



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:29 PM GMT

PDB ID : 1KH3  
Title : Crystal Structure of Thermus thermophilus HB8 Argininosuccinate Synthetase in complex with inhibitor  
Authors : goto, m.; Hirotsu, k.; miyahara, i.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)  
Deposited on : 2001-11-29  
Resolution : 2.15 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](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

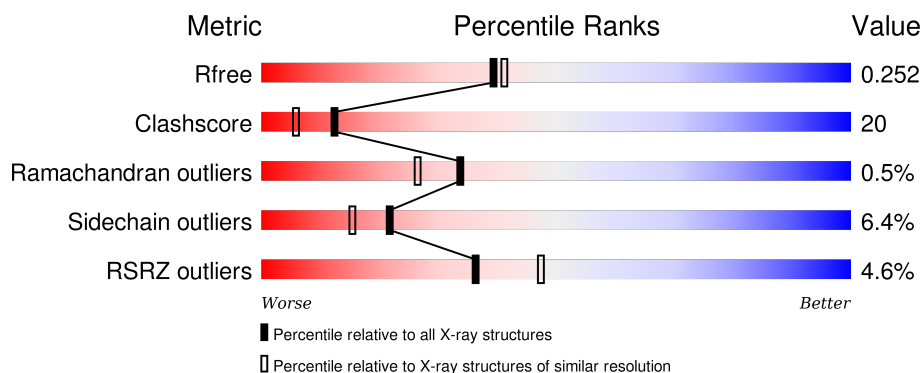
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.15 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	1045 (2.16-2.16)
Clashscore	102246	1152 (2.16-2.16)
Ramachandran outliers	100387	1131 (2.16-2.16)
Sidechain outliers	100360	1131 (2.16-2.16)
RSRZ outliers	91569	1050 (2.16-2.16)

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>2%</div> <div>66% 25% 5%</div> </div>
1	B	400	<div> <div>4%</div> <div>67% 28%</div> </div>
1	C	400	<div> <div>6%</div> <div>65% 28%</div> </div>
1	D	400	<div> <div>6%</div> <div>61% 33%</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	SO4	A	540	-	-	-	X
2	SO4	C	2540	-	-	-	X
2	SO4	D	3540	-	-	-	X
5	ARG	A	2520	-	-	-	X
5	ARG	A	3520	-	-	-	X
5	ARG	A	4520	-	-	-	X
5	ARG	A	520	-	-	-	X

## 2 Entry composition [i](#)

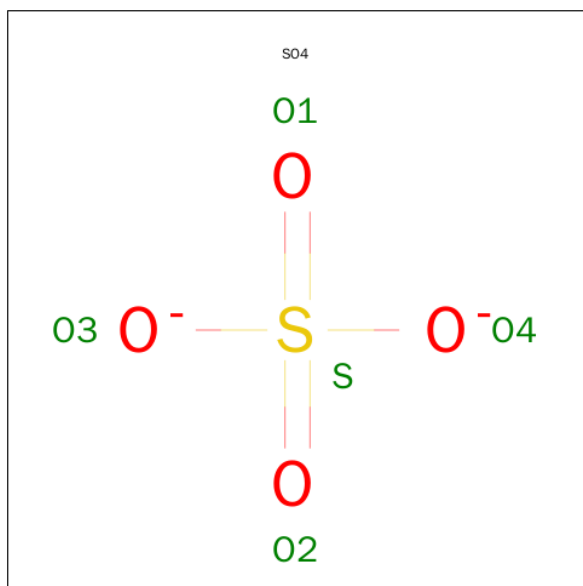
There are 6 unique types of molecules in this entry. The entry contains 13014 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	386	Total	C	N	O	S	0	0	0
			3066	1963	533	561	9			
1	C	386	Total	C	N	O	S	0	0	0
			3066	1963	533	561	9			
1	D	386	Total	C	N	O	S	0	0	0
			3066	1963	533	561	9			

- Molecule 2 is SULFATE ION (three-letter code: SO<sub>4</sub>) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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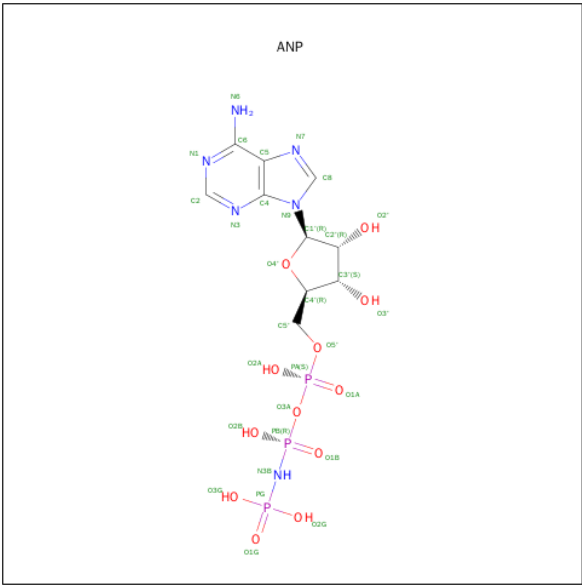
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: C<sub>10</sub>H<sub>17</sub>N<sub>6</sub>O<sub>12</sub>P<sub>3</sub>).



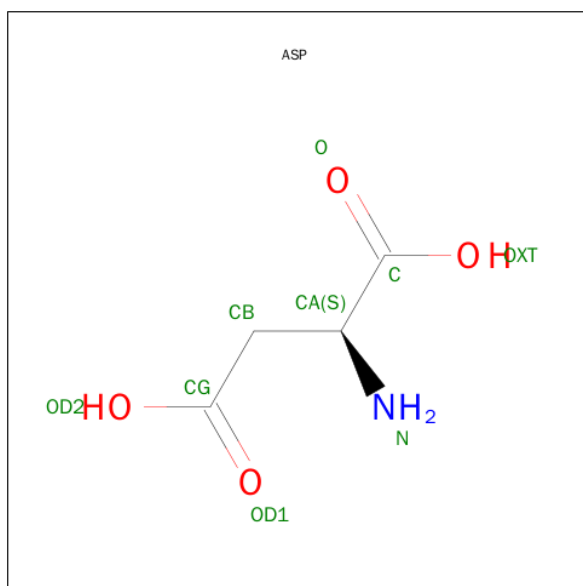
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
4	B	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
4	C	1	Total	C	N	O	P	0	0
			31	10	6	12	3		

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

- Molecule 5 is ARGinine (three-letter code: ARG) (formula:  $C_6H_{15}N_4O_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	8	Total	C	N	O	0	0
			84	40	20	24		

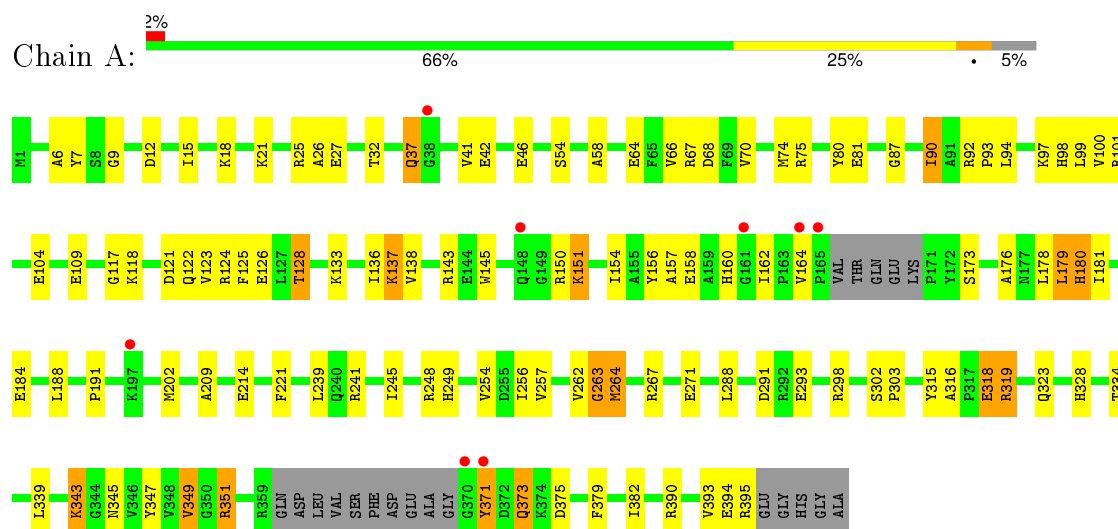
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	160	Total	O	0	0
			160	160		
6	B	161	Total	O	0	0
			161	161		
6	C	125	Total	O	0	0
			125	125		
6	D	121	Total	O	0	0
			121	121		

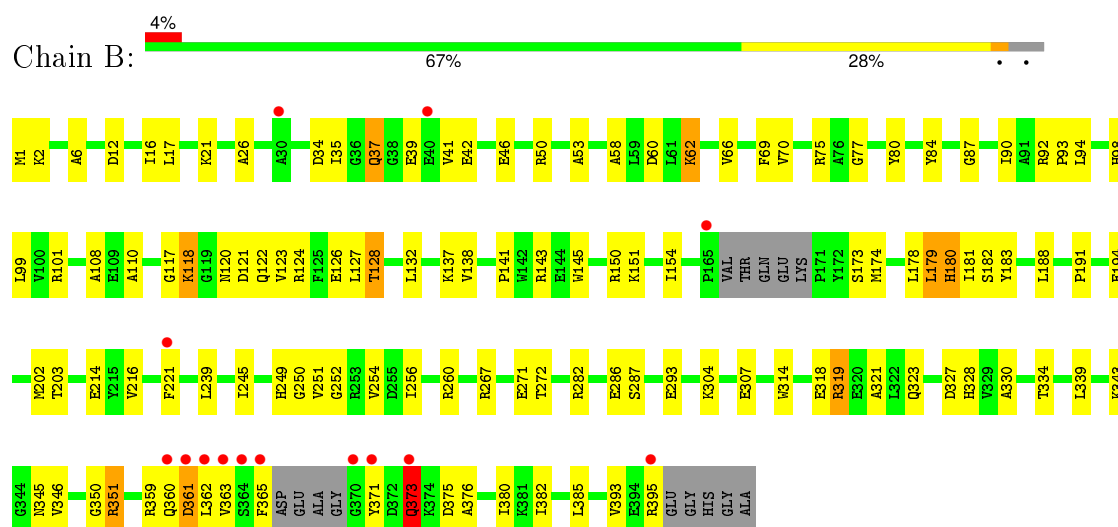
### 3 Residue-property plots [i](#)

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

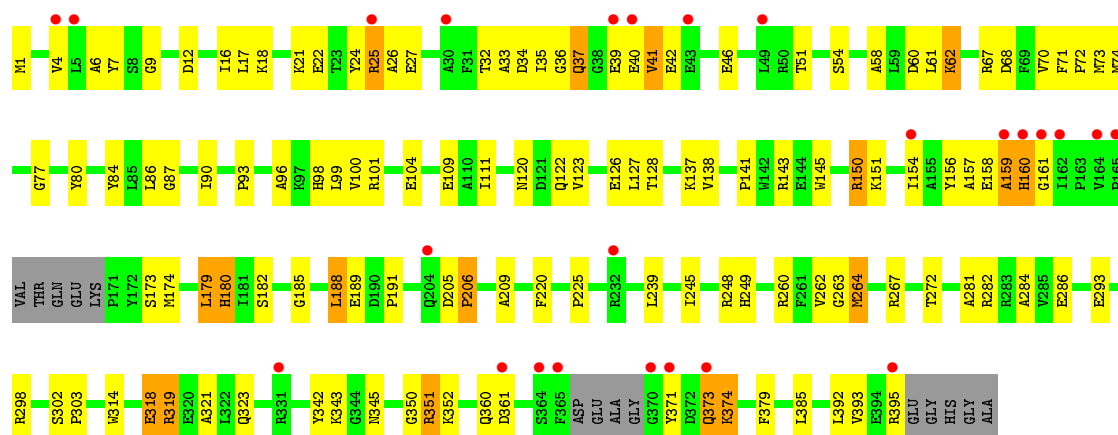


#### • Molecule 1: Argininosuccinate Synthetase

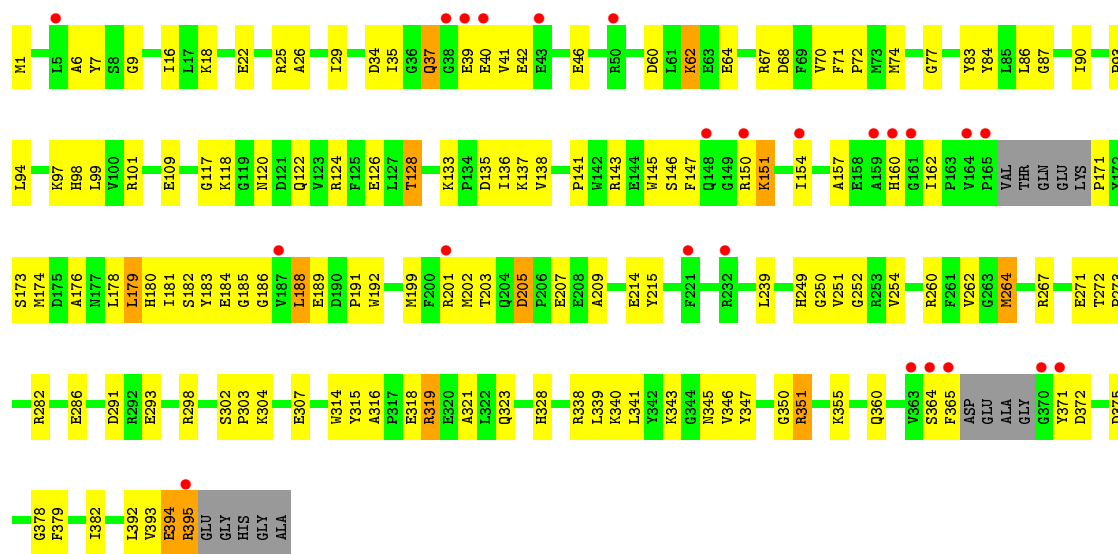


#### • Molecule 1: Argininosuccinate Synthetase





• Molecule 1: Argininosuccinate Synthetase





## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	229.10Å 229.10Å 160.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.15 19.99 – 2.15	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.15) 99.6 (19.99-2.15)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.48 (at 2.15Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.226 , 0.256 0.223 , 0.252	Depositor DCC
$R_{free}$ test set	16698 reflections (9.94%)	DCC
Wilson B-factor (Å <sup>2</sup> )	30.9	Xtriage
Anisotropy	0.298	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 48.0	EDS
Estimated twinning fraction	0.003 for h,-h-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.36$	Xtriage
Outliers	0 of 169389 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	13014	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.57% 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: MG, ANP, SO4

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	1/4169 (0.0%)
1	B	0.38	0/3135	0.63	0/4237
1	C	0.38	0/3135	0.64	1/4237 (0.0%)
1	D	0.38	0/3135	0.62	0/4237
All	All	0.38	0/12490	0.63	2/16880 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	263	GLY	N-CA-C	5.57	127.02	113.10
1	C	263	GLY	N-CA-C	5.15	125.96	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	124	0
1	B	3066	0	3052	124	0
1	C	3066	0	3052	150	0
1	D	3066	0	3052	138	0
2	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	31	0	13	1	0
4	B	31	0	13	1	0
4	C	31	0	13	0	0
4	D	31	0	13	2	0
5	A	84	0	60	17	0
6	A	160	0	0	6	0
6	B	161	0	0	12	0
6	C	125	0	0	7	0
6	D	121	0	0	4	0
All	All	13014	0	12274	485	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 20.

All (485) 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	1.01	1.10
1:A:151:LYS:H	1:A:151:LYS:HD2	1.11	1.10
1:B:151:LYS:HD2	1:B:151:LYS:H	1.14	1.06
1:B:35:ILE:H	1:B:37:GLN:NE2	1.53	1.06
1:A:123:VAL:HG22	1:B:380:ILE:HD11	1.41	1.02
1:A:124:ARG:O	1:A:128:THR:HG22	1.62	0.99
1:D:35:ILE:H	1:D:37:GLN:NE2	1.60	0.99
1:A:262:VAL:HG23	1:A:264:MET:HE1	1.46	0.98
1:C:151:LYS:HD2	1:C:151:LYS:H	1.24	0.98
1:A:262:VAL:HG23	1:A:264:MET:CE	1.93	0.97
1:D:124:ARG:O	1:D:128:THR:HG22	1.67	0.94
1:C:109:GLU:O	1:C:137:LYS:HE2	1.71	0.90
1:A:373:GLN:H	1:A:373:GLN:NE2	1.68	0.89
1:D:151:LYS:N	1:D:151:LYS:HD2	1.86	0.89
1:A:151:LYS:N	1:A:151:LYS:HD2	1.88	0.89
1:B:41:VAL:HG12	1:B:58:ALA:HB1	1.55	0.88
1:D:35:ILE:H	1:D:37:GLN:HE22	1.21	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:118:LYS:H	1:B:118:LYS:HD2	1.39	0.85
1:D:351:ARG:HD2	1:D:351:ARG:N	1.92	0.85
1:A:293:GLU:CB	1:C:318:GLU:HG2	2.07	0.84
1:C:37:GLN:H	1:C:37:GLN:NE2	1.76	0.83
1:B:124:ARG:O	1:B:128:THR:HG23	1.77	0.83
1:A:67:ARG:HH11	1:A:67:ARG:HG2	1.44	0.83
1:D:151:LYS:H	1:D:151:LYS:CD	1.80	0.81
1:B:151:LYS:CD	1:B:151:LYS:H	1.92	0.81
5:A:3520:ARG:HH22	1:D:173:SER:C	1.85	0.80
1:C:122:GLN:HE21	1:C:143:ARG:HH11	1.29	0.80
1:C:35:ILE:H	1:C:37:GLN:NE2	1.77	0.80
1:B:35:ILE:N	1:B:37:GLN:NE2	2.29	0.80
1:C:1:MET:O	1:C:26:ALA:HB1	1.82	0.80
1:A:173:SER:C	5:A:520:ARG:HH22	1.87	0.78
1:C:122:GLN:NE2	1:C:143:ARG:HH11	1.82	0.78
1:A:293:GLU:HB2	1:C:318:GLU:CG	2.13	0.77
1:B:359:ARG:HH21	1:B:362:LEU:HD21	1.48	0.77
1:C:351:ARG:N	1:C:351:ARG:HD2	2.00	0.77
1:B:293:GLU:CB	1:D:318:GLU:HG2	2.15	0.76
1:B:151:LYS:HD2	1:B:151:LYS:N	1.98	0.76
1:A:100:VAL:O	1:A:104:GLU:HG3	1.83	0.76
5:A:4520:ARG:HH22	1:B:173:SER:C	1.89	0.76
5:A:2520:ARG:HH22	1:C:173:SER:C	1.88	0.76
1:B:314:TRP:O	1:B:319:ARG:HD3	1.86	0.75
1:A:293:GLU:HB2	1:C:318:GLU:HG2	1.68	0.75
1:D:83:TYR:HB3	1:D:254:VAL:HG21	1.69	0.74
1:D:18:LYS:HG3	1:D:162:ILE:HD11	1.70	0.74
1:D:319:ARG:HH21	1:D:323:GLN:HE21	1.34	0.73
1:C:37:GLN:H	1:C:37:GLN:HE21	1.36	0.73
1:B:351:ARG:N	1:B:351:ARG:HD2	2.04	0.72
1:A:293:GLU:HB3	1:C:318:GLU:HG2	1.71	0.72
1:B:293:GLU:HB3	1:D:318:GLU:HG2	1.70	0.71
1:B:41:VAL:CG1	1:B:58:ALA:HB1	2.19	0.71
1:C:319:ARG:HH21	1:C:323:GLN:NE2	1.88	0.71
1:A:349:VAL:HG22	1:B:194:GLU:HG3	1.73	0.70
1:C:262:VAL:O	1:C:262:VAL:HG12	1.90	0.69
1:A:318:GLU:HG2	1:C:293:GLU:CB	2.22	0.69
1:B:318:GLU:HG3	1:D:293:GLU:HB2	1.74	0.69
1:B:87:GLY:HA3	1:B:180:HIS:CD2	2.27	0.69
1:D:109:GLU:O	1:D:137:LYS:HE2	1.91	0.69
1:C:21:LYS:O	1:C:25:ARG:HA	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:393:VAL:O	1:D:395:ARG:HG3	1.92	0.69
1:B:118:LYS:N	1:B:118:LYS:HD2	2.09	0.68
1:C:245:ILE:O	1:C:248:ARG:HG2	1.93	0.68
1:C:86:LEU:O	1:C:90:ILE:HG13	1.94	0.68
5:A:4520:ARG:NH2	1:B:173:SER:O	2.25	0.68
1:B:293:GLU:HB2	1:D:318:GLU:CG	2.24	0.68
1:A:109:GLU:O	1:A:137:LYS:HE2	1.94	0.68
1:C:17:LEU:HD13	1:C:51:THR:HB	1.76	0.67
1:D:124:ARG:O	1:D:128:THR:CG2	2.42	0.67
1:A:351:ARG:N	1:A:351:ARG:HD2	2.09	0.67
1:A:93:PRO:HG3	6:A:1201:HOH:O	1.93	0.67
1:A:390:ARG:O	1:A:394:GLU:HG2	1.94	0.67
1:A:262:VAL:HG23	1:A:264:MET:HE3	1.74	0.67
1:C:319:ARG:HH21	1:C:323:GLN:HE21	1.42	0.67
1:A:37:GLN:HG2	6:A:1494:HOH:O	1.95	0.67
1:A:371:TYR:HD1	1:A:375:ASP:HB2	1.60	0.67
1:C:374:LYS:HE2	6:C:1162:HOH:O	1.95	0.66
1:A:319:ARG:HH21	1:A:323:GLN:NE2	1.93	0.66
1:A:318:GLU:HG2	1:C:293:GLU:HB3	1.78	0.66
1:A:319:ARG:HH21	1:A:323:GLN:HE21	1.43	0.66
1:C:156:TYR:O	1:C:159:ALA:HB3	1.96	0.66
1:C:109:GLU:HG3	1:C:137:LYS:HE3	1.77	0.66
1:C:343:LYS:HD2	1:D:343:LYS:HD2	1.78	0.66
1:C:41:VAL:HG12	1:C:58:ALA:HB1	1.78	0.66
1:D:62:LYS:HD3	1:D:62:LYS:H	1.62	0.65
1:B:318:GLU:CG	1:D:293:GLU:HB2	2.26	0.65
1:B:37:GLN:HG2	6:B:1237:HOH:O	1.95	0.64
1:A:373:GLN:H	1:A:373:GLN:HE21	1.46	0.64
1:A:318:GLU:CG	1:C:293:GLU:HB2	2.27	0.64
1:A:42:GLU:O	1:A:46:GLU:HG3	1.96	0.64
1:C:151:LYS:CD	1:C:151:LYS:H	1.99	0.64
6:A:1551:HOH:O	1:B:345:ASN:HB2	1.98	0.64
1:A:151:LYS:CD	1:A:151:LYS:H	1.89	0.64
1:B:216:VAL:HG22	6:B:1549:HOH:O	1.97	0.63
1:B:75:ARG:HG2	1:B:330:ALA:CB	2.28	0.63
1:D:37:GLN:NE2	1:D:37:GLN:H	1.96	0.63
1:D:83:TYR:CB	1:D:254:VAL:HG21	2.28	0.63
1:C:60:ASP:OD1	1:C:62:LYS:HE3	1.99	0.63
1:C:122:GLN:HE21	1:C:143:ARG:NH1	1.95	0.62
1:C:371:TYR:HA	6:C:1158:HOH:O	1.99	0.62
1:B:360:GLN:HA	1:B:363:VAL:HG22	1.79	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:159:ALA:O	1:C:161:GLY:N	2.32	0.62
1:B:35:ILE:H	1:B:37:GLN:HE22	1.42	0.62
1:D:214:GLU:OE1	1:D:249:HIS:HE1	1.82	0.62
1:B:122:GLN:HE21	1:B:143:ARG:HH11	1.48	0.61
1:D:395:ARG:HH11	1:D:395:ARG:HG3	1.66	0.61
1:B:314:TRP:CE3	1:B:319:ARG:HD2	2.36	0.61
1:A:67:ARG:HH11	1:A:67:ARG:CG	2.14	0.61
1:B:42:GLU:O	1:B:46:GLU:HG3	2.00	0.61
1:A:67:ARG:NH1	1:A:68:ASP:OD1	2.33	0.60
1:C:98:HIS:HD2	1:C:101:ARG:NH1	2.00	0.60
1:B:293:GLU:HB2	1:D:318:GLU:HG2	1.81	0.60
1:A:41:VAL:HG12	1:A:58:ALA:HB1	1.84	0.60
6:C:1251:HOH:O	1:D:191:PRO:HG2	2.01	0.60
1:C:343:LYS:HD2	1:D:343:LYS:CD	2.32	0.60
1:B:117:GLY:HA2	1:B:122:GLN:NE2	2.16	0.60
1:B:282:ARG:O	1:B:286:GLU:HG3	2.01	0.60
1:A:67:ARG:HH12	1:A:68:ASP:CG	2.03	0.60
1:C:262:VAL:HG11	1:D:371:TYR:OH	2.02	0.60
1:C:18:LYS:NZ	1:C:160:HIS:O	2.34	0.60
1:B:50:ARG:NH1	6:B:1386:HOH:O	2.34	0.60
1:C:21:LYS:O	1:C:25:ARG:CA	2.50	0.60
1:A:123:VAL:HG22	1:B:380:ILE:CD1	2.25	0.60
1:C:248:ARG:HG3	1:C:249:HIS:CD2	2.37	0.60
1:B:117:GLY:HA2	1:B:122:GLN:HE22	1.67	0.59
1:C:93:PRO:HG3	6:C:1365:HOH:O	2.02	0.59
1:C:150:ARG:O	1:C:154:ILE:HG13	2.00	0.59
1:D:395:ARG:HB2	1:D:395:ARG:CZ	2.32	0.59
1:C:395:ARG:HH11	1:C:395:ARG:HG3	1.67	0.59
1:B:1:MET:O	1:B:26:ALA:HB1	2.03	0.59
1:A:245:ILE:O	1:A:248:ARG:HG2	2.03	0.59
1:B:318:GLU:HG2	1:D:293:GLU:HB3	1.85	0.59
1:D:67:ARG:HD2	1:D:68:ASP:OD1	2.02	0.59
1:A:173:SER:O	5:A:520:ARG:NH2	2.36	0.59
1:D:7:TYR:CE2	1:D:9:GLY:HA2	2.38	0.59
1:C:41:VAL:CG1	1:C:58:ALA:HB1	2.33	0.58
1:D:6:ALA:HB2	1:D:99:LEU:HD11	1.83	0.58
1:C:109:GLU:CG	1:C:137:LYS:HE3	2.33	0.58
1:A:27:GLU:HG3	1:A:54:SER:HB3	1.85	0.58
1:D:93:PRO:HG3	6:D:1516:HOH:O	2.02	0.58
1:B:126:GLU:HG2	1:B:138:VAL:HG11	1.84	0.58
1:B:93:PRO:HA	1:B:128:THR:HG21	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:GLU:HG2	1:A:133:LYS:HZ2	1.69	0.58
1:A:6:ALA:HB2	1:A:99:LEU:HD11	1.86	0.58
1:D:282:ARG:O	1:D:286:GLU:HG3	2.03	0.57
1:B:93:PRO:HG3	6:B:1266:HOH:O	2.04	0.57
1:A:302:SER:HB2	1:A:303:PRO:HD3	1.86	0.57
1:A:264:MET:H	1:A:264:MET:CE	2.18	0.57
1:A:104:GLU:OE2	1:A:133:LYS:NZ	2.33	0.57
1:B:35:ILE:N	1:B:37:GLN:HE22	2.01	0.57
1:A:382:ILE:HD12	1:C:262:VAL:HG13	1.86	0.57
1:A:318:GLU:HG2	1:C:293:GLU:HB2	1.87	0.57
1:B:371:TYR:HD1	1:B:375:ASP:HB2	1.70	0.57
1:A:345:ASN:HB3	1:A:347:TYR:CE1	2.40	0.57
1:C:17:LEU:O	1:C:17:LEU:HD23	2.04	0.56
1:C:18:LYS:HD2	1:C:160:HIS:HB3	1.85	0.56
1:A:98:HIS:HD2	1:A:101:ARG:NH1	2.03	0.56
1:D:360:GLN:HG2	1:D:364:SER:OG	2.05	0.56
1:C:209:ALA:O	1:C:343:LYS:NZ	2.39	0.56
1:C:18:LYS:NZ	1:C:160:HIS:HB3	2.20	0.56
1:B:37:GLN:H	1:B:37:GLN:NE2	2.04	0.56
5:A:3520:ARG:HH21	1:D:184:GLU:HG2	1.71	0.56
1:B:318:GLU:CG	1:D:293:GLU:CB	2.84	0.56
1:D:35:ILE:N	1:D:37:GLN:HE22	1.97	0.55
1:A:373:GLN:N	1:A:373:GLN:NE2	2.48	0.55
1:B:34:ASP:HB2	1:B:41:VAL:HG21	1.87	0.55
1:C:126:GLU:HG2	1:C:138:VAL:HG11	1.88	0.55
1:A:104:GLU:HG2	1:A:133:LYS:NZ	2.20	0.55
1:A:125:PHE:O	1:A:128:THR:HG23	2.07	0.55
1:A:150:ARG:HG3	1:A:154:ILE:HD13	1.89	0.55
1:D:126:GLU:HG2	1:D:138:VAL:HG11	1.87	0.55
1:D:351:ARG:CD	1:D:351:ARG:N	2.69	0.55
1:B:318:GLU:HG2	1:D:293:GLU:CB	2.37	0.55
1:D:360:GLN:HG2	1:D:364:SER:HG	1.71	0.55
1:D:176:ALA:CB	1:D:181:ILE:HG12	2.36	0.55
1:B:359:ARG:NH2	1:B:362:LEU:HD21	2.20	0.55
1:A:382:ILE:HD12	1:C:262:VAL:CG1	2.38	0.54
1:A:87:GLY:HA2	1:A:179:LEU:HD13	1.87	0.54
1:B:6:ALA:HB2	1:B:99:LEU:HD11	1.89	0.54
1:C:21:LYS:O	1:C:25:ARG:N	2.41	0.54
1:C:220:PHE:CE2	1:C:225:PRO:HG3	2.42	0.54
1:D:87:GLY:HA2	1:D:179:LEU:HD13	1.89	0.54
1:A:351:ARG:HD3	1:B:191:PRO:HB2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:98:HIS:CD2	1:C:101:ARG:NH1	2.76	0.54
1:C:393:VAL:O	1:C:395:ARG:NH1	2.40	0.54
1:A:75:ARG:NH2	6:A:1451:HOH:O	2.41	0.54
1:C:373:GLN:HG2	1:C:374:LYS:N	2.22	0.54
1:A:90:ILE:HD11	1:A:179:LEU:HD12	1.90	0.54
1:C:109:GLU:O	1:C:137:LYS:CE	2.51	0.54
1:C:37:GLN:HB2	1:C:39:GLU:HG2	1.90	0.54
1:B:216:VAL:HG21	1:B:245:ILE:HG21	1.90	0.54
1:D:1:MET:O	1:D:26:ALA:HB1	2.08	0.54
1:B:382:ILE:HG13	1:C:379:PHE:HB2	1.90	0.54
1:D:29:ILE:N	1:D:29:ILE:HD12	2.23	0.54
1:D:42:GLU:O	1:D:46:GLU:HG3	2.08	0.54
1:A:214:GLU:OE1	1:A:249:HIS:HE1	1.91	0.54
1:D:22:GLU:CD	1:D:160:HIS:HE2	2.11	0.54
1:B:118:LYS:CD	1:B:118:LYS:N	2.70	0.53
1:D:371:TYR:HB2	1:D:375:ASP:OD1	2.08	0.53
1:D:122:GLN:HE21	1:D:143:ARG:HH11	1.57	0.53
1:B:249:HIS:HD2	6:B:1070:HOH:O	1.90	0.53
1:D:338:ARG:HD3	6:D:1261:HOH:O	2.09	0.53
1:D:345:ASN:HB3	1:D:347:TYR:CE1	2.44	0.53
5:A:3520:ARG:NH2	1:D:173:SER:O	2.42	0.53
5:A:2520:ARG:NH2	1:C:173:SER:O	2.42	0.53
1:D:120:ASN:HD21	1:D:260:ARG:HG2	1.74	0.53
1:B:98:HIS:HD2	1:B:101:ARG:NH1	2.07	0.53
1:D:314:TRP:O	1:D:319:ARG:HD3	2.09	0.53
1:A:98:HIS:CD2	1:A:101:ARG:NH1	2.76	0.53
1:A:395:ARG:HH22	1:D:355:LYS:HE3	1.72	0.53
1:C:360:GLN:OE1	1:D:189:GLU:OE2	2.27	0.53
1:C:374:LYS:HE3	1:D:143:ARG:O	2.08	0.53
1:D:205:ASP:OD1	1:D:207:GLU:OE2	2.27	0.53
1:C:73:MET:SD	1:C:90:ILE:HG21	2.49	0.52
1:A:41:VAL:CG1	1:A:58:ALA:HB1	2.40	0.52
1:A:221:PHE:O	1:A:334:THR:OG1	2.17	0.52
1:B:75:ARG:NH2	6:B:1271:HOH:O	2.42	0.52
1:A:21:LYS:HD2	1:A:26:ALA:O	2.10	0.52
1:B:124:ARG:O	1:B:128:THR:CG2	2.55	0.52
1:D:117:GLY:HA2	1:D:122:GLN:NE2	2.24	0.52
1:D:118:LYS:O	1:D:264:MET:HE1	2.09	0.52
1:A:126:GLU:HG2	1:A:138:VAL:HG11	1.91	0.52
1:B:395:ARG:HH11	1:B:395:ARG:HG3	1.73	0.52
1:C:35:ILE:N	1:C:37:GLN:NE2	2.53	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:LYS:HG2	1:A:136:ILE:HB	1.91	0.52
1:C:150:ARG:HB2	1:C:150:ARG:HH11	1.75	0.52
1:B:180:HIS:ND1	1:B:180:HIS:C	2.62	0.52
1:C:373:GLN:HG2	1:C:374:LYS:H	1.75	0.52
1:A:70:VAL:O	1:A:74:MET:HG3	2.10	0.52
1:D:314:TRP:CE3	1:D:319:ARG:HD2	2.44	0.52
1:B:87:GLY:HA2	1:B:179:LEU:HD13	1.92	0.52
1:D:40:GLU:OE1	1:D:42:GLU:HB2	2.09	0.52
1:D:98:HIS:HD2	1:D:101:ARG:NH1	2.08	0.52
1:A:7:TYR:CE2	1:A:9:GLY:HA2	2.45	0.51
1:B:221:PHE:O	1:B:334:THR:OG1	2.15	0.51
1:B:110:ALA:HB2	1:B:137:LYS:HB3	1.92	0.51
1:C:25:ARG:NH1	1:C:25:ARG:HB3	2.26	0.51
1:D:137:LYS:HD3	1:D:137:LYS:N	2.26	0.51
1:A:137:LYS:N	1:A:137:LYS:HD3	2.26	0.51
1:C:343:LYS:CD	1:D:343:LYS:HD2	2.38	0.51
1:A:249:HIS:HD2	6:A:1275:HOH:O	1.91	0.51
1:C:282:ARG:O	1:C:286:GLU:HG3	2.11	0.51
1:A:379:PHE:HB2	1:D:382:ILE:HG13	1.93	0.51
1:D:319:ARG:HH21	1:D:323:GLN:NE2	2.04	0.51
1:D:70:VAL:O	1:D:74:MET:HG3	2.10	0.51
1:B:34:ASP:HA	1:B:37:GLN:HE22	1.75	0.51
1:C:158:GLU:O	1:C:159:ALA:O	2.29	0.51
1:C:342:TYR:OH	1:D:207:GLU:HA	2.11	0.51
1:D:133:LYS:O	1:D:136:ILE:HG22	2.11	0.51
1:D:117:GLY:HA2	1:D:122:GLN:HE22	1.76	0.51
1:B:371:TYR:HB2	1:B:375:ASP:OD1	2.11	0.50
5:A:3520:ARG:NH2	1:D:184:GLU:HG2	2.25	0.50
1:B:120:ASN:HD21	1:B:260:ARG:HG2	1.76	0.50
1:C:262:VAL:O	1:C:262:VAL:CG1	2.60	0.50
1:B:328:HIS:CD2	1:D:321:ALA:HA	2.46	0.50
1:B:77:GLY:HA2	6:B:1168:HOH:O	2.10	0.50
1:D:350:GLY:C	1:D:351:ARG:HD2	2.31	0.50
1:D:304:LYS:O	1:D:307:GLU:HG2	2.12	0.50
1:C:185:GLY:O	1:C:188:LEU:HB2	2.11	0.50
1:C:302:SER:HB2	1:C:303:PRO:HD3	1.93	0.50
1:C:351:ARG:CD	1:C:351:ARG:N	2.72	0.50
1:C:350:GLY:C	1:C:351:ARG:HD2	2.32	0.50
1:D:339:LEU:N	1:D:339:LEU:HD12	2.26	0.50
1:B:181:ILE:HD12	1:B:202:MET:HB2	1.93	0.50
1:C:25:ARG:CB	1:C:25:ARG:NH1	2.74	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:97:LYS:HD2	1:D:315:TYR:OH	2.12	0.50
1:C:32:THR:CG2	1:C:41:VAL:HG13	2.42	0.49
1:B:319:ARG:HH21	1:B:323:GLN:HE21	1.60	0.49
1:D:16:ILE:HG23	1:D:141:PRO:HG2	1.94	0.49
1:C:34:ASP:HA	1:C:37:GLN:HE22	1.77	0.49
1:B:359:ARG:C	1:B:361:ASP:H	2.16	0.49
1:A:180:HIS:C	1:A:180:HIS:ND1	2.65	0.49
1:D:150:ARG:O	1:D:154:ILE:HD13	2.11	0.49
1:B:350:GLY:C	1:B:351:ARG:HD2	2.32	0.49
1:A:298:ARG:C	1:A:298:ARG:HD2	2.33	0.49
1:D:316:ALA:HB1	1:D:318:GLU:OE2	2.12	0.49
1:B:12:ASP:OD1	1:B:150:ARG:NH2	2.39	0.49
1:C:96:ALA:HB3	1:C:128:THR:HG21	1.92	0.49
1:A:293:GLU:HB2	1:C:318:GLU:HG3	1.92	0.49
1:B:150:ARG:O	1:B:154:ILE:HD13	2.11	0.49
1:B:2:LYS:HB2	1:B:108:ALA:HA	1.95	0.49
1:C:24:TYR:O	1:C:25:ARG:C	2.50	0.49
1:B:90:ILE:HD11	1:B:179:LEU:HD12	1.94	0.49
1:C:6:ALA:HB2	1:C:99:LEU:HD11	1.95	0.49
1:A:191:PRO:HG2	6:B:1561:HOH:O	2.13	0.48
1:A:32:THR:CG2	1:A:41:VAL:HG13	2.44	0.48
1:B:214:GLU:OE1	1:B:249:HIS:HE1	1.96	0.48
1:C:27:GLU:HG3	1:C:54:SER:CB	2.42	0.48
1:B:373:GLN:O	1:B:376:ALA:HB3	2.12	0.48
1:D:209:ALA:O	1:D:343:LYS:NZ	2.47	0.48
1:A:27:GLU:HG3	1:A:54:SER:CB	2.43	0.48
1:D:365:PHE:N	6:D:1390:HOH:O	2.47	0.48
1:D:37:GLN:HE21	1:D:37:GLN:H	1.62	0.48
1:A:371:TYR:OH	1:B:118:LYS:HA	2.13	0.48
1:A:67:ARG:NH1	1:A:68:ASP:CG	2.67	0.48
1:B:351:ARG:CD	1:B:351:ARG:N	2.75	0.48
1:C:70:VAL:O	1:C:74:MET:HG3	2.14	0.48
1:D:298:ARG:HD2	1:D:298:ARG:O	2.14	0.47
1:C:122:GLN:HE22	1:C:143:ARG:HD3	1.79	0.47
1:C:73:MET:SD	1:C:90:ILE:CG2	3.02	0.47
1:D:251:VAL:HG11	1:D:346:VAL:CG2	2.44	0.47
1:C:262:VAL:HB	1:C:264:MET:CE	2.44	0.47
1:D:203:THR:HB	1:D:250:GLY:HA2	1.96	0.47
1:D:171:PRO:O	1:D:186:GLY:HA3	2.14	0.47
1:B:69:PHE:CE1	1:B:94:LEU:HD23	2.48	0.47
1:C:21:LYS:HD2	1:C:26:ALA:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:HIS:HD2	1:A:101:ARG:HH12	1.62	0.47
1:D:215:TYR:CZ	1:D:340:LYS:HD3	2.49	0.47
1:D:41:VAL:HG23	1:D:42:GLU:N	2.29	0.47
1:D:64:GLU:OE1	1:D:98:HIS:HE1	1.97	0.47
1:D:251:VAL:CG1	1:D:252:GLY:N	2.78	0.47
1:B:339:LEU:HD12	1:B:339:LEU:N	2.29	0.47
1:C:32:THR:O	1:C:58:ALA:HA	2.14	0.47
1:D:18:LYS:CG	1:D:162:ILE:HD11	2.42	0.47
1:D:34:ASP:HB2	1:D:41:VAL:HG11	1.95	0.47
1:C:40:GLU:O	1:C:40:GLU:HG3	2.15	0.47
1:B:34:ASP:HB3	1:B:60:ASP:HA	1.97	0.47
1:A:67:ARG:NH1	1:A:67:ARG:HG2	2.22	0.47
1:C:120:ASN:HD21	1:C:260:ARG:HG2	1.80	0.47
1:A:97:LYS:HD2	1:A:315:TYR:OH	2.15	0.47
1:C:33:ALA:HB1	1:C:61:LEU:HD12	1.97	0.47
1:C:392:LEU:O	1:C:395:ARG:HA	2.15	0.47
1:B:101:ARG:HD3	6:B:1166:HOH:O	2.14	0.47
1:B:251:VAL:HG11	1:B:346:VAL:CG2	2.45	0.47
1:A:21:LYS:O	1:A:25:ARG:HA	2.15	0.46
1:A:118:LYS:HD3	1:B:365:PHE:CE2	2.51	0.46
1:C:100:VAL:O	1:C:104:GLU:HG3	2.15	0.46
1:B:16:ILE:HG23	1:B:141:PRO:HG2	1.97	0.46
1:A:154:ILE:HG13	1:A:164:VAL:HG11	1.97	0.46
1:C:314:TRP:O	1:C:319:ARG:HD3	2.15	0.46
6:A:1195:HOH:O	1:B:191:PRO:HG2	2.14	0.46
1:A:343:LYS:HD2	1:B:343:LYS:HD3	1.98	0.46
1:A:371:TYR:CE1	1:D:382:ILE:HD11	2.50	0.46
1:C:37:GLN:N	1:C:37:GLN:HE21	2.09	0.46
1:C:18:LYS:CD	1:C:160:HIS:HB3	2.45	0.46
1:B:254:VAL:CG1	1:B:256:ILE:HG23	2.46	0.46
1:D:262:VAL:HG13	1:D:262:VAL:O	2.13	0.46
1:C:16:ILE:HG23	1:C:141:PRO:HG2	1.97	0.46
1:C:18:LYS:HD2	1:C:160:HIS:CB	2.45	0.46
1:C:101:ARG:NH2	6:C:1285:HOH:O	2.41	0.46
1:C:360:GLN:HG3	1:D:189:GLU:HB3	1.96	0.46
1:A:257:VAL:HB	1:B:287:SER:HB2	1.97	0.46
1:B:321:ALA:HA	1:D:328:HIS:CD2	2.51	0.46
1:C:36:GLY:N	1:C:37:GLN:HE21	2.14	0.46
1:B:393:VAL:C	1:B:395:ARG:N	2.69	0.46
1:C:189:GLU:O	1:C:191:PRO:HD3	2.16	0.46
1:A:181:ILE:HD12	1:A:202:MET:HB2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:TYR:HB3	1:B:84:TYR:HB3	1.98	0.46
1:A:67:ARG:NH1	1:A:67:ARG:CG	2.76	0.46
1:D:319:ARG:NH2	1:D:323:GLN:HE21	2.07	0.46
1:D:98:HIS:HD2	1:D:101:ARG:HH11	1.64	0.46
1:D:252:GLY:O	1:D:271:GLU:HA	2.16	0.45
1:B:182:SER:HB2	6:B:1105:HOH:O	2.16	0.45
1:C:67:ARG:HG2	1:C:68:ASP:OD1	2.17	0.45
1:D:60:ASP:OD2	1:D:62:LYS:HE2	2.16	0.45
1:B:385:LEU:HD22	1:C:371:TYR:HD1	1.80	0.45
1:D:22:GLU:OE2	1:D:160:HIS:NE2	2.40	0.45
1:D:341:LEU:HD22	1:D:346:VAL:HG22	1.98	0.45
1:B:359:ARG:C	1:B:361:ASP:N	2.69	0.45
1:B:122:GLN:NE2	1:B:143:ARG:HH11	2.10	0.45
1:C:179:LEU:O	1:C:272:THR:HG23	2.16	0.45
1:C:373:GLN:HE21	1:C:373:GLN:HB3	1.52	0.45
1:C:40:GLU:O	1:C:42:GLU:N	2.50	0.45
4:A:510:ANP:HNB1	5:A:520:ARG:HH12	1.64	0.45
1:A:349:VAL:HG13	6:B:1076:HOH:O	2.16	0.45
1:D:176:ALA:HB1	1:D:181:ILE:HG12	1.97	0.45
1:A:15:ILE:HG23	1:A:156:TYR:CD2	2.52	0.45
1:C:393:VAL:C	1:C:395:ARG:N	2.67	0.45
1:B:393:VAL:C	1:B:395:ARG:H	2.20	0.45
1:C:27:GLU:CG	1:C:54:SER:OG	2.65	0.45
1:D:84:TYR:CE2	1:D:86:LEU:HA	2.52	0.45
1:B:203:THR:HB	1:B:250:GLY:HA2	1.99	0.45
1:A:157:ALA:HB1	1:A:162:ILE:HB	1.98	0.45
1:B:293:GLU:HB2	1:D:318:GLU:HG3	1.98	0.45
1:C:298:ARG:C	1:C:298:ARG:HD2	2.37	0.45
1:A:154:ILE:O	1:A:158:GLU:HG3	2.17	0.44
1:C:27:GLU:HG3	1:C:54:SER:OG	2.16	0.44
1:A:371:TYR:CD1	1:D:382:ILE:HD11	2.52	0.44
1:B:359:ARG:HH21	1:B:362:LEU:CD2	2.24	0.44
1:C:260:ARG:HD2	1:C:264:MET:SD	2.58	0.44
1:C:60:ASP:OD1	1:C:62:LYS:CE	2.65	0.44
1:D:185:GLY:O	1:D:188:LEU:HB2	2.18	0.44
1:C:80:TYR:HB3	1:C:84:TYR:HB3	1.99	0.44
1:C:25:ARG:CB	1:C:25:ARG:HH11	2.31	0.44
5:A:4520:ARG:HH12	4:B:4510:ANP:HNB1	1.66	0.44
1:A:349:VAL:HG22	1:B:194:GLU:CG	2.44	0.44
1:B:34:ASP:OD1	1:B:62:LYS:HD2	2.18	0.44
1:B:179:LEU:O	1:B:272:THR:HG23	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:392:LEU:O	1:D:395:ARG:HG2	2.18	0.44
1:C:74:MET:HB3	1:C:282:ARG:NH1	2.33	0.44
1:B:21:LYS:HD3	6:B:1417:HOH:O	2.18	0.44
1:A:254:VAL:CG1	1:A:256:ILE:HG23	2.48	0.44
1:C:180:HIS:ND1	1:C:180:HIS:C	2.71	0.44
1:C:395:ARG:NH1	1:C:395:ARG:HG3	2.31	0.44
1:D:181:ILE:HD12	1:D:202:MET:HB2	2.00	0.44
1:C:87:GLY:HA3	1:C:180:HIS:CD2	2.53	0.44
1:B:17:LEU:HD11	1:B:53:ALA:HB2	2.00	0.44
1:C:77:GLY:HA2	6:C:1193:HOH:O	2.17	0.44
1:B:181:ILE:CD1	1:B:202:MET:HB2	2.48	0.43
5:A:3520:ARG:NH2	1:D:173:SER:C	2.64	0.43
1:C:90:ILE:HD11	1:C:179:LEU:HD12	2.01	0.43
1:A:262:VAL:CG2	1:A:264:MET:HE1	2.33	0.43
1:B:174:MET:HG2	1:B:183:TYR:CD2	2.53	0.43
1:D:60:ASP:OD1	1:D:62:LYS:HD3	2.17	0.43
1:D:6:ALA:HB3	4:D:3510:ANP:O2'	2.18	0.43
1:D:151:LYS:N	1:D:151:LYS:CD	2.60	0.43
1:C:17:LEU:C	1:C:17:LEU:HD23	2.38	0.43
1:C:71:PHE:HB2	1:C:72:PRO:HD3	2.00	0.43
1:C:123:VAL:HG12	1:C:127:LEU:CD1	2.48	0.43
1:A:191:PRO:HB2	1:B:351:ARG:HD3	2.01	0.43
1:A:12:ASP:OD1	1:A:150:ARG:NH2	2.47	0.43
5:A:2520:ARG:HG2	1:C:182:SER:OG	2.19	0.43
1:B:98:HIS:CD2	1:B:101:ARG:NH1	2.86	0.43
5:A:2520:ARG:HH22	1:C:174:MET:N	2.17	0.43
1:C:18:LYS:HZ3	1:C:160:HIS:HB3	1.84	0.43
1:A:241:ARG:O	1:A:245:ILE:HG13	2.19	0.43
1:B:75:ARG:HD2	1:B:327:ASP:OD2	2.19	0.42
1:D:181:ILE:HG22	1:D:183:TYR:CE1	2.54	0.42
1:A:209:ALA:HB3	1:A:343:LYS:HG2	2.01	0.42
1:D:378:GLY:O	1:D:382:ILE:HG12	2.19	0.42
1:C:282:ARG:NH2	6:C:1320:HOH:O	2.52	0.42
1:A:92:ARG:HE	1:A:121:ASP:CG	2.22	0.42
1:A:64:GLU:OE1	1:A:98:HIS:HE1	2.01	0.42
1:B:251:VAL:CG1	1:B:252:GLY:N	2.82	0.42
1:D:199:MET:O	1:D:201:ARG:NH1	2.52	0.42
1:C:205:ASP:O	1:C:206:PRO:C	2.58	0.42
1:C:151:LYS:HD2	1:C:151:LYS:N	2.09	0.42
1:C:98:HIS:CD2	1:C:101:ARG:HH11	2.36	0.42
1:C:39:GLU:O	1:C:41:VAL:N	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:157:ALA:HB1	1:D:162:ILE:HB	2.01	0.42
1:A:87:GLY:HA3	1:A:180:HIS:CD2	2.55	0.42
1:A:393:VAL:C	1:A:395:ARG:N	2.73	0.42
1:A:117:GLY:C	1:A:122:GLN:HE22	2.22	0.42
1:C:352:LYS:HE3	1:D:192:TRP:CH2	2.54	0.42
1:A:176:ALA:HB1	1:A:181:ILE:HG12	2.02	0.42
1:A:379:PHE:HA	1:D:379:PHE:HA	2.01	0.42
1:A:288:LEU:HD11	1:A:351:ARG:HG2	2.01	0.42
1:D:302:SER:HB2	1:D:303:PRO:HD3	2.01	0.42
1:B:304:LYS:O	1:B:307:GLU:HG2	2.18	0.42
1:B:395:ARG:HG3	1:B:395:ARG:NH1	2.34	0.42
1:C:12:ASP:OD1	1:C:150:ARG:NH2	2.51	0.42
1:D:74:MET:HB3	1:D:282:ARG:CZ	2.49	0.42
1:D:40:GLU:OE2	1:D:40:GLU:HA	2.19	0.42
1:C:4:VAL:O	1:C:111:ILE:HA	2.20	0.42
1:C:12:ASP:O	1:C:16:ILE:HG13	2.19	0.41
1:A:339:LEU:N	1:A:339:LEU:HD12	2.35	0.41
1:D:179:LEU:O	1:D:272:THR:HG23	2.20	0.41
1:B:123:VAL:HG12	1:B:127:LEU:HD12	2.01	0.41
1:A:263:GLY:HA3	1:C:385:LEU:HD23	2.03	0.41
1:D:174:MET:HA	1:D:182:SER:O	2.20	0.41
1:B:319:ARG:HH21	1:B:323:GLN:NE2	2.18	0.41
1:B:128:THR:O	1:B:132:LEU:HG	2.20	0.41
1:D:394:GLU:O	1:D:395:ARG:C	2.58	0.41
1:C:393:VAL:C	1:C:395:ARG:H	2.21	0.41
1:A:18:LYS:HE3	1:A:160:HIS:HB3	2.03	0.41
1:D:395:ARG:NH1	1:D:395:ARG:HG3	2.35	0.41
1:A:180:HIS:HB2	1:A:271:GLU:O	2.21	0.41
1:D:84:TYR:CZ	1:D:86:LEU:HA	2.55	0.41
1:A:328:HIS:CD2	1:C:321:ALA:HA	2.56	0.41
1:A:184:GLU:OE2	5:A:520:ARG:HG3	2.21	0.41
1:A:316:ALA:HB1	1:A:318:GLU:OE2	2.20	0.41
1:A:66:VAL:HA	1:A:70:VAL:HB	2.01	0.41
1:C:174:MET:HA	1:C:182:SER:O	2.21	0.41
1:A:318:GLU:HG3	1:C:293:GLU:HB2	2.03	0.41
1:B:180:HIS:HB2	1:B:271:GLU:O	2.20	0.41
1:D:394:GLU:O	1:D:395:ARG:O	2.38	0.41
5:A:3530:ASP:N	4:D:3510:ANP:O2B	2.54	0.41
1:A:150:ARG:HG3	1:A:154:ILE:CD1	2.49	0.41
1:D:77:GLY:HA2	6:D:1450:HOH:O	2.21	0.41
1:D:135:ASP:OD2	1:D:135:ASP:N	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:71:PHE:HB2	1:D:72:PRO:HD3	2.03	0.41
1:C:154:ILE:O	1:C:158:GLU:HG3	2.21	0.41
1:C:157:ALA:O	1:C:158:GLU:C	2.59	0.41
1:C:7:TYR:CE2	1:C:9:GLY:HA2	2.56	0.41
1:C:73:MET:CE	1:C:90:ILE:HG12	2.52	0.40
1:A:254:VAL:HG13	1:A:256:ILE:HG23	2.02	0.40
1:A:245:ILE:O	1:A:248:ARG:CG	2.69	0.40
1:A:176:ALA:CB	1:A:181:ILE:HG12	2.51	0.40
1:B:92:ARG:HE	1:B:121:ASP:CG	2.25	0.40
1:D:146:SER:O	1:D:147:PHE:C	2.59	0.40
1:D:39:GLU:OE1	1:D:39:GLU:HA	2.22	0.40
1:B:66:VAL:HA	1:B:70:VAL:HB	2.02	0.40
1:A:191:PRO:O	1:B:350:GLY:HA2	2.21	0.40
1:D:395:ARG:NH1	1:D:395:ARG:CG	2.82	0.40
1:A:80:TYR:CD2	1:A:81:GLU:HG3	2.56	0.40
1:A:143:ARG:HD2	1:B:373:GLN:O	2.22	0.40
1:C:281:ALA:O	1:C:284:ALA:HB3	2.22	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	374/400 (94%)	362 (97%)	11 (3%)	1 (0%)	46	42
1	B	380/400 (95%)	366 (96%)	13 (3%)	1 (0%)	46	42
1	C	380/400 (95%)	356 (94%)	20 (5%)	4 (1%)	17	10
1	D	380/400 (95%)	367 (97%)	12 (3%)	1 (0%)	46	42
All	All	1514/1600 (95%)	1451 (96%)	56 (4%)	7 (0%)	34	26

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	159	ALA
1	C	160	HIS
1	C	22	GLU
1	C	41	VAL
1	A	343	LYS
1	B	373	GLN
1	D	273	PRO

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	304/319 (95%)	283 (93%)	21 (7%)	19	13
1	B	310/319 (97%)	294 (95%)	16 (5%)	29	23
1	C	310/319 (97%)	290 (94%)	20 (6%)	21	15
1	D	310/319 (97%)	288 (93%)	22 (7%)	18	12
All	All	1234/1276 (97%)	1155 (94%)	79 (6%)	22	15

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	GLN
1	A	90	ILE
1	A	94	LEU
1	A	128	THR
1	A	137	LYS
1	A	145	TRP
1	A	151	LYS
1	A	178	LEU
1	A	179	LEU
1	A	180	HIS
1	A	188	LEU
1	A	239	LEU
1	A	264	MET
1	A	267	ARG
1	A	291	ASP

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Mol	Chain	Res	Type
1	A	318	GLU
1	A	319	ARG
1	A	349	VAL
1	A	351	ARG
1	A	371	TYR
1	A	373	GLN
1	B	37	GLN
1	B	39	GLU
1	B	62	LYS
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	267	ARG
1	B	319	ARG
1	B	351	ARG
1	B	361	ASP
1	B	373	GLN
1	C	25	ARG
1	C	37	GLN
1	C	46	GLU
1	C	62	LYS
1	C	145	TRP
1	C	150	ARG
1	C	179	LEU
1	C	180	HIS
1	C	188	LEU
1	C	206	PRO
1	C	239	LEU
1	C	264	MET
1	C	267	ARG
1	C	318	GLU
1	C	319	ARG
1	C	345	ASN
1	C	351	ARG
1	C	361	ASP
1	C	373	GLN
1	C	374	LYS

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Mol	Chain	Res	Type
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	151	LYS
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	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 (28) 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
1	B	37	GLN
1	B	98	HIS
1	B	122	GLN
1	B	148	GLN
1	B	249	HIS
1	B	323	GLN
1	B	345	ASN
1	B	360	GLN
1	C	37	GLN
1	C	98	HIS

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Mol	Chain	Res	Type
1	C	122	GLN
1	C	323	GLN
1	C	360	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	297	GLN
1	D	323	GLN
1	D	360	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](#)

Of 20 ligands modelled in this entry, 4 are monoatomic - leaving 16 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
5	ARG	A	2520	-	5,11,11	0.78	0	3,13,13	1.24	1 (33%)
5	ASP	A	2530	-	2,8,8	0.20	0	0,10,10	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	ARG	A	3520	-	5,11,11	0.58	0	3,13,13	1.21	0
5	ASP	A	3530	-	2,8,8	0.42	0	0,10,10	0.00	-
5	ARG	A	4520	-	5,11,11	0.77	0	3,13,13	1.65	1 (33%)
5	ASP	A	4530	-	2,8,8	0.27	0	0,10,10	0.00	-
4	ANP	A	510	3	27,33,33	2.63	8 (29%)	30,52,52	2.58	9 (30%)
5	ARG	A	520	-	5,11,11	0.66	0	3,13,13	1.38	1 (33%)
5	ASP	A	530	-	2,8,8	0.30	0	0,10,10	0.00	-
2	SO4	A	540	-	4,4,4	3.19	2 (50%)	6,6,6	0.98	0
4	ANP	B	4510	3	27,33,33	2.40	10 (37%)	30,52,52	2.48	9 (30%)
2	SO4	B	4540	-	4,4,4	3.24	2 (50%)	6,6,6	1.04	0
4	ANP	C	2510	3	27,33,33	2.44	9 (33%)	30,52,52	2.64	9 (30%)
2	SO4	C	2540	-	4,4,4	3.23	2 (50%)	6,6,6	1.01	0
4	ANP	D	3510	3	27,33,33	2.59	9 (33%)	30,52,52	2.64	9 (30%)
2	SO4	D	3540	-	4,4,4	3.27	2 (50%)	6,6,6	0.98	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	ARG	A	2520	-	-	0/5/11/11	0/0/0/0
5	ASP	A	2530	-	-	0/2/8/8	0/0/0/0
5	ARG	A	3520	-	-	0/5/11/11	0/0/0/0
5	ASP	A	3530	-	-	0/2/8/8	0/0/0/0
5	ARG	A	4520	-	-	0/5/11/11	0/0/0/0
5	ASP	A	4530	-	-	0/2/8/8	0/0/0/0
4	ANP	A	510	3	-	0/12/38/38	0/3/3/3
5	ARG	A	520	-	-	0/5/11/11	0/0/0/0
5	ASP	A	530	-	-	0/2/8/8	0/0/0/0
2	SO4	A	540	-	-	0/0/0/0	0/0/0/0
4	ANP	B	4510	3	-	0/12/38/38	0/3/3/3
2	SO4	B	4540	-	-	0/0/0/0	0/0/0/0
4	ANP	C	2510	3	-	1/12/38/38	0/3/3/3
2	SO4	C	2540	-	-	0/0/0/0	0/0/0/0
4	ANP	D	3510	3	-	0/12/38/38	0/3/3/3
2	SO4	D	3540	-	-	0/0/0/0	0/0/0/0

All (44) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	3540	SO4	O3-S	-4.48	1.31	1.47
2	C	2540	SO4	O3-S	-4.47	1.31	1.47
2	B	4540	SO4	O3-S	-4.44	1.31	1.47
2	A	540	SO4	O3-S	-4.35	1.31	1.47
4	A	510	ANP	PB-O2B	-4.24	1.44	1.56
4	C	2510	ANP	PB-O2B	-3.92	1.45	1.56
4	B	4510	ANP	PB-O2B	-3.87	1.45	1.56
4	D	3510	ANP	PB-O2B	-3.38	1.47	1.56
4	C	2510	ANP	PG-O3G	-2.20	1.50	1.56
4	B	4510	ANP	PG-O2G	-2.20	1.50	1.56
4	B	4510	ANP	PG-O3G	-2.11	1.50	1.56
4	D	3510	ANP	PG-O3G	-2.04	1.51	1.56
4	B	4510	ANP	C2-N1	2.13	1.37	1.33
4	C	2510	ANP	C2-N1	2.21	1.38	1.33
4	B	4510	ANP	C2-N3	2.34	1.36	1.32
4	A	510	ANP	C2-N1	2.44	1.38	1.33
4	D	3510	ANP	C2-N1	2.45	1.38	1.33
4	C	2510	ANP	C2-N3	2.54	1.36	1.32
4	D	3510	ANP	O4'-C1'	2.66	1.44	1.41
4	D	3510	ANP	C2-N3	2.72	1.37	1.32
4	A	510	ANP	C2-N3	2.73	1.37	1.32
4	C	2510	ANP	O4'-C1'	2.78	1.44	1.41
4	B	4510	ANP	O4'-C1'	2.95	1.44	1.41
4	A	510	ANP	O4'-C1'	3.30	1.45	1.41
4	A	510	ANP	PG-N3B	3.87	1.73	1.63
4	C	2510	ANP	PG-N3B	3.87	1.73	1.63
4	C	2510	ANP	PB-O1B	4.05	1.50	1.46
4	B	4510	ANP	PG-N3B	4.07	1.74	1.63
4	B	4510	ANP	PB-O1B	4.39	1.51	1.46
2	C	2540	SO4	O1-S	4.52	1.62	1.47
2	A	540	SO4	O1-S	4.53	1.62	1.47
2	B	4540	SO4	O1-S	4.59	1.62	1.47
4	C	2510	ANP	PG-O1G	4.64	1.51	1.46
2	D	3540	SO4	O1-S	4.66	1.63	1.47
4	D	3510	ANP	PG-N3B	4.68	1.75	1.63
4	A	510	ANP	PG-O1G	4.68	1.51	1.46
4	D	3510	ANP	PG-O1G	5.28	1.52	1.46
4	A	510	ANP	PB-O1B	5.41	1.52	1.46
4	B	4510	ANP	PB-O3A	5.46	1.65	1.59
4	B	4510	ANP	PG-O1G	5.57	1.52	1.46
4	D	3510	ANP	PB-O1B	5.63	1.52	1.46
4	D	3510	ANP	PB-O3A	6.78	1.67	1.59
4	C	2510	ANP	PB-O3A	7.05	1.67	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	510	ANP	PB-O3A	7.37	1.68	1.59

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	2510	ANP	N3-C2-N1	-6.92	123.60	128.89
4	A	510	ANP	N3-C2-N1	-6.76	123.72	128.89
4	D	3510	ANP	N3-C2-N1	-6.69	123.77	128.89
4	B	4510	ANP	N3-C2-N1	-6.53	123.89	128.89
4	A	510	ANP	O1B-PB-N3B	-5.31	103.76	111.90
4	D	3510	ANP	O1B-PB-N3B	-5.24	103.86	111.90
4	C	2510	ANP	O1B-PB-N3B	-5.22	103.90	111.90
4	D	3510	ANP	O1G-PG-N3B	-4.89	104.40	111.90
4	B	4510	ANP	O1B-PB-N3B	-4.72	104.65	111.90
4	C	2510	ANP	O1G-PG-N3B	-4.65	104.77	111.90
4	A	510	ANP	O1G-PG-N3B	-3.90	105.92	111.90
4	B	4510	ANP	O1G-PG-N3B	-3.46	106.60	111.90
4	A	510	ANP	O5'-C5'-C4'	-2.86	98.59	109.12
4	D	3510	ANP	O4'-C4'-C3'	-2.85	99.41	105.15
4	D	3510	ANP	O5'-C5'-C4'	-2.66	99.31	109.12
5	A	4520	ARG	CG-CD-NE	-2.64	103.21	112.35
4	B	4510	ANP	O5'-C5'-C4'	-2.63	99.43	109.12
4	A	510	ANP	O4'-C4'-C3'	-2.50	100.10	105.15
4	C	2510	ANP	O4'-C4'-C3'	-2.47	100.17	105.15
4	B	4510	ANP	O4'-C4'-C3'	-2.43	100.25	105.15
4	C	2510	ANP	O5'-C5'-C4'	-2.34	100.49	109.12
5	A	520	ARG	CG-CD-NE	-2.16	104.86	112.35
5	A	2520	ARG	CG-CD-NE	-2.06	105.21	112.35
4	C	2510	ANP	O3A-PB-N3B	3.05	114.82	106.44
4	D	3510	ANP	O4'-C1'-N9	3.21	114.81	108.10
4	B	4510	ANP	O3A-PB-N3B	3.22	115.29	106.44
4	B	4510	ANP	O4'-C1'-N9	3.22	114.85	108.10
4	C	2510	ANP	O4'-C1'-N9	3.25	114.90	108.10
4	A	510	ANP	O4'-C1'-N9	3.28	114.96	108.10
4	A	510	ANP	O3A-PB-N3B	3.29	115.50	106.44
4	B	4510	ANP	O3A-PA-O5'	3.32	111.74	102.94
4	D	3510	ANP	O3A-PB-N3B	3.40	115.80	106.44
4	D	3510	ANP	O3A-PA-O5'	3.58	112.44	102.94
4	C	2510	ANP	O3A-PA-O5'	3.76	112.91	102.94
4	A	510	ANP	O3A-PA-O5'	4.18	114.03	102.94
4	A	510	ANP	C4'-O4'-C1'	6.59	116.96	109.72
4	D	3510	ANP	C4'-O4'-C1'	7.12	117.54	109.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	4510	ANP	C4'-O4'-C1'	7.14	117.56	109.72
4	C	2510	ANP	C4'-O4'-C1'	7.22	117.65	109.72

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	2510	ANP	O1G-PG-N3B-PB

There are no ring outliers.

8 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	2520	ARG	4	0
5	A	3520	ARG	5	0
5	A	3530	ASP	1	0
5	A	4520	ARG	3	0
4	A	510	ANP	1	0
5	A	520	ARG	4	0
4	B	4510	ANP	1	0
4	D	3510	ANP	2	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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%)	-0.05	8 (2%) 67 74	20, 28, 52, 76	0
1	B	386/400 (96%)	-0.09	14 (3%) 46 57	19, 28, 49, 67	0
1	C	386/400 (96%)	0.15	25 (6%) 22 31	21, 33, 60, 84	0
1	D	386/400 (96%)	0.16	24 (6%) 24 33	22, 33, 55, 77	0
All	All	1538/1600 (96%)	0.04	71 (4%) 36 47	19, 30, 55, 84	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	370	GLY	9.3
1	C	165	PRO	8.0
1	A	371	TYR	6.7
1	D	370	GLY	5.6
1	A	165	PRO	5.6
1	B	365	PHE	5.0
1	D	159	ALA	4.8
1	A	164	VAL	4.5
1	D	165	PRO	4.3
1	B	165	PRO	4.1
1	B	371	TYR	4.0
1	B	363	VAL	3.8
1	C	25	ARG	3.8
1	A	38	GLY	3.8
1	D	365	PHE	3.6
1	B	370	GLY	3.6
1	C	373	GLN	3.6
1	D	154	ILE	3.5
1	C	370	GLY	3.5
1	D	371	TYR	3.5
1	B	364	SER	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	148	GLN	3.4
1	C	39	GLU	3.4
1	B	373	GLN	3.4
1	D	40	GLU	3.2
1	D	39	GLU	3.2
1	C	365	PHE	3.0
1	C	40	GLU	3.0
1	B	360	GLN	3.0
1	D	38	GLY	3.0
1	C	371	TYR	2.9
1	C	395	ARG	2.8
1	D	364	SER	2.8
1	B	362	LEU	2.8
1	D	164	VAL	2.8
1	C	364	SER	2.7
1	D	201	ARG	2.7
1	C	49	LEU	2.7
1	B	395	ARG	2.7
1	C	30	ALA	2.7
1	C	43	GLU	2.7
1	C	164	VAL	2.6
1	D	161	GLY	2.5
1	C	160	HIS	2.5
1	C	232	ARG	2.5
1	B	361	ASP	2.4
1	D	221	PHE	2.4
1	D	150	ARG	2.4
1	D	232	ARG	2.4
1	A	161	GLY	2.4
1	D	363	VAL	2.4
1	C	361	ASP	2.3
1	D	187	VAL	2.3
1	C	5	LEU	2.3
1	C	4	VAL	2.3
1	C	159	ALA	2.3
1	C	204	GLN	2.2
1	D	148	GLN	2.2
1	B	40	GLU	2.2
1	B	221	PHE	2.2
1	D	5	LEU	2.1
1	C	331	ARG	2.1
1	C	161	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	160	HIS	2.1
1	C	162	ILE	2.1
1	A	197	LYS	2.1
1	C	154	ILE	2.0
1	B	30	ALA	2.0
1	D	43	GLU	2.0
1	D	50	ARG	2.0
1	D	395	ARG	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
5	ARG	A	3520	12/12	0.78	0.29	6.95	28,33,53,58	0
5	ARG	A	520	12/12	0.81	0.26	6.67	19,28,49,52	0
5	ARG	A	2520	12/12	0.79	0.27	6.30	24,38,55,56	0
5	ARG	A	4520	12/12	0.83	0.22	6.07	20,29,45,49	0
2	SO4	A	540	5/5	0.88	0.21	3.86	48,61,63,64	5
2	SO4	D	3540	5/5	0.85	0.21	3.15	61,74,78,81	0
2	SO4	C	2540	5/5	0.73	0.26	2.96	75,81,85,86	5
2	SO4	B	4540	5/5	0.95	0.13	1.35	32,51,58,60	0
4	ANP	B	4510	31/31	0.92	0.13	0.36	20,31,40,41	0
4	ANP	D	3510	31/31	0.86	0.14	0.28	33,41,60,62	0
4	ANP	A	510	31/31	0.92	0.13	0.14	25,32,41,44	0
4	ANP	C	2510	31/31	0.89	0.13	-0.08	25,35,47,55	0
5	ASP	A	2530	9/9	0.94	0.11	-0.15	20,25,30,30	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
5	ASP	A	3530	9/9	0.95	0.10	-0.38	26,28,32,34	0
5	ASP	A	530	9/9	0.97	0.09	-0.47	24,29,30,38	0
5	ASP	A	4530	9/9	0.97	0.07	-0.60	25,27,29,29	0
3	MG	D	3550	1/1	0.82	0.12	-	55,55,55,55	0
3	MG	A	550	1/1	0.80	0.09	-	42,42,42,42	0
3	MG	C	2550	1/1	0.89	0.08	-	43,43,43,43	0
3	MG	B	4550	1/1	0.92	0.05	-	38,38,38,38	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.