



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:10 PM GMT

PDB ID : 1J21  
Title : Crystal Structure of Thermus thermophilus HB8 Argininosuccinate Synthetase  
in complex with ATP and citrulline  
Authors : Goto, M.; Hirotsu, K.; RIKEN Structural Genomics/Proteomics Initiative  
(RSGI)  
Deposited on : 2002-12-24  
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

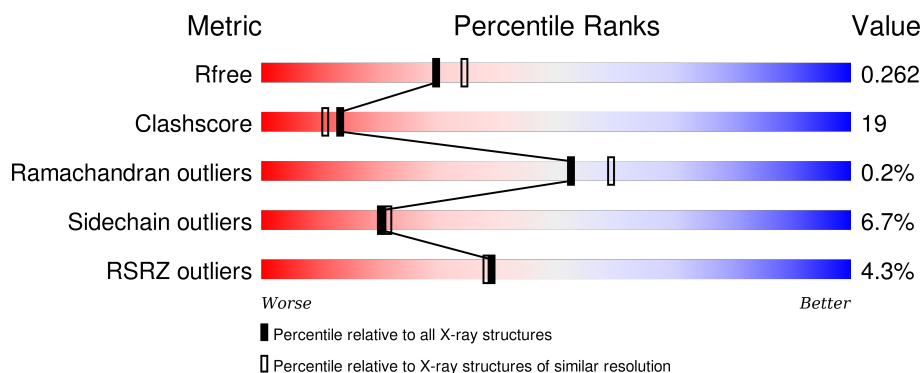
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	400	<div> <div>3%</div> <div>65%</div> <div>26%</div> <div>5%</div> </div>
1	B	400	<div> <div>2%</div> <div>68%</div> <div>24%</div> <div>5%</div> </div>
1	C	400	<div> <div>7%</div> <div>65%</div> <div>26%</div> <div>5%</div> </div>
1	D	400	<div> <div>5%</div> <div>63%</div> <div>29%</div> <div>5%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ATP	A	510	-	-	-	X
2	ATP	B	1510	-	-	-	X
2	ATP	D	3510	-	-	-	X
3	CIR	A	520	-	-	-	X
3	CIR	B	1520	-	-	-	X
3	CIR	C	2520	-	-	-	X
3	CIR	D	3520	-	-	-	X

## 2 Entry composition [i](#)

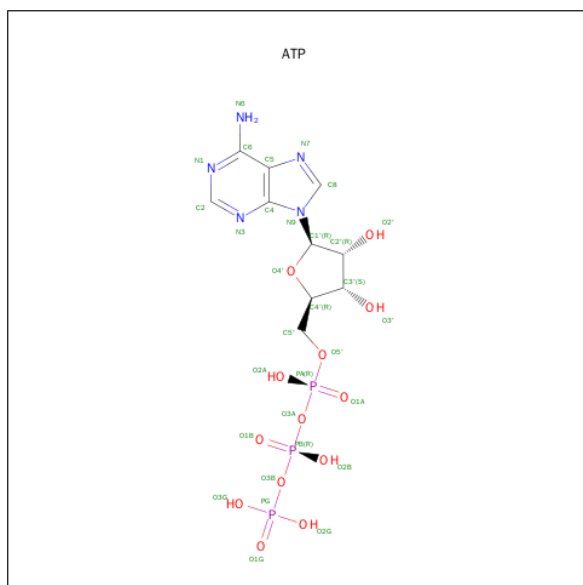
There are 4 unique types of molecules in this entry. The entry contains 12707 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Argininosuccinate Synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	380	Total	C	N	O	S	0	0	0
			3017	1931	526	551	9			
1	B	380	Total	C	N	O	S	0	0	0
			3017	1931	526	551	9			
1	C	380	Total	C	N	O	S	0	0	0
			3017	1931	526	551	9			
1	D	380	Total	C	N	O	S	0	0	0
			3017	1931	526	551	9			

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



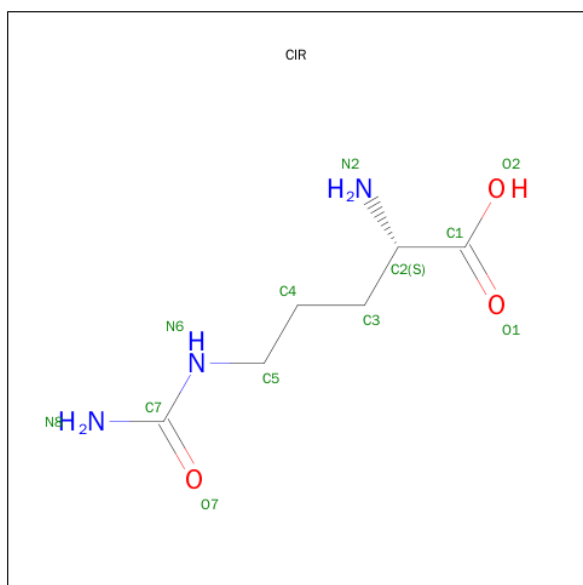
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

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

- Molecule 3 is CITRULLINE (three-letter code: CIR) (formula:  $C_6H_{13}N_3O_3$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			12	6	3	3		
3	B	1	Total	C	N	O	0	0
			12	6	3	3		
3	C	1	Total	C	N	O	0	0
			12	6	3	3		
3	D	1	Total	C	N	O	0	0
			12	6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	121	Total	O	0	0
			121	121		
4	B	134	Total	O	0	0
			134	134		
4	C	101	Total	O	0	0
			101	101		

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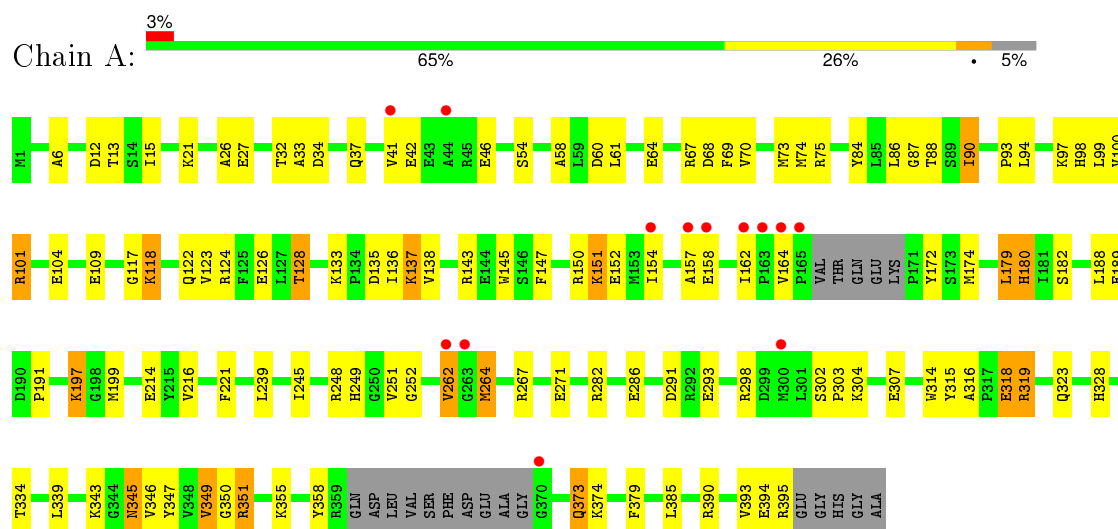
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	111	Total	O	0	0
			111	111		

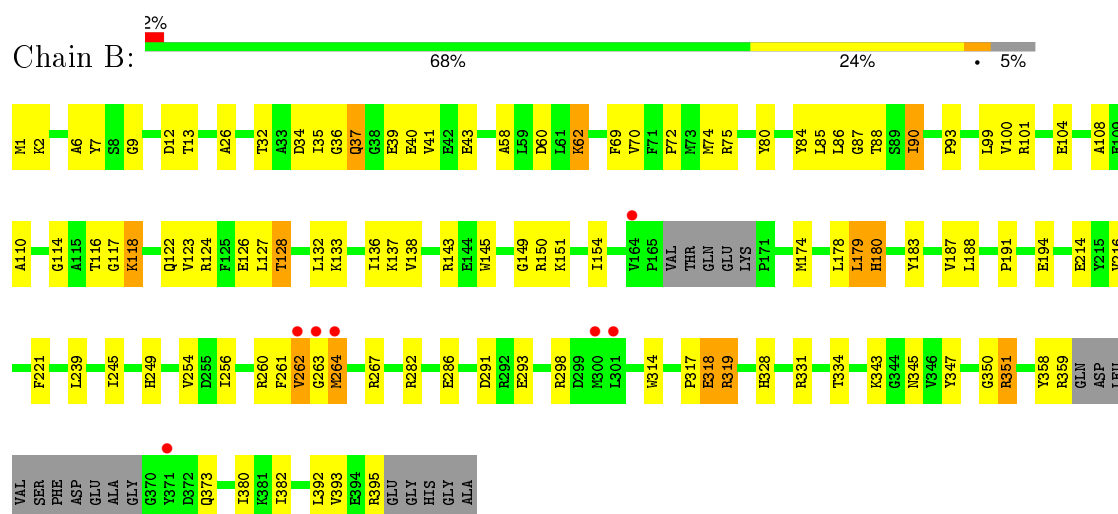
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Argininosuccinate Synthetase

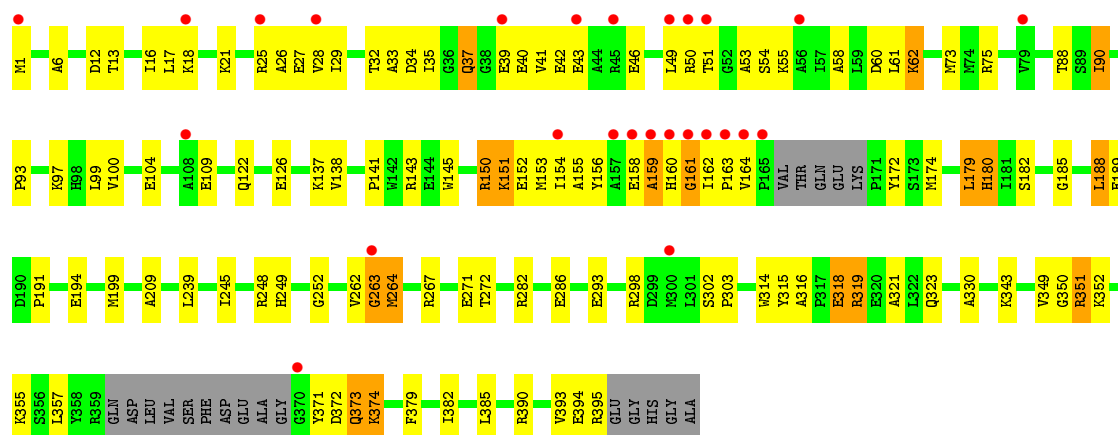


#### • Molecule 1: Argininosuccinate Synthetase

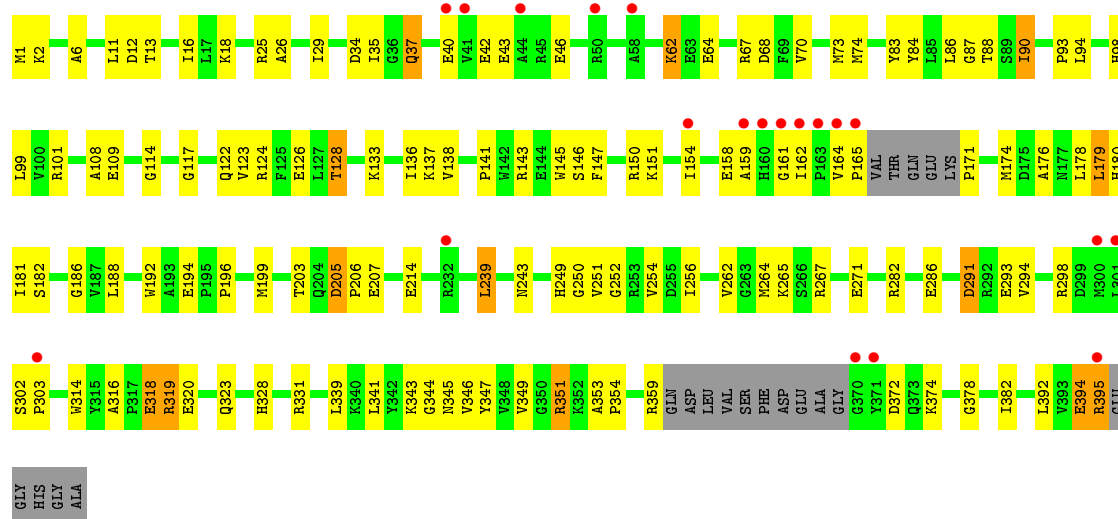


#### • Molecule 1: Argininosuccinate Synthetase





### • Molecule 1: Argininosuccinate Synthetase





## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	229.20Å 229.20Å 161.25Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.20 48.66 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.20) 99.4 (48.66-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.10 (at 2.20Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.230 , 0.259 0.233 , 0.262	Depositor DCC
$R_{free}$ test set	15719 reflections (9.96%)	DCC
Wilson B-factor (Å <sup>2</sup> )	35.2	Xtriage
Anisotropy	0.129	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 60.0	EDS
Estimated twinning fraction	0.001 for h,-h-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.53$ , $\langle L^2 \rangle = 0.37$	Xtriage
Outliers	0 of 159320 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	12707	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.92% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: CIR, 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	0.37	0/3085	0.62	0/4169
1	B	0.38	0/3085	0.63	0/4169
1	C	0.38	0/3085	0.64	1/4169 (0.0%)
1	D	0.38	0/3085	0.63	0/4169
All	All	0.38	0/12340	0.63	1/16676 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	263	GLY	N-CA-C	6.79	130.09	113.10

There are no chirality outliers.

There are no planarity outliers.

### 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	3017	0	3006	127	0
1	B	3017	0	3006	120	0
1	C	3017	0	3006	134	0
1	D	3017	0	3006	126	0
2	A	31	0	10	3	0
2	B	31	0	10	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	31	0	10	2	0
2	D	31	0	12	7	0
3	A	12	0	12	1	0
3	B	12	0	12	2	0
3	C	12	0	12	2	0
3	D	12	0	12	1	0
4	A	121	0	0	8	0
4	B	134	0	0	7	0
4	C	101	0	0	3	0
4	D	111	0	0	5	0
All	All	12707	0	12114	457	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 19.

All (457) 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:D:151:LYS:H	1:D:151:LYS:HD2	1.06	1.16
1:C:151:LYS:HD2	1:C:151:LYS:H	1.02	1.12
1:B:101:ARG:HH11	1:B:101:ARG:HB3	1.17	1.08
1:A:151:LYS:H	1:A:151:LYS:HD2	1.15	1.06
1:B:151:LYS:HD2	1:B:151:LYS:H	1.15	1.06
1:C:18:LYS:HG3	1:C:162:ILE:HD11	1.35	1.04
1:A:123:VAL:HG22	1:B:380:ILE:HD11	1.43	0.98
1:C:109:GLU:O	1:C:137:LYS:HE2	1.62	0.98
1:B:101:ARG:HB3	1:B:101:ARG:NH1	1.81	0.94
1:B:35:ILE:H	1:B:37:GLN:NE2	1.66	0.94
1:C:35:ILE:H	1:C:37:GLN:NE2	1.66	0.93
1:A:124:ARG:O	1:A:128:THR:HG22	1.68	0.93
1:C:151:LYS:CD	1:C:151:LYS:H	1.82	0.91
1:D:35:ILE:H	1:D:37:GLN:NE2	1.69	0.90
1:D:151:LYS:H	1:D:151:LYS:CD	1.84	0.89
1:D:351:ARG:HD2	1:D:351:ARG:N	1.88	0.88
1:C:151:LYS:HD2	1:C:151:LYS:N	1.88	0.88
1:A:293:GLU:CB	1:C:318:GLU:HG2	2.05	0.86
1:D:124:ARG:O	1:D:128:THR:HG23	1.75	0.85
1:A:151:LYS:N	1:A:151:LYS:HD2	1.92	0.84
1:B:151:LYS:CD	1:B:151:LYS:H	1.91	0.82
1:D:150:ARG:O	1:D:154:ILE:HD13	1.80	0.82
1:B:293:GLU:CB	1:D:318:GLU:HG2	2.10	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:VAL:HB	1:A:264:MET:HE1	1.62	0.81
1:A:351:ARG:N	1:A:351:ARG:HD2	1.95	0.81
1:B:351:ARG:N	1:B:351:ARG:HD2	1.97	0.80
1:B:151:LYS:HD2	1:B:151:LYS:N	1.97	0.80
1:B:293:GLU:HB2	1:D:318:GLU:HG2	1.63	0.80
1:B:41:VAL:HG12	1:B:58:ALA:HB1	1.63	0.80
1:B:124:ARG:O	1:B:128:THR:HG23	1.82	0.80
1:A:373:GLN:H	1:A:373:GLN:NE2	1.80	0.79
1:D:151:LYS:N	1:D:151:LYS:HD2	1.92	0.79
1:C:122:GLN:NE2	1:C:143:ARG:HH11	1.82	0.77
1:A:293:GLU:HB2	1:C:318:GLU:HG2	1.64	0.77
1:B:101:ARG:HH11	1:B:101:ARG:CB	1.97	0.77
1:A:293:GLU:HB2	1:C:318:GLU:CG	2.14	0.76
1:C:351:ARG:N	1:C:351:ARG:HD2	2.01	0.76
1:B:34:ASP:HB2	1:B:41:VAL:HG21	1.66	0.76
1:B:100:VAL:O	1:B:104:GLU:HG3	1.85	0.76
1:A:151:LYS:CD	1:A:151:LYS:H	1.92	0.76
1:C:373:GLN:HE21	1:C:373:GLN:H	1.33	0.76
1:D:319:ARG:HH21	1:D:323:GLN:HE21	1.34	0.75
1:B:40:GLU:HG3	1:B:43:GLU:HG3	1.70	0.74
1:D:40:GLU:OE1	1:D:42:GLU:HB2	1.87	0.73
1:A:374:LYS:NZ	1:D:374:LYS:HG2	2.03	0.73
1:A:262:VAL:HG23	1:A:264:MET:SD	2.27	0.73
1:A:318:GLU:HG2	1:C:293:GLU:CB	2.19	0.72
1:A:293:GLU:HB3	1:C:318:GLU:HG2	1.71	0.72
1:C:109:GLU:HG3	1:C:137:LYS:HE3	1.71	0.72
1:C:1:MET:O	1:C:26:ALA:HB1	1.89	0.71
1:C:122:GLN:HE21	1:C:143:ARG:HH11	1.35	0.71
1:B:293:GLU:HB2	1:D:318:GLU:CG	2.20	0.71
1:C:319:ARG:HH21	1:C:323:GLN:HE21	1.36	0.71
1:C:319:ARG:HH21	1:C:323:GLN:NE2	1.89	0.71
1:A:264:MET:N	1:A:264:MET:SD	2.64	0.71
1:C:373:GLN:NE2	1:C:373:GLN:H	1.88	0.70
1:B:13:THR:HG23	2:B:1510:ATP:O3G	1.92	0.70
1:B:93:PRO:HG3	4:B:1346:HOH:O	1.92	0.69
1:A:41:VAL:HG12	1:A:58:ALA:HB1	1.73	0.69
1:A:93:PRO:HG3	4:A:1220:HOH:O	1.92	0.69
1:C:37:GLN:H	1:C:37:GLN:NE2	1.90	0.69
1:A:100:VAL:O	1:A:104:GLU:HG3	1.93	0.69
1:D:1:MET:O	1:D:26:ALA:HB1	1.94	0.68
1:B:314:TRP:CE3	1:B:319:ARG:HD2	2.29	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:GLU:CG	1:C:293:GLU:HB2	2.24	0.67
1:B:35:ILE:N	1:B:37:GLN:NE2	2.41	0.67
1:A:150:ARG:HG3	1:A:154:ILE:HD13	1.77	0.66
1:B:314:TRP:O	1:B:319:ARG:HD3	1.95	0.66
1:A:318:GLU:HG2	1:C:293:GLU:HB2	1.78	0.66
1:B:37:GLN:H	1:B:37:GLN:NE2	1.94	0.66
1:B:1:MET:O	1:B:26:ALA:HB1	1.96	0.66
1:A:27:GLU:HG3	1:A:54:SER:CB	2.27	0.65
1:D:83:TYR:HB3	1:D:254:VAL:HG21	1.78	0.65
1:A:27:GLU:HG3	1:A:54:SER:HB3	1.77	0.65
1:C:93:PRO:HG3	4:C:1461:HOH:O	1.96	0.65
1:D:6:ALA:HB2	1:D:99:LEU:HD11	1.77	0.64
1:A:318:GLU:HG2	1:C:293:GLU:HB3	1.80	0.64
1:D:40:GLU:HG3	1:D:43:GLU:HG3	1.79	0.64
1:C:245:ILE:O	1:C:248:ARG:HG2	1.98	0.64
1:D:42:GLU:O	1:D:46:GLU:HG3	1.98	0.63
1:A:143:ARG:HG2	1:B:373:GLN:HB3	1.81	0.63
1:D:37:GLN:NE2	1:D:37:GLN:H	1.97	0.63
1:D:205:ASP:OD1	1:D:207:GLU:OE2	2.17	0.63
1:B:41:VAL:CG1	1:B:58:ALA:HB1	2.29	0.63
1:B:293:GLU:HB3	1:D:318:GLU:HG2	1.80	0.63
1:D:93:PRO:HG3	4:D:1266:HOH:O	1.97	0.63
1:D:328:HIS:HA	1:D:331:ARG:HH12	1.63	0.63
1:D:341:LEU:HD22	1:D:346:VAL:HG22	1.80	0.63
1:A:13:THR:HG23	2:A:510:ATP:O3G	1.99	0.62
1:C:13:THR:HG23	2:C:2510:ATP:O3G	1.98	0.62
1:D:35:ILE:H	1:D:37:GLN:HE22	1.45	0.62
1:B:34:ASP:HA	1:B:37:GLN:HE22	1.65	0.62
1:A:395:ARG:HG3	1:A:395:ARG:HH11	1.65	0.62
1:A:191:PRO:HG2	4:B:1371:HOH:O	1.98	0.61
1:A:374:LYS:HZ1	1:D:374:LYS:HG2	1.66	0.61
1:A:349:VAL:HG13	4:A:1112:HOH:O	2.00	0.61
1:A:385:LEU:HD23	1:C:263:GLY:HA3	1.82	0.61
1:D:162:ILE:O	1:D:164:VAL:HG13	2.01	0.61
1:A:88:THR:HG21	3:A:520:CIR:H42	1.83	0.61
1:B:282:ARG:O	1:B:286:GLU:HG3	2.01	0.60
1:D:124:ARG:O	1:D:128:THR:CG2	2.49	0.60
1:D:319:ARG:HH21	1:D:323:GLN:NE2	1.98	0.60
1:C:6:ALA:HB1	2:C:2510:ATP:N3	2.16	0.60
1:A:282:ARG:HD2	4:A:1293:HOH:O	2.01	0.60
1:C:25:ARG:NH1	1:C:25:ARG:HB3	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:50:ARG:O	1:C:50:ARG:HG2	2.01	0.60
1:A:75:ARG:NH2	4:A:1227:HOH:O	2.35	0.60
1:C:150:ARG:O	1:C:154:ILE:HG13	2.01	0.60
1:D:70:VAL:O	1:D:74:MET:HG3	2.01	0.59
1:A:104:GLU:HG2	1:A:133:LYS:NZ	2.17	0.59
1:C:32:THR:CG2	1:C:41:VAL:HG13	2.33	0.59
1:A:41:VAL:CG1	1:A:58:ALA:HB1	2.32	0.59
1:A:154:ILE:O	1:A:158:GLU:HG3	2.03	0.59
1:B:318:GLU:CG	1:D:293:GLU:HB2	2.33	0.59
1:C:35:ILE:N	1:C:37:GLN:NE2	2.45	0.59
1:C:154:ILE:O	1:C:158:GLU:HG3	2.03	0.59
1:D:29:ILE:HD12	1:D:29:ILE:N	2.18	0.59
1:A:104:GLU:HG2	1:A:133:LYS:HZ2	1.67	0.58
1:D:62:LYS:HD3	1:D:62:LYS:H	1.67	0.58
1:A:133:LYS:HG2	1:A:136:ILE:HB	1.86	0.58
1:A:262:VAL:C	1:A:264:MET:HE3	2.24	0.58
1:C:75:ARG:HG2	1:C:330:ALA:CB	2.34	0.58
1:A:150:ARG:HG3	1:A:154:ILE:CD1	2.33	0.58
1:D:6:ALA:HB3	2:D:3510:ATP:O2'	2.02	0.58
1:D:126:GLU:HG2	1:D:138:VAL:HG11	1.84	0.58
1:A:390:ARG:O	1:A:394:GLU:HG2	2.03	0.58
1:B:118:LYS:H	1:B:118:LYS:HD2	1.67	0.58
1:C:33:ALA:HB1	1:C:61:LEU:HD12	1.85	0.58
1:A:302:SER:HB2	1:A:303:PRO:HD3	1.86	0.57
1:C:35:ILE:H	1:C:37:GLN:HE22	1.50	0.57
1:A:373:GLN:H	1:A:373:GLN:CD	2.06	0.57
1:B:262:VAL:HG23	1:B:264:MET:SD	2.45	0.57
1:C:374:LYS:HE2	4:C:1306:HOH:O	2.04	0.57
1:C:262:VAL:O	1:C:262:VAL:HG12	2.05	0.57
1:A:345:ASN:HB3	1:A:347:TYR:CE1	2.39	0.57
1:D:13:THR:HG23	2:D:3510:ATP:O3G	2.02	0.57
1:D:345:ASN:HB3	1:D:347:TYR:CE1	2.40	0.57
1:B:117:GLY:HA2	1:B:122:GLN:NE2	2.19	0.57
1:C:156:TYR:O	1:C:159:ALA:HB3	2.04	0.57
1:A:21:LYS:HD2	1:A:26:ALA:O	2.05	0.56
1:C:122:GLN:HE21	1:C:143:ARG:NH1	2.01	0.56
1:B:260:ARG:HD2	1:B:264:MET:SD	2.44	0.56
1:A:157:ALA:HB1	1:A:162:ILE:HB	1.86	0.56
1:A:42:GLU:O	1:A:46:GLU:HG3	2.06	0.56
1:A:98:HIS:HD2	1:A:101:ARG:NH1	2.03	0.56
1:B:93:PRO:HA	1:B:128:THR:HG21	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:318:GLU:HG2	1:D:293:GLU:HB3	1.88	0.56
1:A:343:LYS:HD2	1:B:343:LYS:HD3	1.88	0.56
1:C:21:LYS:HD2	1:C:26:ALA:O	2.06	0.56
1:A:314:TRP:O	1:A:319:ARG:HD3	2.06	0.56
1:A:262:VAL:CB	1:A:264:MET:HE1	2.34	0.56
1:A:262:VAL:O	1:A:264:MET:HE3	2.06	0.56
1:B:395:ARG:HH11	1:B:395:ARG:HG3	1.70	0.56
1:B:249:HIS:HD2	4:B:1058:HOH:O	1.88	0.55
1:A:214:GLU:OE1	1:A:249:HIS:HE1	1.88	0.55
1:C:179:LEU:O	1:C:272:THR:HG23	2.06	0.55
1:C:314:TRP:O	1:C:319:ARG:HD3	2.07	0.55
1:A:349:VAL:HG22	1:B:194:GLU:HG3	1.88	0.55
1:A:262:VAL:O	1:C:382:ILE:HG23	2.06	0.55
1:D:109:GLU:O	1:D:137:LYS:HE2	2.06	0.55
1:B:122:GLN:HE21	1:B:143:ARG:HH11	1.55	0.55
1:C:395:ARG:HG3	1:C:395:ARG:HH11	1.71	0.55
1:A:21:LYS:HD3	4:A:1393:HOH:O	2.06	0.55
1:C:34:ASP:HA	1:C:37:GLN:HE22	1.72	0.54
1:B:12:ASP:HB2	2:B:1510:ATP:O1G	2.07	0.54
1:C:264:MET:CE	1:C:264:MET:H	2.20	0.54
1:C:189:GLU:O	1:C:191:PRO:HD3	2.07	0.54
1:D:64:GLU:OE1	1:D:98:HIS:HE1	1.90	0.54
1:C:12:ASP:OD1	1:C:150:ARG:NH2	2.40	0.54
1:D:122:GLN:HE21	1:D:143:ARG:HH11	1.54	0.54
1:D:164:VAL:HB	1:D:165:PRO:HD2	1.89	0.54
1:B:214:GLU:OE1	1:B:249:HIS:HE1	1.90	0.54
1:B:117:GLY:HA2	1:B:122:GLN:HE22	1.72	0.54
1:C:282:ARG:HD2	4:C:1310:HOH:O	2.08	0.54
1:B:216:VAL:HG21	1:B:245:ILE:HG21	1.90	0.54
1:B:318:GLU:HG2	1:D:293:GLU:CB	2.38	0.54
1:B:126:GLU:HG2	1:B:138:VAL:HG11	1.89	0.54
1:D:214:GLU:OE1	1:D:249:HIS:HE1	1.90	0.54
1:A:221:PHE:O	1:A:334:THR:OG1	2.19	0.54
1:B:393:VAL:C	1:B:395:ARG:H	2.12	0.53
1:B:1:MET:HB3	1:B:26:ALA:HB2	1.91	0.53
1:A:339:LEU:N	1:A:339:LEU:HD12	2.24	0.53
1:B:104:GLU:HG2	1:B:133:LYS:NZ	2.23	0.53
1:C:88:THR:HG21	3:C:2520:CIR:H42	1.90	0.53
1:A:191:PRO:HD3	1:B:358:TYR:CE2	2.44	0.53
1:D:314:TRP:O	1:D:319:ARG:HD3	2.07	0.53
1:B:124:ARG:O	1:B:128:THR:CG2	2.54	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:262:VAL:O	1:C:264:MET:HE3	2.09	0.53
1:D:316:ALA:HB1	1:D:318:GLU:OE2	2.08	0.53
1:B:133:LYS:O	1:B:136:ILE:HG22	2.08	0.53
1:B:395:ARG:HH22	1:C:355:LYS:NZ	2.06	0.53
1:A:358:TYR:HE2	1:B:191:PRO:HD3	1.73	0.53
1:B:75:ARG:NH2	4:B:1219:HOH:O	2.41	0.53
1:B:318:GLU:HG3	1:D:293:GLU:HB2	1.90	0.52
1:B:118:LYS:O	1:B:264:MET:HE1	2.09	0.52
1:C:37:GLN:HB2	1:C:39:GLU:HG2	1.90	0.52
1:C:126:GLU:HG2	1:C:138:VAL:HG11	1.91	0.52
1:D:98:HIS:HD2	1:D:101:ARG:NH1	2.08	0.52
1:B:87:GLY:HA3	1:B:180:HIS:CD2	2.44	0.52
1:B:393:VAL:O	1:B:395:ARG:NH1	2.42	0.52
1:C:372:ASP:C	1:C:372:ASP:OD2	2.49	0.52
1:B:60:ASP:OD1	1:B:62:LYS:HD3	2.09	0.52
1:A:264:MET:HG3	1:C:385:LEU:HD21	1.90	0.52
1:C:12:ASP:OD2	1:C:150:ARG:NH2	2.42	0.52
1:C:21:LYS:O	1:C:25:ARG:HA	2.10	0.51
1:B:87:GLY:HA2	1:B:179:LEU:HD13	1.92	0.51
1:D:87:GLY:HA2	1:D:179:LEU:HD13	1.92	0.51
1:B:37:GLN:HG2	4:B:1334:HOH:O	2.09	0.51
1:B:351:ARG:N	1:B:351:ARG:CD	2.72	0.51
1:C:209:ALA:O	1:C:343:LYS:NZ	2.43	0.51
1:A:216:VAL:HG22	4:A:1194:HOH:O	2.10	0.51
1:D:339:LEU:N	1:D:339:LEU:HD12	2.25	0.51
1:C:12:ASP:O	1:C:16:ILE:HG13	2.10	0.51
1:D:171:PRO:O	1:D:186:GLY:HA3	2.11	0.51
1:C:25:ARG:CB	1:C:25:ARG:HH11	2.24	0.51
1:C:6:ALA:HB2	1:C:99:LEU:HD11	1.92	0.51
1:D:133:LYS:O	1:D:136:ILE:HG22	2.11	0.51
1:C:100:VAL:O	1:C:104:GLU:HG3	2.11	0.51
1:A:319:ARG:HH21	1:A:323:GLN:NE2	2.08	0.51
1:C:40:GLU:CG	1:C:43:GLU:HG3	2.41	0.51
1:D:351:ARG:CD	1:D:351:ARG:N	2.69	0.51
1:B:116:THR:OG1	1:B:118:LYS:HD2	2.11	0.51
1:D:83:TYR:CB	1:D:254:VAL:HG21	2.40	0.50
1:B:180:HIS:ND1	1:B:180:HIS:C	2.63	0.50
1:A:172:TYR:CZ	1:A:199:MET:HG3	2.46	0.50
1:B:37:GLN:H	1:B:37:GLN:HE21	1.60	0.50
1:C:150:ARG:HB2	1:C:150:ARG:HH11	1.76	0.50
1:A:97:LYS:HD2	1:A:315:TYR:OH	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:137:LYS:N	1:C:137:LYS:HD3	2.27	0.50
1:B:393:VAL:C	1:B:395:ARG:N	2.63	0.50
1:C:27:GLU:HG3	1:C:54:SER:OG	2.11	0.50
1:A:109:GLU:O	1:A:137:LYS:HE2	2.11	0.50
1:B:80:TYR:HB3	1:B:84:TYR:HB3	1.93	0.50
1:B:40:GLU:CG	1:B:43:GLU:HG3	2.40	0.50
1:C:25:ARG:HH11	1:C:25:ARG:HB3	1.77	0.50
1:D:137:LYS:HD3	1:D:137:LYS:N	2.27	0.50
1:D:16:ILE:HG23	1:D:141:PRO:HG2	1.94	0.50
1:C:27:GLU:HG3	1:C:54:SER:CB	2.41	0.50
1:D:18:LYS:HG3	1:D:162:ILE:HD11	1.93	0.50
1:D:394:GLU:O	1:D:395:ARG:C	2.49	0.50
1:A:358:TYR:CE2	1:B:191:PRO:HD3	2.46	0.49
1:C:42:GLU:O	1:C:46:GLU:HG3	2.12	0.49
1:A:245:ILE:O	1:A:248:ARG:HG2	2.11	0.49
1:A:319:ARG:HH21	1:A:323:GLN:HE21	1.59	0.49
1:D:67:ARG:HD2	1:D:68:ASP:OD1	2.13	0.49
1:D:114:GLY:N	2:D:3510:ATP:O3'	2.46	0.49
1:A:249:HIS:HD2	4:A:1302:HOH:O	1.94	0.49
1:C:393:VAL:C	1:C:395:ARG:H	2.13	0.49
1:C:393:VAL:C	1:C:395:ARG:N	2.65	0.49
1:A:298:ARG:C	1:A:298:ARG:HD2	2.33	0.49
1:A:73:MET:CE	1:A:90:ILE:HG12	2.42	0.49
1:C:37:GLN:H	1:C:37:GLN:HE21	1.57	0.49
1:C:51:THR:HG22	1:C:51:THR:O	2.12	0.49
1:B:187:VAL:HG12	4:B:1373:HOH:O	2.13	0.49
1:B:350:GLY:C	1:B:351:ARG:HD2	2.33	0.48
1:A:355:LYS:HE2	1:D:394:GLU:OE2	2.13	0.48
1:A:61:LEU:HD13	1:A:94:LEU:HD11	1.96	0.48
1:C:151:LYS:CD	1:C:151:LYS:N	2.61	0.48
1:C:282:ARG:O	1:C:286:GLU:HG3	2.13	0.48
1:D:316:ALA:O	1:D:320:GLU:HG3	2.14	0.48
1:C:248:ARG:HG3	1:C:249:HIS:CD2	2.48	0.48
1:A:147:PHE:HD2	1:A:152:GLU:HG2	1.78	0.48
1:D:203:THR:HB	1:D:250:GLY:HA2	1.94	0.48
1:A:98:HIS:CD2	1:A:101:ARG:NH1	2.81	0.48
1:D:158:GLU:O	1:D:161:GLY:N	2.41	0.48
1:D:328:HIS:HA	1:D:331:ARG:NH1	2.27	0.48
1:C:390:ARG:O	1:C:394:GLU:HG2	2.14	0.48
1:B:32:THR:CG2	1:B:41:VAL:HG13	2.44	0.48
1:C:12:ASP:CG	1:C:150:ARG:HH21	2.17	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:302:SER:HB2	1:D:303:PRO:HD3	1.94	0.48
1:B:150:ARG:O	1:B:154:ILE:HD13	2.14	0.47
1:A:123:VAL:HG22	1:B:380:ILE:CD1	2.29	0.47
1:A:6:ALA:HB1	2:A:510:ATP:N3	2.29	0.47
4:A:1410:HOH:O	1:B:191:PRO:HG2	2.14	0.47
1:D:176:ALA:CB	1:D:181:ILE:HG12	2.43	0.47
1:B:74:MET:HB3	1:B:282:ARG:CZ	2.44	0.47
1:B:328:HIS:HA	1:B:331:ARG:HH12	1.79	0.47
1:B:88:THR:HG21	3:B:1520:CIR:H42	1.96	0.47
1:C:1:MET:CG	1:C:26:ALA:HB2	2.45	0.47
1:A:154:ILE:HG13	1:A:164:VAL:HG11	1.97	0.47
1:D:154:ILE:N	1:D:154:ILE:HD12	2.30	0.47
1:C:350:GLY:C	1:C:351:ARG:HD2	2.35	0.47
1:A:87:GLY:HA2	1:A:179:LEU:HD13	1.96	0.47
1:B:123:VAL:HG12	1:B:127:LEU:HD12	1.97	0.47
1:C:162:ILE:O	1:C:164:VAL:HG13	2.15	0.47
1:D:314:TRP:CE3	1:D:319:ARG:HD2	2.49	0.47
1:C:252:GLY:O	1:C:271:GLU:HA	2.15	0.47
1:A:350:GLY:C	1:A:351:ARG:HD2	2.35	0.47
1:D:250:GLY:O	1:D:344:GLY:HA2	2.14	0.47
1:C:21:LYS:CD	1:C:28:VAL:HG23	2.45	0.47
1:C:16:ILE:HG23	1:C:141:PRO:HG2	1.96	0.47
1:A:126:GLU:HG2	1:A:138:VAL:HG11	1.97	0.47
1:C:319:ARG:NH2	1:C:323:GLN:HE21	2.10	0.46
1:A:385:LEU:HD23	1:C:263:GLY:CA	2.44	0.46
1:C:262:VAL:O	1:C:264:MET:CE	2.63	0.46
1:D:2:LYS:HB2	1:D:108:ALA:HA	1.96	0.46
1:D:164:VAL:HB	1:D:165:PRO:CD	2.44	0.46
1:D:122:GLN:NE2	1:D:143:ARG:HH11	2.13	0.46
1:D:151:LYS:N	1:D:151:LYS:CD	2.63	0.46
1:B:318:GLU:CG	1:D:293:GLU:CB	2.93	0.46
1:C:1:MET:HG3	1:C:26:ALA:CB	2.46	0.46
1:C:40:GLU:HG3	1:C:43:GLU:HG3	1.96	0.46
1:C:152:GLU:O	1:C:155:ALA:HB3	2.16	0.46
1:D:262:VAL:O	1:D:262:VAL:CG1	2.64	0.46
1:C:49:LEU:HA	1:C:53:ALA:HB3	1.98	0.46
1:B:110:ALA:HB2	1:B:137:LYS:HB3	1.97	0.46
1:C:109:GLU:CG	1:C:137:LYS:HE3	2.44	0.46
1:B:395:ARG:NH2	1:C:355:LYS:NZ	2.64	0.46
1:C:371:TYR:OH	1:D:262:VAL:HG11	2.15	0.46
1:A:67:ARG:HG2	1:A:68:ASP:OD1	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:12:ASP:HB2	2:D:3510:ATP:O3G	2.16	0.46
1:A:282:ARG:O	1:A:286:GLU:HG3	2.16	0.46
1:A:137:LYS:N	1:A:137:LYS:HD3	2.31	0.46
1:C:17:LEU:HD23	1:C:17:LEU:O	2.16	0.46
1:B:317:PRO:HB2	1:D:294:VAL:CG2	2.45	0.46
1:D:154:ILE:N	1:D:154:ILE:CD1	2.79	0.46
1:D:239:LEU:HD22	1:D:243:ASN:HD21	1.79	0.46
1:C:180:HIS:ND1	1:C:180:HIS:C	2.69	0.46
1:A:293:GLU:HB2	1:C:318:GLU:HG3	1.95	0.46
1:D:174:MET:HA	1:D:182:SER:O	2.16	0.45
1:A:314:TRP:CE3	1:A:319:ARG:HD2	2.51	0.45
1:B:69:PHE:C	1:B:72:PRO:HD2	2.36	0.45
1:D:341:LEU:CD2	1:D:346:VAL:HG22	2.46	0.45
1:C:162:ILE:O	1:C:163:PRO:C	2.54	0.45
1:C:122:GLN:HE22	1:C:143:ARG:HD3	1.82	0.45
1:B:122:GLN:NE2	1:B:143:ARG:HH11	2.14	0.45
1:A:33:ALA:HB1	1:A:61:LEU:HD12	1.99	0.45
1:A:123:VAL:HA	1:B:380:ILE:HD12	1.98	0.45
1:C:298:ARG:C	1:C:298:ARG:HD2	2.37	0.45
1:C:351:ARG:CD	1:C:351:ARG:N	2.78	0.45
1:A:34:ASP:HB3	1:A:60:ASP:HA	1.98	0.45
1:C:159:ALA:O	1:C:161:GLY:N	2.50	0.45
1:D:88:THR:HG21	3:D:3520:CIR:H42	1.99	0.45
1:D:395:ARG:HH11	1:D:395:ARG:HG3	1.81	0.45
1:D:122:GLN:HG3	1:D:123:VAL:N	2.31	0.45
1:A:70:VAL:O	1:A:74:MET:HG3	2.17	0.45
1:A:374:LYS:HG3	1:D:374:LYS:HE3	1.98	0.45
1:A:32:THR:CG2	1:A:41:VAL:HG13	2.47	0.44
1:B:7:TYR:CE2	1:B:9:GLY:HA2	2.53	0.44
1:B:35:ILE:N	1:B:37:GLN:HE22	2.14	0.44
1:B:317:PRO:HB2	1:D:294:VAL:HG21	1.98	0.44
1:A:180:HIS:HB2	1:A:271:GLU:O	2.18	0.44
1:B:221:PHE:O	1:B:334:THR:OG1	2.25	0.44
1:C:343:LYS:HD2	1:D:343:LYS:HD3	1.99	0.44
1:C:185:GLY:O	1:C:188:LEU:HB2	2.18	0.44
1:C:302:SER:HB2	1:C:303:PRO:HD3	1.99	0.44
1:D:73:MET:CE	1:D:90:ILE:HG12	2.48	0.44
1:D:298:ARG:HD2	1:D:298:ARG:O	2.17	0.44
1:B:328:HIS:CE1	1:B:331:ARG:HH22	2.36	0.44
1:C:349:VAL:HB	1:D:194:GLU:HG3	2.00	0.44
1:A:34:ASP:HB2	1:A:41:VAL:HG21	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:150:ARG:HA	1:C:153:MET:HE3	2.00	0.44
1:A:117:GLY:HA2	1:A:122:GLN:OE1	2.17	0.44
1:C:41:VAL:HG12	1:C:58:ALA:HB1	1.99	0.44
1:A:6:ALA:HB2	1:A:99:LEU:HD11	2.00	0.44
1:A:69:PHE:CE1	1:A:94:LEU:HD23	2.53	0.44
1:D:378:GLY:O	1:D:382:ILE:HG12	2.18	0.44
1:B:174:MET:HG2	1:B:183:TYR:CD2	2.53	0.43
1:D:34:ASP:HA	1:D:37:GLN:HE22	1.83	0.43
1:A:15:ILE:HD11	1:A:157:ALA:HB2	2.00	0.43
1:C:34:ASP:HB2	1:C:41:VAL:HG21	2.00	0.43
1:D:395:ARG:NH1	1:D:395:ARG:HG3	2.34	0.43
1:A:135:ASP:OD2	1:A:135:ASP:N	2.50	0.43
1:A:264:MET:HG3	1:C:385:LEU:CD2	2.49	0.43
1:C:159:ALA:C	1:C:161:GLY:H	2.22	0.43
1:D:249:HIS:HD2	4:D:1218:HOH:O	2.01	0.43
1:D:84:TYR:CE2	1:D:86:LEU:HA	2.53	0.43
1:C:172:TYR:CZ	1:C:199:MET:HG3	2.53	0.43
1:B:298:ARG:HD2	1:B:298:ARG:C	2.38	0.43
1:B:35:ILE:H	1:B:37:GLN:HE22	1.57	0.43
1:D:40:GLU:OE2	1:D:40:GLU:HA	2.18	0.43
1:A:349:VAL:CG2	1:B:194:GLU:HG3	2.47	0.43
1:A:94:LEU:HD13	1:A:94:LEU:C	2.39	0.43
1:B:114:GLY:HA3	2:B:1510:ATP:O3'	2.19	0.43
1:C:40:GLU:HB3	1:C:43:GLU:CD	2.39	0.43
1:D:86:LEU:O	1:D:90:ILE:HG13	2.19	0.43
1:A:174:MET:HA	1:A:182:SER:O	2.19	0.43
1:B:86:LEU:O	1:B:90:ILE:HG13	2.19	0.43
1:C:73:MET:CE	1:C:90:ILE:HG12	2.49	0.43
1:D:114:GLY:CA	2:D:3510:ATP:O3'	2.67	0.43
1:A:27:GLU:CG	1:A:54:SER:OG	2.67	0.43
1:C:75:ARG:HG2	1:C:330:ALA:HB1	2.00	0.43
1:A:64:GLU:OE1	1:A:98:HIS:HE1	2.02	0.43
1:D:98:HIS:CD2	1:D:101:ARG:NH1	2.86	0.43
1:C:298:ARG:O	1:C:298:ARG:HD2	2.18	0.43
1:B:149:GLY:HA3	4:B:1432:HOH:O	2.19	0.43
1:D:254:VAL:HG22	1:D:256:ILE:HG23	2.00	0.43
1:B:116:THR:CB	1:B:118:LYS:HD2	2.49	0.43
1:A:189:GLU:O	1:A:191:PRO:HD3	2.19	0.42
1:D:392:LEU:O	1:D:395:ARG:CB	2.67	0.42
1:D:205:ASP:HA	1:D:206:PRO:HD3	1.96	0.42
1:B:392:LEU:O	1:B:395:ARG:HA	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:LYS:CB	1:A:118:LYS:NZ	2.81	0.42
1:A:84:TYR:CE2	1:A:86:LEU:HA	2.55	0.42
1:C:352:LYS:HE3	1:D:192:TRP:CH2	2.54	0.42
1:D:35:ILE:H	1:D:37:GLN:HE21	1.61	0.42
1:A:395:ARG:CG	1:A:395:ARG:HH11	2.31	0.42
1:B:70:VAL:O	1:B:74:MET:HG3	2.20	0.42
1:D:251:VAL:HG11	1:D:346:VAL:CG2	2.49	0.42
1:A:379:PHE:HB2	1:D:382:ILE:HG13	2.01	0.42
1:C:12:ASP:CG	1:C:150:ARG:NH2	2.73	0.42
1:B:122:GLN:HE22	1:B:143:ARG:HD3	1.83	0.42
1:B:261:PHE:C	1:B:263:GLY:H	2.23	0.42
1:D:114:GLY:HA3	2:D:3510:ATP:O3'	2.19	0.42
1:A:373:GLN:N	1:A:373:GLN:NE2	2.60	0.42
1:A:328:HIS:CD2	1:C:321:ALA:HA	2.55	0.42
1:C:60:ASP:OD1	1:C:62:LYS:CE	2.67	0.42
1:A:150:ARG:O	1:A:151:LYS:C	2.58	0.42
1:A:316:ALA:HB1	1:A:318:GLU:OE2	2.19	0.42
1:B:123:VAL:HG12	1:B:127:LEU:CD1	2.50	0.42
1:B:36:GLY:N	1:B:37:GLN:HE21	2.18	0.42
1:A:180:HIS:ND1	1:A:180:HIS:C	2.73	0.42
1:C:194:GLU:HG3	1:D:349:VAL:HB	2.02	0.42
1:C:357:LEU:HB2	1:D:265:LYS:CE	2.50	0.42
1:B:6:ALA:HB2	1:B:99:LEU:HD11	2.01	0.42
1:C:316:ALA:HB1	1:C:318:GLU:OE2	2.20	0.41
1:D:158:GLU:O	1:D:159:ALA:C	2.57	0.41
1:D:128:THR:HG22	4:D:1389:HOH:O	2.20	0.41
1:B:395:ARG:HG3	1:B:395:ARG:NH1	2.35	0.41
1:A:304:LYS:O	1:A:307:GLU:HG2	2.20	0.41
1:D:6:ALA:HB3	2:D:3510:ATP:HO2'	1.85	0.41
1:C:88:THR:HG21	3:C:2520:CIR:C4	2.50	0.41
1:B:345:ASN:HB3	1:B:347:TYR:CE1	2.54	0.41
1:B:254:VAL:CG1	1:B:256:ILE:HG23	2.50	0.41
1:C:29:ILE:HD12	1:C:55:LYS:HB3	2.03	0.41
1:B:60:ASP:OD1	1:B:62:LYS:CE	2.69	0.41
1:A:251:VAL:CG1	1:A:252:GLY:N	2.83	0.41
1:A:394:GLU:HA	1:A:394:GLU:OE1	2.19	0.41
1:B:2:LYS:HB2	1:B:108:ALA:HA	2.02	0.41
1:D:12:ASP:OD1	1:D:150:ARG:NH2	2.37	0.41
1:C:264:MET:N	1:C:264:MET:SD	2.94	0.41
1:D:90:ILE:H	1:D:90:ILE:HG13	1.61	0.41
1:D:146:SER:O	1:D:147:PHE:C	2.57	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:174:MET:HA	1:C:182:SER:O	2.21	0.41
1:B:85:LEU:HA	1:B:85:LEU:HD23	1.90	0.41
1:D:11:LEU:HD22	1:D:165:PRO:HD3	2.03	0.41
1:C:97:LYS:HD2	1:C:315:TYR:OH	2.21	0.41
1:A:393:VAL:O	1:A:395:ARG:NH1	2.51	0.41
1:B:328:HIS:HA	1:B:331:ARG:NH1	2.35	0.41
1:D:282:ARG:O	1:D:286:GLU:HG3	2.21	0.41
1:D:291:ASP:HB2	4:D:1396:HOH:O	2.21	0.41
1:B:128:THR:O	1:B:132:LEU:HG	2.20	0.41
1:A:12:ASP:HB2	2:A:510:ATP:O1G	2.21	0.41
1:D:196:PRO:HG2	1:D:199:MET:HB2	2.03	0.41
1:D:29:ILE:CD1	1:D:29:ILE:N	2.84	0.40
1:D:117:GLY:HA2	1:D:122:GLN:NE2	2.36	0.40
1:D:252:GLY:O	1:D:271:GLU:HA	2.21	0.40
1:C:60:ASP:OD1	1:C:62:LYS:HE3	2.21	0.40
1:A:251:VAL:HG11	1:A:346:VAL:CG2	2.51	0.40
1:D:353:ALA:HA	1:D:354:PRO:HD3	1.92	0.40
1:B:382:ILE:HG13	1:C:379:PHE:HB2	2.04	0.40
1:D:359:ARG:HG2	4:D:1365:HOH:O	2.21	0.40
1:B:41:VAL:HG12	1:B:58:ALA:CB	2.42	0.40
1:B:118:LYS:N	1:B:118:LYS:HD2	2.36	0.40
1:B:88:THR:CB	3:B:1520:CIR:H42	2.52	0.40
1:A:84:TYR:CZ	1:A:86:LEU:HA	2.56	0.40
1:A:197:LYS:HB2	1:A:197:LYS:HE3	1.83	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	374/400 (94%)	361 (96%)	13 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	374/400 (94%)	362 (97%)	12 (3%)	0	100	100
1	C	374/400 (94%)	354 (95%)	17 (4%)	3 (1%)	24	22
1	D	374/400 (94%)	361 (96%)	13 (4%)	0	100	100
All	All	1496/1600 (94%)	1438 (96%)	55 (4%)	3 (0%)	52	59

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	159	ALA
1	C	160	HIS
1	C	161	GLY

### 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	304/319 (95%)	281 (92%)	23 (8%)	16	16
1	B	304/319 (95%)	284 (93%)	20 (7%)	21	22
1	C	304/319 (95%)	287 (94%)	17 (6%)	26	29
1	D	304/319 (95%)	282 (93%)	22 (7%)	18	18
All	All	1216/1276 (95%)	1134 (93%)	82 (7%)	20	21

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	GLN
1	A	90	ILE
1	A	101	ARG
1	A	118	LYS
1	A	128	THR
1	A	137	LYS
1	A	145	TRP
1	A	151	LYS

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Mol	Chain	Res	Type
1	A	179	LEU
1	A	180	HIS
1	A	188	LEU
1	A	197	LYS
1	A	239	LEU
1	A	262	VAL
1	A	264	MET
1	A	267	ARG
1	A	291	ASP
1	A	318	GLU
1	A	319	ARG
1	A	345	ASN
1	A	349	VAL
1	A	351	ARG
1	A	373	GLN
1	B	37	GLN
1	B	39	GLU
1	B	62	LYS
1	B	90	ILE
1	B	118	LYS
1	B	128	THR
1	B	145	TRP
1	B	178	LEU
1	B	179	LEU
1	B	180	HIS
1	B	188	LEU
1	B	239	LEU
1	B	262	VAL
1	B	264	MET
1	B	267	ARG
1	B	291	ASP
1	B	318	GLU
1	B	319	ARG
1	B	351	ARG
1	B	359	ARG
1	C	37	GLN
1	C	62	LYS
1	C	90	ILE
1	C	145	TRP
1	C	150	ARG
1	C	151	LYS
1	C	179	LEU

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Mol	Chain	Res	Type
1	C	180	HIS
1	C	188	LEU
1	C	239	LEU
1	C	264	MET
1	C	267	ARG
1	C	318	GLU
1	C	319	ARG
1	C	351	ARG
1	C	373	GLN
1	C	374	LYS
1	D	25	ARG
1	D	37	GLN
1	D	62	LYS
1	D	90	ILE
1	D	94	LEU
1	D	128	THR
1	D	145	TRP
1	D	178	LEU
1	D	179	LEU
1	D	180	HIS
1	D	188	LEU
1	D	205	ASP
1	D	239	LEU
1	D	264	MET
1	D	267	ARG
1	D	291	ASP
1	D	318	GLU
1	D	319	ARG
1	D	351	ARG
1	D	372	ASP
1	D	394	GLU
1	D	395	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	HIS
1	A	148	GLN
1	A	249	HIS
1	A	323	GLN
1	A	345	ASN
1	A	373	GLN

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Mol	Chain	Res	Type
1	B	37	GLN
1	B	98	HIS
1	B	122	GLN
1	B	180	HIS
1	B	249	HIS
1	B	323	GLN
1	B	328	HIS
1	B	373	GLN
1	C	37	GLN
1	C	98	HIS
1	C	122	GLN
1	C	323	GLN
1	C	373	GLN
1	D	37	GLN
1	D	98	HIS
1	D	122	GLN
1	D	148	GLN
1	D	249	HIS
1	D	323	GLN
1	D	373	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

8 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
2	ATP	A	510	-	24,33,33	2.44	8 (33%)	31,52,52	3.58	13 (41%)
3	CIR	A	520	-	8,11,11	0.99	0	7,13,13	1.59	1 (14%)
2	ATP	B	1510	-	24,33,33	2.40	8 (33%)	31,52,52	3.67	13 (41%)
3	CIR	B	1520	-	8,11,11	1.00	1 (12%)	7,13,13	1.55	2 (28%)
2	ATP	C	2510	-	24,33,33	2.36	8 (33%)	31,52,52	3.46	12 (38%)
3	CIR	C	2520	-	8,11,11	1.07	1 (12%)	7,13,13	1.66	1 (14%)
2	ATP	D	3510	-	24,33,33	2.37	7 (29%)	31,52,52	3.34	12 (38%)
3	CIR	D	3520	-	8,11,11	0.97	0	7,13,13	1.69	2 (28%)

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	510	-	-	0/18/38/38	0/3/3/3
3	CIR	A	520	-	-	0/7/11/11	0/0/0/0
2	ATP	B	1510	-	-	0/18/38/38	0/3/3/3
3	CIR	B	1520	-	-	0/7/11/11	0/0/0/0
2	ATP	C	2510	-	-	0/18/38/38	0/3/3/3
3	CIR	C	2520	-	-	0/7/11/11	0/0/0/0
2	ATP	D	3510	-	-	0/18/38/38	0/3/3/3
3	CIR	D	3520	-	-	0/7/11/11	0/0/0/0

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	510	ATP	O5'-C5'	-5.66	1.21	1.44
2	B	1510	ATP	O5'-C5'	-5.48	1.22	1.44
2	C	2510	ATP	O5'-C5'	-5.37	1.22	1.44
2	D	3510	ATP	O5'-C5'	-4.96	1.24	1.44
2	A	510	ATP	PA-O5'	-3.36	1.43	1.59
2	C	2510	ATP	PA-O5'	-3.29	1.43	1.59
2	B	1510	ATP	PA-O5'	-3.23	1.44	1.59
2	D	3510	ATP	PA-O5'	-2.86	1.45	1.59
2	D	3510	ATP	C3'-C4'	-2.79	1.45	1.53
2	C	2510	ATP	C3'-C4'	-2.73	1.45	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1510	ATP	C3'-C4'	-2.73	1.45	1.53
2	A	510	ATP	C3'-C4'	-2.70	1.45	1.53
2	A	510	ATP	C5'-C4'	-2.58	1.43	1.51
2	B	1510	ATP	C8-N7	-2.44	1.29	1.34
2	C	2510	ATP	C8-N7	-2.38	1.30	1.34
2	C	2510	ATP	C5'-C4'	-2.35	1.44	1.51
2	D	3510	ATP	C8-N7	-2.29	1.30	1.34
2	B	1510	ATP	C5'-C4'	-2.26	1.44	1.51
2	A	510	ATP	C8-N7	-2.07	1.30	1.34
3	B	1520	CIR	C7-N6	-2.00	1.31	1.34
2	B	1510	ATP	C2-N3	2.00	1.35	1.32
3	C	2520	CIR	C5-N6	2.02	1.50	1.46
2	A	510	ATP	C2-N3	2.08	1.35	1.32
2	C	2510	ATP	C2-N3	2.23	1.36	1.32
2	D	3510	ATP	C2-N3	2.55	1.36	1.32
2	C	2510	ATP	C4-N3	4.64	1.42	1.35
2	A	510	ATP	C4-N3	4.79	1.42	1.35
2	B	1510	ATP	C4-N3	4.86	1.42	1.35
2	D	3510	ATP	C4-N3	5.34	1.43	1.35
2	B	1510	ATP	O4'-C1'	5.51	1.48	1.41
2	C	2510	ATP	O4'-C1'	5.58	1.48	1.41
2	D	3510	ATP	O4'-C1'	5.64	1.48	1.41
2	A	510	ATP	O4'-C1'	6.16	1.49	1.41

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1510	ATP	C5'-C4'-C3'	-6.94	87.68	115.21
2	A	510	ATP	C5'-C4'-C3'	-6.62	88.95	115.21
2	C	2510	ATP	C5'-C4'-C3'	-6.51	89.37	115.21
2	C	2510	ATP	O3A-PA-O5'	-6.42	85.91	102.94
2	B	1510	ATP	O3A-PA-O5'	-6.37	86.05	102.94
2	A	510	ATP	O3A-PA-O5'	-6.15	86.63	102.94
2	D	3510	ATP	C5'-C4'-C3'	-6.14	90.85	115.21
2	D	3510	ATP	O3A-PA-O5'	-5.72	87.77	102.94
2	A	510	ATP	C4'-O4'-C1'	-4.27	105.02	109.72
2	B	1510	ATP	C4'-O4'-C1'	-3.85	105.49	109.72
2	D	3510	ATP	N3-C2-N1	-3.63	126.12	128.89
2	A	510	ATP	N3-C2-N1	-3.61	126.13	128.89
2	D	3510	ATP	O2A-PA-O5'	-3.60	90.31	108.46
2	C	2510	ATP	C4'-O4'-C1'	-3.60	105.77	109.72
2	B	1510	ATP	N3-C2-N1	-3.54	126.18	128.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2510	ATP	N3-C2-N1	-3.47	126.24	128.89
2	B	1510	ATP	O2B-PB-O3A	-3.44	89.49	105.09
2	A	510	ATP	O2B-PB-O3A	-3.39	89.72	105.09
2	A	510	ATP	O2A-PA-O5'	-3.38	91.40	108.46
2	C	2510	ATP	O2B-PB-O3A	-3.35	89.87	105.09
2	D	3510	ATP	C4'-O4'-C1'	-3.31	106.08	109.72
2	C	2510	ATP	O2A-PA-O5'	-3.03	93.19	108.46
2	D	3510	ATP	O2B-PB-O3A	-3.00	91.50	105.09
2	B	1510	ATP	O2A-PA-O5'	-2.67	94.99	108.46
2	B	1510	ATP	O5'-PA-O1A	-2.50	99.92	109.62
2	C	2510	ATP	O5'-PA-O1A	-2.24	100.92	109.62
3	D	3520	CIR	O7-C7-N8	-2.10	119.71	123.30
3	B	1520	CIR	O7-C7-N8	-2.04	119.81	123.30
2	D	3510	ATP	O3'-C3'-C4'	2.08	117.28	111.05
2	B	1510	ATP	O3'-C3'-C2'	2.13	118.75	111.83
2	A	510	ATP	O3'-C3'-C4'	2.19	117.64	111.05
2	A	510	ATP	O3'-C3'-C2'	2.21	119.01	111.83
2	B	1510	ATP	O3'-C3'-C4'	2.24	117.77	111.05
2	C	2510	ATP	O2A-PA-O3A	2.33	115.65	105.09
2	A	510	ATP	O2A-PA-O3A	2.47	116.30	105.09
2	D	3510	ATP	O2A-PA-O3A	2.56	116.72	105.09
3	A	520	CIR	N8-C7-N6	2.63	119.39	116.17
3	B	1520	CIR	N8-C7-N6	2.67	119.43	116.17
2	D	3510	ATP	C4-C5-N7	2.69	111.95	109.48
3	C	2520	CIR	N8-C7-N6	2.83	119.63	116.17
2	A	510	ATP	C4-C5-N7	2.85	112.10	109.48
3	D	3520	CIR	N8-C7-N6	2.89	119.71	116.17
2	B	1510	ATP	C4-C5-N7	2.93	112.18	109.48
2	C	2510	ATP	C4-C5-N7	3.05	112.28	109.48
2	C	2510	ATP	O2B-PB-O3B	3.14	119.36	105.09
2	D	3510	ATP	O4'-C4'-C3'	3.18	111.55	105.15
2	D	3510	ATP	O2B-PB-O3B	3.24	119.81	105.09
2	A	510	ATP	O2B-PB-O3B	3.31	120.11	105.09
2	A	510	ATP	O4'-C4'-C3'	3.36	111.92	105.15
2	B	1510	ATP	O2B-PB-O3B	3.38	120.42	105.09
2	B	1510	ATP	O4'-C4'-C3'	3.49	112.17	105.15
2	C	2510	ATP	O4'-C4'-C3'	3.60	112.39	105.15
2	D	3510	ATP	O5'-C5'-C4'	12.90	156.69	109.12
2	C	2510	ATP	O5'-C5'-C4'	13.25	157.96	109.12
2	A	510	ATP	O5'-C5'-C4'	13.96	160.59	109.12
2	B	1510	ATP	O5'-C5'-C4'	14.63	163.06	109.12

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	510	ATP	3	0
3	A	520	CIR	1	0
2	B	1510	ATP	3	0
3	B	1520	CIR	2	0
2	C	2510	ATP	2	0
3	C	2520	CIR	2	0
2	D	3510	ATP	7	0
3	D	3520	CIR	1	0

## 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 ⓘ

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	380/400 (95%)	0.00	13 (3%) 49 47	24, 33, 55, 76	0
1	B	380/400 (95%)	-0.14	7 (1%) 71 70	23, 32, 52, 71	0
1	C	380/400 (95%)	0.26	26 (6%) 20 20	23, 37, 65, 83	0
1	D	380/400 (95%)	0.16	20 (5%) 30 29	24, 37, 61, 88	0
All	All	1520/1600 (95%)	0.07	66 (4%) 39 38	23, 35, 60, 88	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	165	PRO	7.7
1	D	165	PRO	6.7
1	A	165	PRO	4.9
1	C	263	GLY	4.9
1	C	162	ILE	4.4
1	C	39	GLU	4.1
1	D	164	VAL	4.0
1	C	159	ALA	3.8
1	C	18	LYS	3.8
1	D	159	ALA	3.8
1	C	25	ARG	3.6
1	C	161	GLY	3.4
1	C	49	LEU	3.3
1	A	370	GLY	3.1
1	C	28	VAL	3.0
1	C	108	ALA	3.0
1	D	154	ILE	3.0
1	D	371	TYR	3.0
1	C	154	ILE	2.9
1	A	162	ILE	2.9
1	D	41	VAL	2.9

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	D	163	PRO	2.9
1	C	79	VAL	2.9
1	C	163	PRO	2.8
1	A	263	GLY	2.8
1	D	44	ALA	2.7
1	C	50	ARG	2.7
1	C	164	VAL	2.7
1	A	164	VAL	2.7
1	A	154	ILE	2.7
1	A	262	VAL	2.6
1	D	161	GLY	2.6
1	B	264	MET	2.6
1	C	160	HIS	2.5
1	D	162	ILE	2.5
1	B	262	VAL	2.4
1	D	58	ALA	2.4
1	C	370	GLY	2.4
1	B	263	GLY	2.3
1	D	50	ARG	2.3
1	D	303	PRO	2.3
1	D	160	HIS	2.3
1	B	164	VAL	2.3
1	A	300	MET	2.2
1	C	51	THR	2.2
1	C	43	GLU	2.2
1	D	301	LEU	2.2
1	A	158	GLU	2.2
1	C	157	ALA	2.2
1	B	371	TYR	2.2
1	B	301	LEU	2.2
1	D	40	GLU	2.2
1	C	45	ARG	2.2
1	D	370	GLY	2.2
1	B	300	MET	2.2
1	C	1	MET	2.1
1	C	158	GLU	2.1
1	D	232	ARG	2.1
1	D	300	MET	2.1
1	D	395	ARG	2.1
1	A	163	PRO	2.1
1	A	41	VAL	2.1
1	A	157	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	300	MET	2.0
1	A	44	ALA	2.0
1	C	56	ALA	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	CIR	D	3520	12/12	0.85	0.24	5.35	28,40,60,61	0
3	CIR	A	520	12/12	0.86	0.21	4.30	27,37,47,49	0
2	ATP	B	1510	31/31	0.88	0.16	4.13	27,41,83,86	0
3	CIR	C	2520	12/12	0.84	0.20	2.85	27,38,49,50	0
2	ATP	A	510	31/31	0.86	0.16	2.75	36,44,90,94	0
3	CIR	B	1520	12/12	0.88	0.18	2.61	24,32,44,45	0
2	ATP	D	3510	31/31	0.82	0.22	2.11	50,58,94,99	0
2	ATP	C	2510	31/31	0.52	0.28	1.68	0,45,95,97	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.