALA:O	2:F:225:LEU:CB	2.68	0.42
1:D:174:LYS:HE2	1:D:174:LYS:CA	2.48	0.42
2:F:74:ALA:O	2:F:75:ASN:C	2.63	0.42
2:F:292:THR:C	2:F:294:HIS:H	2.27	0.42
1:D:30:ASN:H	1:D:30:ASN:ND2	2.09	0.42
2:E:344:LEU:HD13	2:E:395:LEU:HD22	2.01	0.42
2:F:101:ARG:O	2:F:104:THR:HB	2.20	0.42
1:A:78:LEU:O	1:A:82:TRP:HB2	2.20	0.42
1:B:5:SER:HA	1:B:13:VAL:O	2.19	0.42
1:D:100:GLU:HG3	1:D:173:TYR:HB2	2.00	0.42
1:D:134:ARG:NE	4:D:2023:HOH:O	2.53	0.42
1:D:157:ASP:OD2	1:D:164:HIS:NE2	2.52	0.42
1:B:95:LEU:N	1:B:95:LEU:CD1	2.82	0.42
1:C:134:ARG:CD	4:C:2017:HOH:O	2.67	0.42
1:C:152:LEU:HD22	1:C:166:HIS:HE1	1.84	0.42
1:B:170:GLU:CG	1:B:171:LEU:N	2.83	0.42
1:C:3:ILE:HD11	1:C:46:PHE:O	2.20	0.42
1:D:44:ALA:HA	1:D:96:ALA:O	2.20	0.42
2:E:76:ALA:HA	2:E:77:PRO:HD3	1.91	0.42
2:E:158:ALA:O	2:E:162:LYS:HG3	2.20	0.42
1:C:68:GLN:O	1:C:70:HIS:N	2.52	0.42
2:E:81:VAL:HG11	2:E:99:ILE:HG12	2.01	0.42
2:F:243:ALA:O	2:F:246:ALA:N	2.52	0.42
2:F:282:LEU:HA	2:F:282:LEU:HD12	1.80	0.42
2:E:27:VAL:HG13	2:E:70:LEU:HG	2.02	0.41
2:E:54:LEU:HB3	2:E:329:ARG:HD3	2.02	0.41
1:A:149:GLU:OE1	1:A:168:ILE:HD11	2.21	0.41
1:C:36:ARG:C	1:C:37:LEU:HD23	2.45	0.41
1:C:56:LEU:HD21	1:C:91:LEU:HD11	2.02	0.41
2:E:61:VAL:C	3:E:500:ANP:C2	2.93	0.41
2:E:394:ARG:NH1	2:E:443:LEU:O	2.48	0.41
2:F:96:VAL:CG1	2:F:281:LEU:HD12	2.50	0.41
2:F:97:ASP:OD1	2:F:97:ASP:N	2.52	0.41
2:F:344:LEU:HD12	2:F:344:LEU:HA	1.82	0.41
2:F:387:THR:OG1	2:F:388:GLU:N	2.54	0.41
3:F:501:ANP:N3	3:F:501:ANP:O4'	2.53	0.41
1:A:8:ARG:O	1:A:11:HIS:HB2	2.20	0.41
1:B:72:VAL:O	1:B:75:ALA:HB3	2.19	0.41
1:D:21:THR:HG22	1:D:22:LEU:N	2.36	0.41
2:E:34:ARG:NH1	2:E:301:HIS:O	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:55:MET:O	2:E:307:SER:HA	2.20	0.41
2:E:152:PRO:C	2:E:156:ARG:N	2.57	0.41
2:F:235:ASN:HB2	2:F:238:GLU:OE2	2.20	0.41
1:A:14:ILE:HG13	1:A:43:ILE:HD12	2.01	0.41
1:D:11:HIS:CE1	1:D:172:SER:OG	2.73	0.41
2:E:152:PRO:HB2	2:E:156:ARG:HB2	2.02	0.41
2:F:362:GLU:HG2	2:F:410:ALA:C	2.45	0.41
1:B:64:LEU:HD23	1:B:74:ALA:HB2	2.02	0.41
2:F:240:LYS:O	2:F:240:LYS:HG2	2.21	0.41
1:B:7:ARG:NE	1:B:118:ASP:OD2	2.41	0.41
1:C:73:LYS:HD2	1:C:76:VAL:CG1	2.50	0.41
1:D:148:ALA:O	1:D:152:LEU:HB2	2.20	0.41
2:E:120:ARG:O	2:E:124:GLU:HG2	2.20	0.41
2:F:291:SER:OG	2:F:296:MET:HE1	2.21	0.41
2:F:335:LEU:HD22	2:F:339:ASP:HB3	2.02	0.41
1:B:82:TRP:CE2	1:B:108:GLY:HA2	2.55	0.41
1:C:28:LYS:HG2	1:C:30:ASN:ND2	2.34	0.41
1:D:80:LYS:HE3	1:D:80:LYS:HB3	1.89	0.41
1:D:100:GLU:HG3	1:D:173:TYR:CB	2.51	0.41
2:E:269:GLY:N	2:E:270:PRO:HD2	2.35	0.41
2:E:351:ILE:HG23	4:E:2072:HOH:O	2.20	0.41
2:F:348:ASN:O	2:F:349:ALA:C	2.62	0.41
2:F:375:ARG:HB3	2:F:425:VAL:HG11	2.03	0.41
1:A:61:GLU:O	1:A:64:LEU:HB2	2.21	0.41
1:C:95:LEU:N	1:C:95:LEU:CD1	2.83	0.41
2:E:108:VAL:C	2:E:110:MET:H	2.29	0.41
2:F:257:GLU:O	2:F:257:GLU:CG	2.69	0.41
2:F:432:LEU:HD12	2:F:432:LEU:H	1.85	0.41
1:B:147:ILE:O	1:B:148:ALA:C	2.63	0.41
1:C:132:ALA:HB1	1:C:154:ILE:HD12	2.03	0.41
1:D:3:ILE:HB	1:D:122:ILE:HG12	2.02	0.41
1:D:22:LEU:CD1	1:D:27:MET:HE2	2.51	0.41
2:E:352:THR:HG22	2:E:353:VAL:N	2.36	0.41
2:F:222:MET:O	2:F:226:ILE:HG12	2.20	0.41
1:A:53:ALA:O	1:A:54:PHE:C	2.64	0.41
1:A:138:GLU:C	1:A:139:ASN:ND2	2.79	0.41
1:C:134:ARG:HD3	4:C:2017:HOH:O	2.21	0.41
2:E:257:GLU:O	2:E:257:GLU:CG	2.69	0.41
1:A:94:LEU:HD23	1:A:94:LEU:HA	1.87	0.40
1:B:5:SER:C	1:B:6:VAL:HG23	2.46	0.40
1:B:37:LEU:O	1:B:38:TYR:C	2.64	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:57:PHE:CE1	1:C:95:LEU:HG	2.56	0.40
1:D:99:ASP:OD1	1:D:99:ASP:C	2.64	0.40
1:B:56:LEU:HD13	1:B:95:LEU:HD11	2.03	0.40
1:B:152:LEU:HD23	1:B:152:LEU:O	2.21	0.40
2:F:96:VAL:CG2	2:F:280:ASP:HB3	2.49	0.40
2:E:63:LYS:HB2	3:E:500:ANP:O1B	2.22	0.40
2:F:121:TYR:O	2:F:125:GLU:HB2	2.22	0.40
2:F:357:ALA:O	2:F:360:ALA:HB3	2.20	0.40
2:E:329:ARG:HG3	2:E:329:ARG:HH11	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	172/175 (98%)	158 (92%)	13 (8%)	1 (1%)	21	51
1	B	172/175 (98%)	144 (84%)	23 (13%)	5 (3%)	3	13
1	C	172/175 (98%)	153 (89%)	16 (9%)	3 (2%)	7	25
1	D	172/175 (98%)	148 (86%)	22 (13%)	2 (1%)	10	34
2	E	398/449 (89%)	348 (87%)	41 (10%)	9 (2%)	5	18
2	F	399/449 (89%)	342 (86%)	43 (11%)	14 (4%)	3	10
All	All	1485/1598 (93%)	1293 (87%)	158 (11%)	34 (2%)	5	18

All (34) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	116	GLU
2	E	92	VAL
2	E	139	ALA

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Mol	Chain	Res	Type
2	F	92	VAL
2	F	130	ARG
2	F	139	ALA
2	F	230	ALA
1	B	101	THR
1	B	116	GLU
1	D	69	GLY
2	E	116	ILE
2	E	212	LYS
2	F	125	GLU
2	F	212	LYS
2	F	226	ILE
2	F	300	ASP
1	B	71	LEU
1	B	142	LEU
1	C	115	PRO
2	E	141	ASN
2	E	300	ASP
2	F	109	LYS
2	F	141	ASN
2	F	228	GLU
1	A	71	LEU
1	B	146	GLU
2	F	236	PRO
1	C	69	GLY
1	D	116	GLU
2	E	265	GLY
2	E	227	GLU
2	E	138	PRO
2	F	138	PRO
2	F	265	GLY

5.3.2 Protein sidechains [i](#)

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	136/136 (100%)	128 (94%)	8 (6%)	18	48
1	B	136/136 (100%)	124 (91%)	12 (9%)	9	29
1	C	136/136 (100%)	123 (90%)	13 (10%)	8	26
1	D	136/136 (100%)	122 (90%)	14 (10%)	7	23
2	E	340/383 (89%)	284 (84%)	56 (16%)	2	8
2	F	339/383 (88%)	281 (83%)	58 (17%)	2	7
All	All	1223/1310 (93%)	1062 (87%)	161 (13%)	4	14

All (161) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	THR
1	A	4	VAL
1	A	30	ASN
1	A	71	LEU
1	A	136	LEU
1	A	137	LEU
1	A	152	LEU
1	A	160	ILE
1	B	1	THR
1	B	4	VAL
1	B	22	LEU
1	B	30	ASN
1	B	72	VAL
1	B	95	LEU
1	B	107	THR
1	B	116	GLU
1	B	136	LEU
1	B	152	LEU
1	B	160	ILE
1	B	174	LYS
1	C	1	THR
1	C	4	VAL
1	C	30	ASN
1	C	37	LEU
1	C	58	GLU
1	C	71	LEU
1	C	72	VAL

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Mol	Chain	Res	Type
1	C	97	VAL
1	C	111	ASP
1	C	116	GLU
1	C	136	LEU
1	C	152	LEU
1	C	160	ILE
1	D	1	THR
1	D	4	VAL
1	D	30	ASN
1	D	36	ARG
1	D	37	LEU
1	D	72	VAL
1	D	91	LEU
1	D	97	VAL
1	D	111	ASP
1	D	114	GLN
1	D	116	GLU
1	D	136	LEU
1	D	152	LEU
1	D	160	ILE
2	E	1	HIS
2	E	11	SER
2	E	13	LEU
2	E	27	VAL
2	E	31	LEU
2	E	37	ARG
2	E	41	ASN
2	E	59	THR
2	E	68	ARG
2	E	70	LEU
2	E	94	LYS
2	E	103	LEU
2	E	108	VAL
2	E	117	GLU
2	E	130	ARG
2	E	140	LYS
2	E	142	ASN
2	E	146	THR
2	E	148	GLN
2	E	149	GLN
2	E	167	GLN
2	E	168	LEU

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Mol	Chain	Res	Type
2	E	172	GLU
2	E	173	ILE
2	E	174[A]	GLU
2	E	174[B]	GLU
2	E	212	LYS
2	E	216	LEU
2	E	223	LYS
2	E	238	GLU
2	E	240	LYS
2	E	241	GLN
2	E	266	GLU
2	E	281	LEU
2	E	296	MET
2	E	311	GLN
2	E	312	ILE
2	E	318	LEU
2	E	326	LEU
2	E	337	THR
2	E	344	LEU
2	E	352	THR
2	E	355	TYR
2	E	361	THR
2	E	366	ILE
2	E	375	ARG
2	E	378	GLU
2	E	385	GLU
2	E	386	SER
2	E	388	GLU
2	E	389	ASN
2	E	413	LEU
2	E	419	THR
2	E	423	ASP
2	E	432	LEU
2	E	438	LEU
2	F	1	HIS
2	F	11	SER
2	F	13	LEU
2	F	26	SER
2	F	27	VAL
2	F	31	LEU
2	F	37	ARG
2	F	41	ASN

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Mol	Chain	Res	Type
2	F	49	THR
2	F	59	THR
2	F	68	ARG
2	F	70	LEU
2	F	94	LYS
2	F	103	LEU
2	F	108	VAL
2	F	121	TYR
2	F	130	ARG
2	F	140	LYS
2	F	142	ASN
2	F	146	THR
2	F	148	GLN
2	F	149	GLN
2	F	167	GLN
2	F	168	LEU
2	F	172	GLU
2	F	173	ILE
2	F	174	GLU
2	F	212	LYS
2	F	216	LEU
2	F	223	LYS
2	F	237	GLU
2	F	239	LEU
2	F	266	GLU
2	F	281	LEU
2	F	291	SER
2	F	292	THR
2	F	296	MET
2	F	311	GLN
2	F	312	ILE
2	F	318	LEU
2	F	326	LEU
2	F	344	LEU
2	F	352	THR
2	F	355	TYR
2	F	361	THR
2	F	366	ILE
2	F	375	ARG
2	F	378	GLU
2	F	385	GLU
2	F	388	GLU

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Mol	Chain	Res	Type
2	F	389	ASN
2	F	404	GLU
2	F	413	LEU
2	F	419	THR
2	F	423	ASP
2	F	438	LEU
2	F	439	SER
2	F	442	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (49) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	11	HIS
1	A	30	ASN
1	A	109	ASN
1	A	114	GLN
1	A	139	ASN
1	A	164	HIS
1	B	11	HIS
1	B	30	ASN
1	B	109	ASN
1	B	114	GLN
1	C	9	ASN
1	C	24	ASN
1	C	30	ASN
1	C	109	ASN
1	C	114	GLN
1	C	117	ASN
1	D	9	ASN
1	D	11	HIS
1	D	30	ASN
1	D	114	GLN
2	E	22	ASN
2	E	33	ASN
2	E	41	ASN
2	E	75	ASN
2	E	114	GLN
2	E	119	ASN
2	E	157	GLN
2	E	241	GLN
2	E	311	GLN

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Mol	Chain	Res	Type
2	E	348	ASN
2	E	382	GLN
2	E	384	ASN
2	E	389	ASN
2	E	416	GLN
2	F	16	HIS
2	F	22	ASN
2	F	33	ASN
2	F	41	ASN
2	F	75	ASN
2	F	114	GLN
2	F	119	ASN
2	F	301	HIS
2	F	348	ASN
2	F	365	ASN
2	F	382	GLN
2	F	384	ASN
2	F	389	ASN
2	F	416	GLN

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](#)

2 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	$\# Z > 2$
3	ANP	E	500	-	33,33,33	6.00	23 (69%)	45,52,52	2.80	15 (33%)
3	ANP	F	501	-	33,33,33	5.17	27 (81%)	45,52,52	2.77	20 (44%)

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
3	ANP	E	500	-	-	8/18/38/38	0/3/3/3
3	ANP	F	501	-	-	9/18/38/38	0/3/3/3

All (50) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	500	ANP	PG-O1G	18.11	1.73	1.46
3	F	501	ANP	PG-O1G	14.34	1.67	1.46
3	E	500	ANP	C8-N9	13.98	1.61	1.37
3	F	501	ANP	C8-N9	10.81	1.55	1.37
3	E	500	ANP	C5-C4	9.86	1.56	1.39
3	F	501	ANP	C2-N1	9.25	1.50	1.33
3	E	500	ANP	PG-N3B	9.24	1.87	1.63
3	F	501	ANP	C5-C4	8.66	1.54	1.39
3	E	500	ANP	C2-N1	8.24	1.48	1.33
3	F	501	ANP	PG-O3G	7.88	1.77	1.56
3	E	500	ANP	C4-N9	-7.43	1.22	1.37
3	E	500	ANP	C1'-N9	6.75	1.65	1.46
3	E	500	ANP	PG-O3G	6.57	1.74	1.56
3	E	500	ANP	C2-N3	-6.19	1.22	1.33
3	F	501	ANP	C1'-N9	6.00	1.62	1.46
3	E	500	ANP	PB-O2B	-5.82	1.41	1.56
3	F	501	ANP	C8-N7	5.76	1.42	1.31
3	F	501	ANP	PB-O2B	-5.39	1.42	1.56
3	E	500	ANP	C8-N7	5.37	1.42	1.31
3	F	501	ANP	C6-N1	5.02	1.49	1.35
3	E	500	ANP	PB-O3A	-4.81	1.53	1.59
3	F	501	ANP	C4-N9	-4.68	1.28	1.37
3	F	501	ANP	PA-O3A	-4.60	1.54	1.59
3	E	500	ANP	C6-N1	4.48	1.48	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	500	ANP	PG-O2G	4.40	1.68	1.56
3	E	500	ANP	PB-O1B	-4.28	1.39	1.46
3	F	501	ANP	PB-N3B	-4.15	1.52	1.63
3	F	501	ANP	PG-O2G	4.13	1.67	1.56
3	F	501	ANP	C2-N3	-4.05	1.26	1.33
3	F	501	ANP	O3'-C3'	-4.00	1.33	1.43
3	E	500	ANP	PA-O1A	-3.76	1.37	1.50
3	E	500	ANP	C2'-C3'	-3.69	1.43	1.53
3	F	501	ANP	C4-N3	3.64	1.41	1.34
3	E	500	ANP	O3'-C3'	-3.45	1.34	1.43
3	F	501	ANP	C5-C6	-3.43	1.31	1.41
3	F	501	ANP	C2'-C3'	-3.39	1.44	1.53
3	F	501	ANP	PG-N3B	3.36	1.72	1.63
3	E	500	ANP	C4-N3	3.02	1.40	1.34
3	F	501	ANP	O4'-C1'	3.01	1.49	1.42
3	F	501	ANP	C6-N6	-2.99	1.26	1.34
3	E	500	ANP	O4'-C1'	2.95	1.48	1.42
3	F	501	ANP	PA-O5'	-2.66	1.49	1.59
3	E	500	ANP	PB-N3B	-2.59	1.56	1.63
3	E	500	ANP	PA-O3A	-2.53	1.56	1.59
3	F	501	ANP	PB-O1B	-2.49	1.42	1.46
3	F	501	ANP	PA-O1A	-2.43	1.42	1.50
3	F	501	ANP	C5-N7	-2.35	1.34	1.39
3	F	501	ANP	O5'-C5'	-2.30	1.35	1.44
3	F	501	ANP	PB-O3A	-2.21	1.56	1.59
3	E	500	ANP	C3'-C4'	2.19	1.58	1.53

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	500	ANP	C5-C4-N3	7.01	136.37	126.72
3	F	501	ANP	O1G-PG-N3B	-6.80	101.76	111.77
3	E	500	ANP	C4-C5-N7	6.44	117.95	110.58
3	F	501	ANP	C5-C4-N3	6.25	135.32	126.72
3	E	500	ANP	C6-C5-C4	-6.23	108.67	117.18
3	E	500	ANP	N3-C4-N9	-6.10	116.78	127.17
3	F	501	ANP	C4-C5-N7	5.88	117.30	110.58
3	E	500	ANP	C5-N7-C8	-5.80	94.33	103.45
3	F	501	ANP	N3-C4-N9	-5.36	118.05	127.17
3	E	500	ANP	O4'-C1'-N9	5.21	118.09	108.09
3	F	501	ANP	O4'-C1'-N9	4.52	116.77	108.09
3	F	501	ANP	C2'-C3'-C4'	4.41	111.12	102.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	500	ANP	O2B-PB-O1B	4.34	119.17	109.87
3	F	501	ANP	C6-C5-C4	-4.19	111.46	117.18
3	E	500	ANP	C2'-C3'-C4'	4.18	110.69	102.61
3	F	501	ANP	C5-N7-C8	-4.06	97.07	103.45
3	F	501	ANP	O2B-PB-O3A	3.83	117.43	104.64
3	F	501	ANP	O1B-PB-N3B	-3.79	106.19	111.77
3	E	500	ANP	C1'-N9-C8	3.78	135.49	127.09
3	F	501	ANP	C1'-N9-C8	3.76	135.44	127.09
3	E	500	ANP	O3'-C3'-C4'	-3.32	101.54	111.08
3	E	500	ANP	O2G-PG-O3G	3.28	116.39	107.59
3	F	501	ANP	O2G-PG-O3G	2.87	115.30	107.59
3	E	500	ANP	C4-N9-C1'	-2.73	120.25	126.63
3	F	501	ANP	PA-O5'-C5'	2.70	136.83	121.35
3	F	501	ANP	C5-C6-N6	-2.63	116.78	123.29
3	F	501	ANP	N6-C6-N1	2.45	123.83	118.38
3	F	501	ANP	C2-N3-C4	-2.40	105.98	111.83
3	F	501	ANP	C4-N9-C8	-2.34	103.28	105.74
3	F	501	ANP	C4-N9-C1'	-2.29	121.28	126.63
3	F	501	ANP	O3A-PB-N3B	-2.22	100.44	106.59
3	F	501	ANP	O5'-PA-O1A	2.19	117.61	108.94
3	E	500	ANP	PA-O5'-C5'	2.19	133.89	121.35
3	E	500	ANP	C2-N1-C6	2.18	122.31	118.73
3	E	500	ANP	O5'-PA-O1A	2.12	117.36	108.94

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	500	ANP	PB-N3B-PG-O1G
3	E	500	ANP	PG-N3B-PB-O1B
3	E	500	ANP	PG-N3B-PB-O3A
3	E	500	ANP	PA-O3A-PB-O1B
3	F	501	ANP	PB-N3B-PG-O1G
3	F	501	ANP	PG-N3B-PB-O1B
3	F	501	ANP	PG-N3B-PB-O3A
3	F	501	ANP	PA-O3A-PB-O2B
3	E	500	ANP	O4'-C1'-N9-C8
3	F	501	ANP	O4'-C1'-N9-C8
3	E	500	ANP	O4'-C4'-C5'-O5'
3	F	501	ANP	O4'-C4'-C5'-O5'
3	E	500	ANP	O4'-C1'-N9-C4
3	F	501	ANP	O4'-C1'-N9-C4

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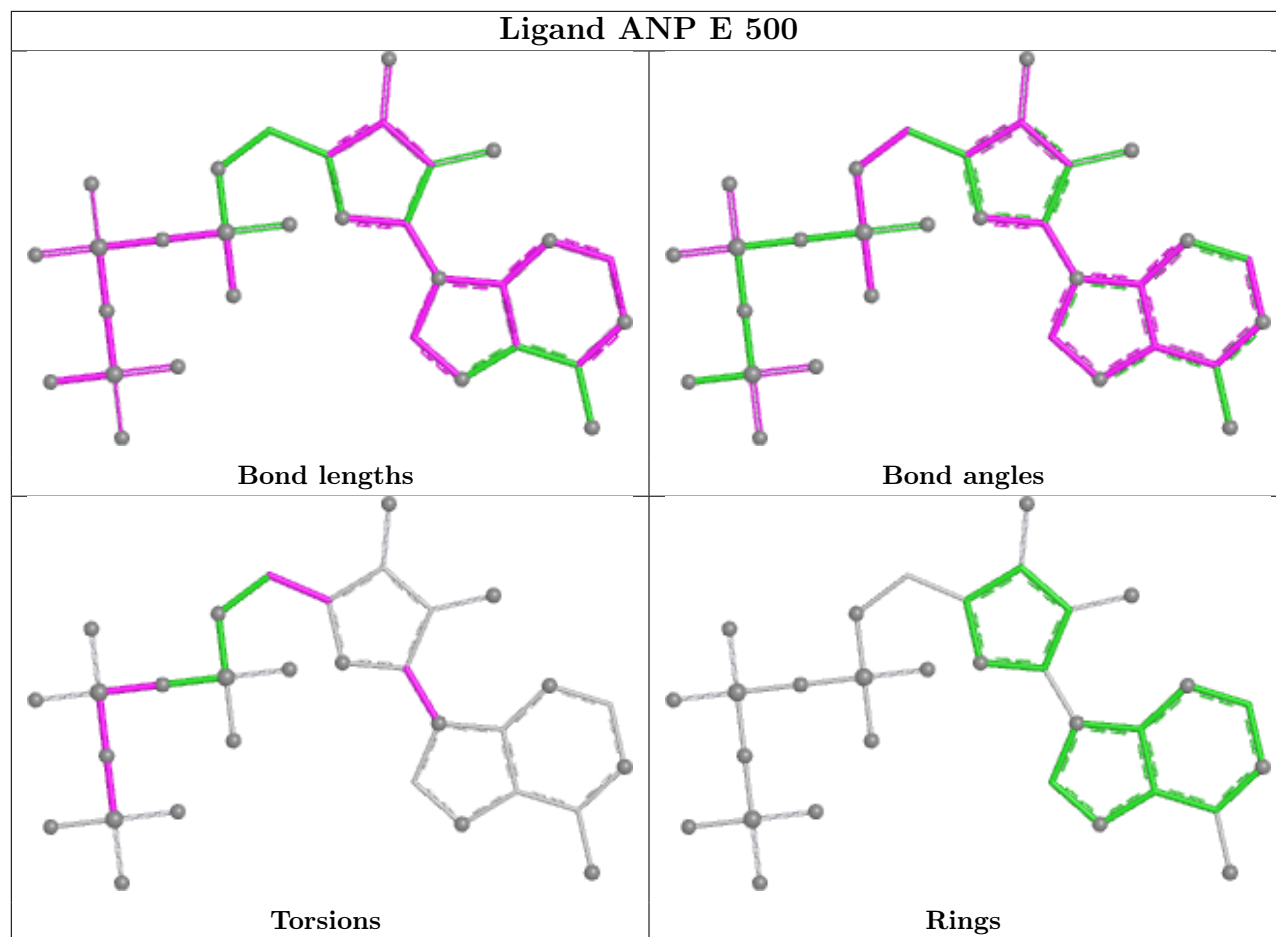
Mol	Chain	Res	Type	Atoms
3	E	500	ANP	C3'-C4'-C5'-O5'
3	F	501	ANP	C3'-C4'-C5'-O5'
3	F	501	ANP	PA-O3A-PB-O1B

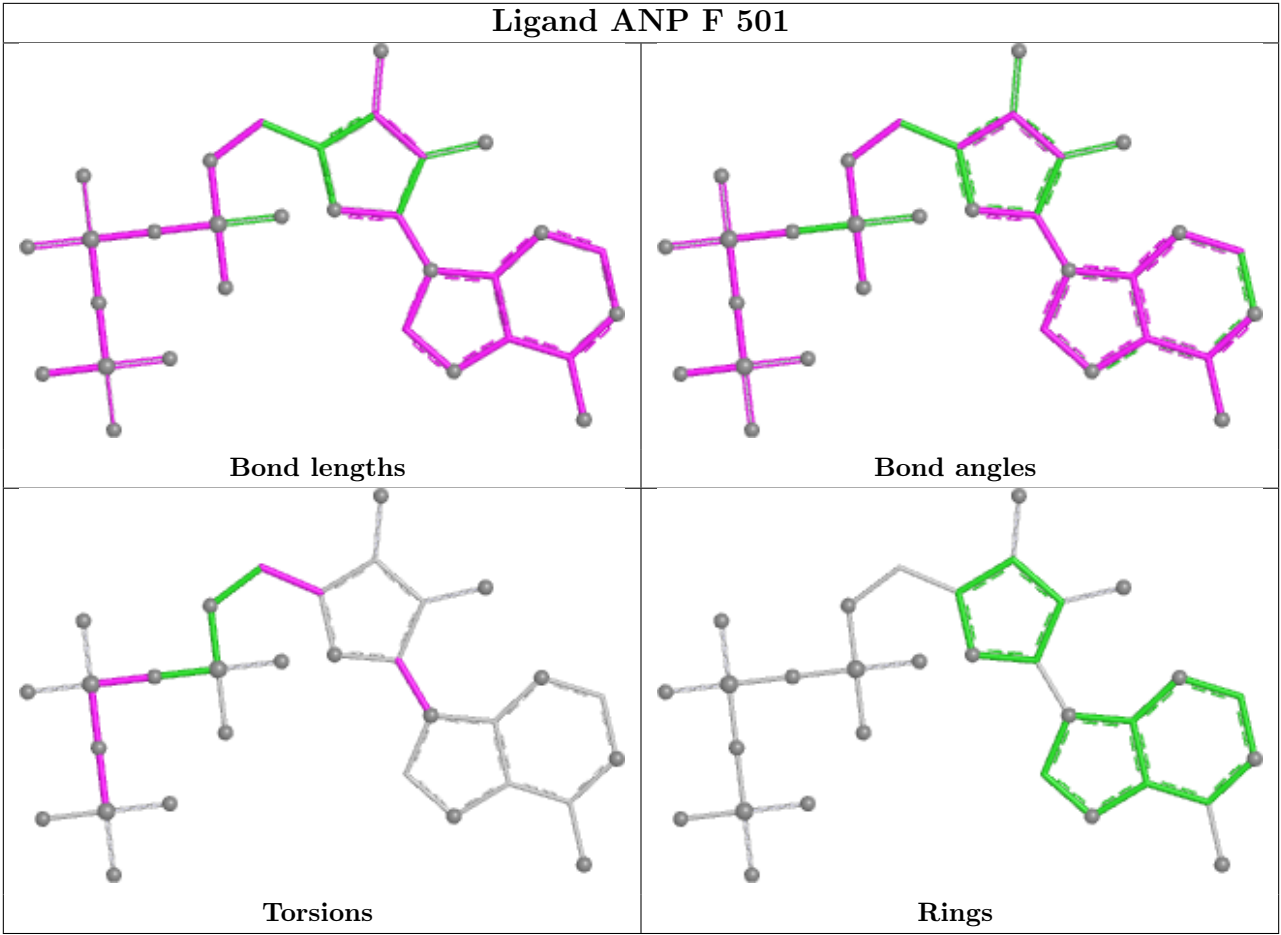
There are no ring outliers.

2 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	500	ANP	7	0
3	F	501	ANP	9	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 ⓘ

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	F	6
2	E	4

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	E	153:SER	C	154:ALA	N	2.67
1	F	153:SER	C	154:ALA	N	2.67
1	E	220:ASP	C	221:ALA	N	2.60
1	F	220:ASP	C	221:ALA	N	2.60

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Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	E	165:GLU	C	166:GLY	N	2.22
1	F	165:GLU	C	166:GLY	N	2.22
1	F	224:LEU	C	225:LEU	N	1.62
1	E	138:PRO	C	139:ALA	N	1.18
1	F	138:PRO	C	139:ALA	N	1.18
1	F	129:GLU	C	130:ARG	N	1.14

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	174/175 (99%)	0.16	5 (2%)	53	43	36, 63, 101, 108	4 (2%)
1	B	174/175 (99%)	0.19	3 (1%)	69	60	39, 65, 102, 108	4 (2%)
1	C	174/175 (99%)	0.48	4 (2%)	61	51	48, 77, 106, 113	3 (1%)
1	D	174/175 (99%)	0.58	11 (6%)	26	19	46, 80, 106, 114	4 (2%)
2	E	382/449 (85%)	0.23	22 (5%)	29	22	0, 60, 114, 124	6 (1%)
2	F	352/449 (78%)	0.23	27 (7%)	19	14	35, 58, 122, 132	4 (1%)
All	All	1430/1598 (89%)	0.29	72 (5%)	34	26	0, 65, 113, 132	25 (1%)

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	153	SER	9.1
2	F	89	VAL	8.1
2	E	174[A]	GLU	7.3
2	F	123	ALA	6.8
2	F	221	ALA	6.2
2	F	225	LEU	6.0
2	F	222	MET	5.4
2	E	152	PRO	5.2
2	E	91	TYR	4.6
2	E	89	VAL	4.5
2	F	91	TYR	4.4
2	E	270	PRO	4.3
2	E	90	GLY	4.3
1	D	53	ALA	4.2
2	F	90	GLY	4.1
2	F	119	ASN	4.1
2	F	226	ILE	4.0
2	F	270	PRO	4.0
1	A	110	GLY	3.9

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Mol	Chain	Res	Type	RSRZ
2	E	151	GLU	3.9
2	E	218	ILE	3.7
1	D	70	HIS	3.7
1	C	83	ARG	3.6
2	E	222	MET	3.4
1	A	83	ARG	3.3
2	F	224	LEU	3.2
1	D	83	ARG	3.2
2	F	126	LEU	3.1
2	E	219	LYS	3.1
2	F	92	VAL	3.1
1	A	69	GLY	3.1
2	F	267	SER	3.1
1	C	110	GLY	3.1
2	E	267	SER	3.0
2	E	139	ALA	3.0
2	F	223	LYS	3.0
2	E	119	ASN	2.9
2	E	92	VAL	2.9
2	E	138	PRO	2.8
2	E	225	LEU	2.6
2	F	88	GLU	2.6
1	B	83	ARG	2.6
1	D	44	ALA	2.6
1	A	54	PHE	2.5
2	E	264	ARG	2.5
2	F	264	ARG	2.5
2	F	266	GLU	2.5
2	F	230	ALA	2.5
2	F	231	ALA	2.5
1	D	76	VAL	2.4
2	F	240	LYS	2.4
1	B	69	GLY	2.3
2	E	266	GLU	2.3
1	D	67	HIS	2.3
1	A	68	GLN	2.2
2	E	236	PRO	2.2
2	E	118	LYS	2.2
1	D	86	ARG	2.2
1	C	118	ASP	2.2
2	F	122	ARG	2.1
2	F	247	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	72	VAL	2.1
2	F	118	LYS	2.1
1	D	55	THR	2.1
2	E	165	GLU	2.1
1	D	87	MET	2.0
2	F	111	VAL	2.0
1	B	46	PHE	2.0
1	D	62	ARG	2.0
2	F	115	ALA	2.0
2	F	121	TYR	2.0
1	D	38	TYR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

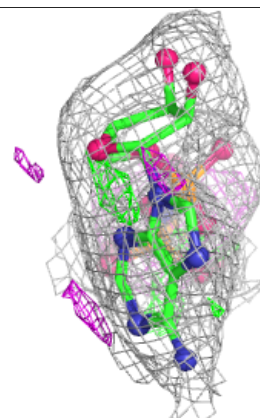
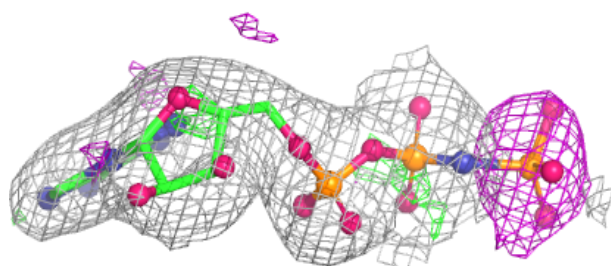
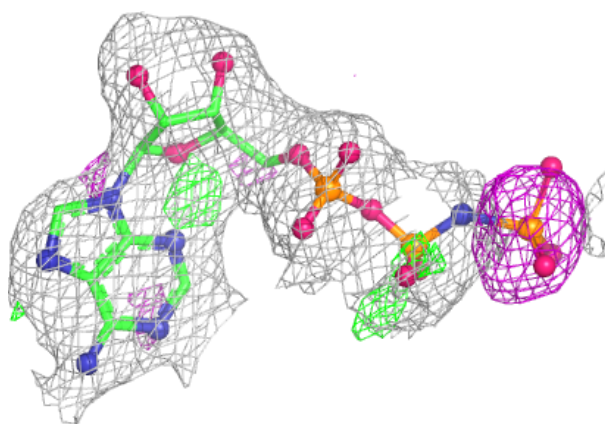
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	ANP	E	500	31/31	0.86	0.14	41,46,77,78	0
3	ANP	F	501	31/31	0.88	0.13	47,50,79,80	0

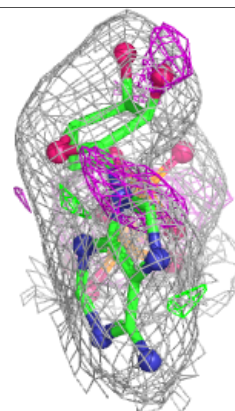
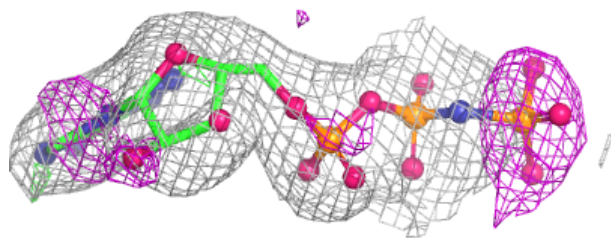
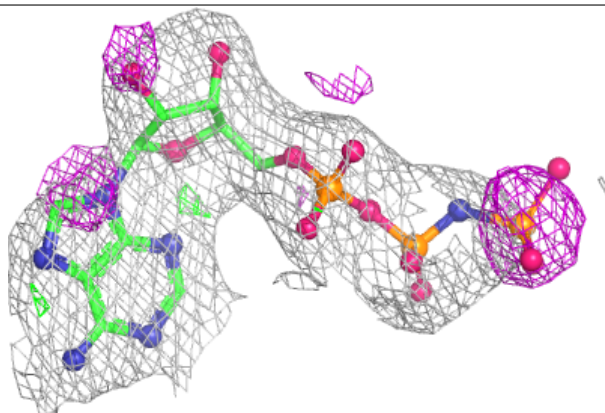
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ANP E 500:

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

**Electron density around ANP F 501:**

$2mF_o-DF_c$ (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.