



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 11:37 AM GMT

PDB ID : 1J20
Title : Crystal Structure of Thermus thermophilus HB8 Argininosuccinate Synthetase in complex with product
Authors : Goto, M.; Hirotsu, K.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2002-12-24
Resolution : 2.00 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

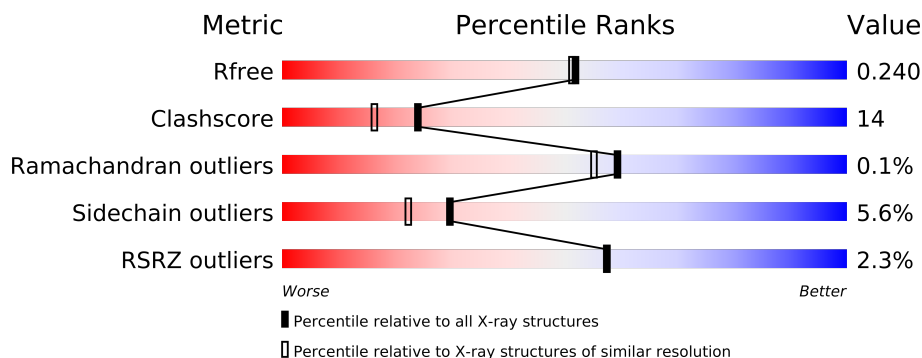
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.00 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}	66092	4888 (2.00-2.00)
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (2.00-2.00)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	400	
1	B	400	
1	C	400	
1	D	400	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	AMP	A	510	-	X
3	AMP	B	2510	-	X
3	AMP	C	3510	-	X
3	AMP	D	4510	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
4	AS1	A	520	-	X
4	AS1	B	2520	-	X
4	AS1	C	3520	-	X
4	AS1	D	4520	-	X

2 Entry composition

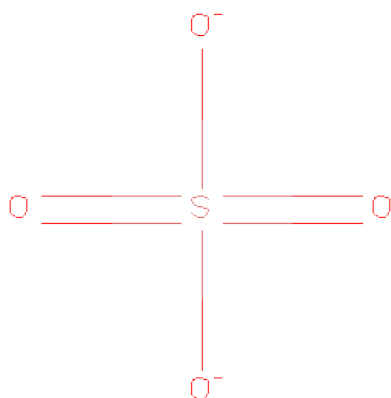
There are 5 unique types of molecules in this entry. The entry contains 13318 atoms, of which 0 are hydrogen and 0 are deuterium.

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	386	Total	C	N	O	S	0	0	0
			3066	1963	533	561	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: SO4) (formula: O₄S).



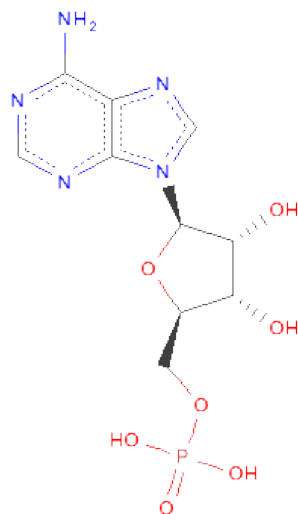
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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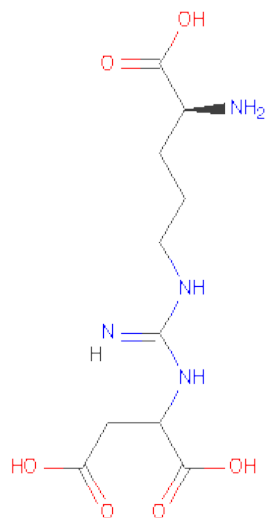
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula: $C_{10}H_{14}N_5O_7P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
3	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
3	C	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
3	D	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 4 is ARGININOSUCCINATE (three-letter code: AS1) (formula: $C_{10}H_{18}N_4O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			20	10	4	6		
4	B	1	Total	C	N	O	0	0
			20	10	4	6		
4	C	1	Total	C	N	O	0	0
			20	10	4	6		
4	D	1	Total	C	N	O	0	0
			20	10	4	6		

- Molecule 5 is water.

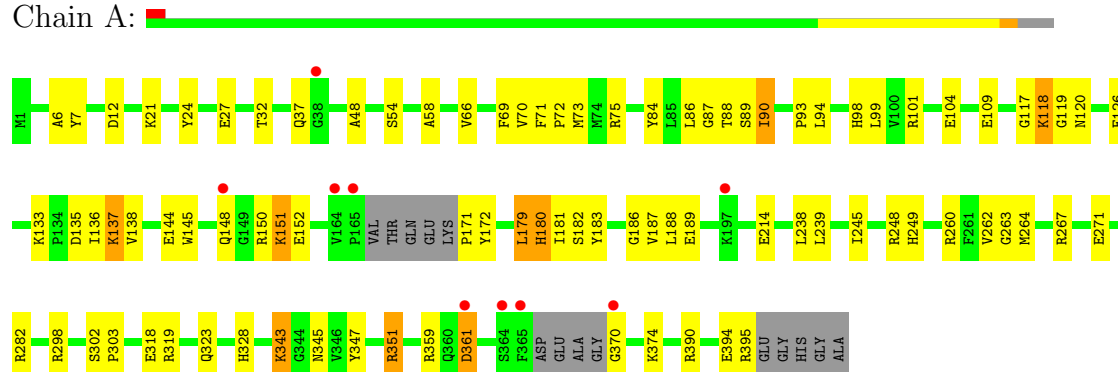
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	230	Total	O	0	0
			230	230		
5	B	238	Total	O	0	0
			238	238		
5	C	194	Total	O	0	0
			194	194		
5	D	200	Total	O	0	0
			200	200		

3 Residue-property plots

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.

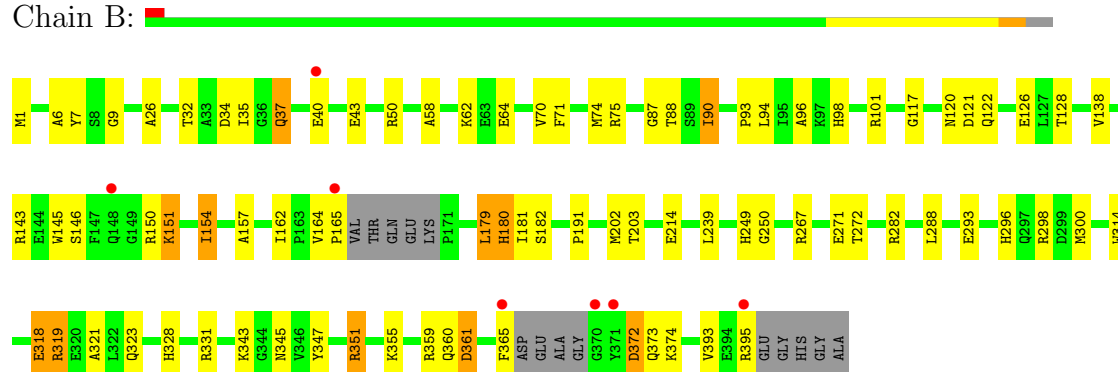
• Molecule 1: Argininosuccinate Synthetase

Chain A:



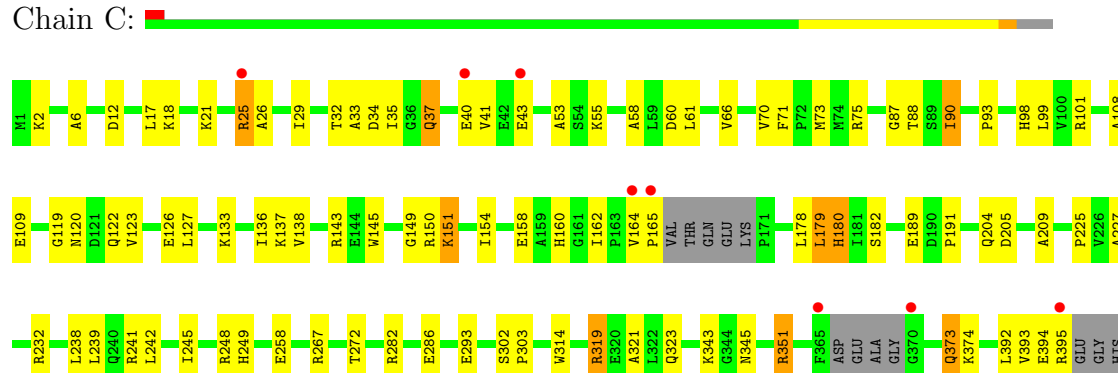
• Molecule 1: Argininosuccinate Synthetase

Chain B:



• Molecule 1: Argininosuccinate Synthetase

Chain C:

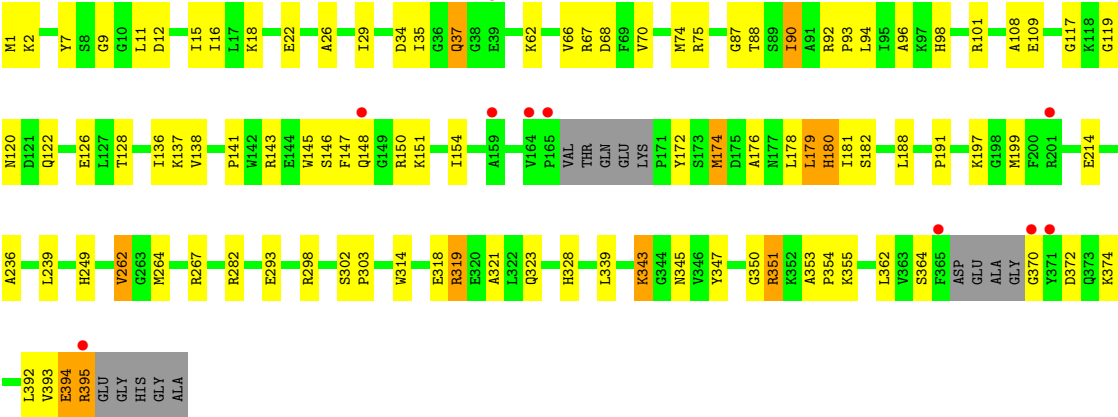


GLY

ALA

● Molecule 1: Argininosuccinate Synthetase

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	228.78Å 228.78Å 159.41Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.00 19.96 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.00) 99.5 (19.96-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 2.01Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.212 , 0.239 0.213 , 0.240	Depositor DCC
R_{free} test set	20669 reflections (9.95%)	DCC
Wilson B-factor (Å ²)	28.7	Xtriage
Anisotropy	0.003	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 48.6	EDS
Estimated twinning fraction	0.003 for h,-h-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Outliers	0 of 208941 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13318	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: AMP, AS1, 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	0.40	0/3135	0.66	1/4237 (0.0%)
1	B	0.39	0/3135	0.67	0/4237
1	C	0.41	0/3135	0.67	0/4237
1	D	0.40	0/3135	0.67	0/4237
All	All	0.40	0/12540	0.67	1/16948 (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	A	263	GLY	N-CA-C	5.16	126.00	113.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3066	0	3052	78	0
1	B	3066	0	3052	95	0
1	C	3066	0	3052	103	0
1	D	3066	0	3052	96	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
3	A	23	0	12	2	0
3	B	23	0	12	3	0
3	C	23	0	12	1	0
3	D	23	0	12	2	0
4	A	20	0	13	8	0
4	B	20	0	13	6	0
4	C	20	0	13	9	0
4	D	20	0	11	8	0
5	A	230	0	0	12	0
5	B	238	0	0	8	0
5	C	194	0	0	7	0
5	D	200	0	0	5	0
All	All	13318	0	12306	352	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 14.

All (352) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:A:520:AS1:CB	4:A:520:AS1:CA	1.80	1.59
4:C:3520:AS1:CB	4:C:3520:AS1:CA	1.80	1.56
4:B:2520:AS1:CB	4:B:2520:AS1:CA	1.80	1.55
4:D:4520:AS1:CB	4:D:4520:AS1:CA	1.81	1.54
1:A:151:LYS:H	1:A:151:LYS:HD2	0.99	1.11
1:B:151:LYS:HD2	1:B:151:LYS:H	1.18	1.09
1:D:151:LYS:H	1:D:151:LYS:HD2	1.12	1.07
1:B:98:HIS:HD2	1:B:101:ARG:HH21	1.05	1.03
1:B:35:ILE:H	1:B:37:GLN:NE2	1.60	0.99
1:C:154:ILE:HG23	1:C:164:VAL:HG11	1.44	0.98
1:A:151:LYS:CD	1:A:151:LYS:H	1.79	0.94
1:A:151:LYS:N	1:A:151:LYS:HD2	1.81	0.92
1:A:98:HIS:HD2	1:A:101:ARG:HH21	1.15	0.90
1:D:351:ARG:HD2	1:D:351:ARG:N	1.87	0.89
1:C:35:ILE:H	1:C:37:GLN:NE2	1.69	0.88
1:A:359:ARG:NH2	5:A:1286:HOH:O	2.06	0.87
1:B:318:GLU:HG3	1:D:293:GLU:HB2	1.55	0.86
1:D:35:ILE:H	1:D:37:GLN:NE2	1.76	0.84
1:C:37:GLN:H	1:C:37:GLN:NE2	1.75	0.83

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:151:LYS:H	1:D:151:LYS:CD	1.91	0.83
1:B:181:ILE:HG13	1:B:202:MET:HE1	1.61	0.82
1:B:151:LYS:CD	1:B:151:LYS:H	1.93	0.82
1:A:351:ARG:N	1:A:351:ARG:HD2	1.95	0.80
1:C:37:GLN:H	1:C:37:GLN:HE21	1.28	0.80
1:B:98:HIS:CD2	1:B:101:ARG:HH21	1.97	0.79
1:B:126:GLU:HG2	1:B:138:VAL:HG11	1.65	0.78
1:C:18:LYS:HG3	1:C:162:ILE:HD11	1.66	0.78
1:B:318:GLU:HG3	1:D:293:GLU:CB	2.13	0.78
1:C:319:ARG:HH21	1:C:323:GLN:HE21	1.31	0.78
1:A:126:GLU:HG2	1:A:138:VAL:HG11	1.67	0.77
1:D:395:ARG:HH11	1:D:395:ARG:HG3	1.51	0.76
1:C:245:ILE:O	1:C:248:ARG:HG2	1.86	0.76
1:D:37:GLN:NE2	1:D:37:GLN:H	1.84	0.76
1:A:319:ARG:HH11	1:A:323:GLN:HE21	1.30	0.75
1:B:37:GLN:H	1:B:37:GLN:NE2	1.84	0.75
1:C:343:LYS:HD2	1:D:343:LYS:HD2	1.67	0.75
1:D:151:LYS:N	1:D:151:LYS:HD2	1.97	0.75
1:B:293:GLU:HB3	1:D:318:GLU:HG2	1.68	0.74
1:D:98:HIS:HD2	1:D:101:ARG:HH21	1.33	0.74
1:C:164:VAL:HG12	1:C:165:PRO:HD2	1.69	0.73
1:B:314:TRP:O	1:B:319:ARG:HD3	1.88	0.73
1:B:40:GLU:HG3	1:B:43:GLU:HG3	1.72	0.72
1:B:181:ILE:HG13	1:B:202:MET:CE	2.19	0.71
1:C:151:LYS:CD	1:C:151:LYS:H	2.04	0.70
1:B:35:ILE:H	1:B:37:GLN:HE22	1.40	0.70
1:D:109:GLU:O	1:D:137:LYS:HE2	1.91	0.70
4:A:520:AS1:HA	4:A:520:AS1:CB	2.17	0.70
4:C:3520:AS1:CB	4:C:3520:AS1:HA	2.17	0.69
1:C:351:ARG:N	1:C:351:ARG:HD2	2.07	0.69
1:A:318:GLU:CG	1:C:293:GLU:HB2	2.23	0.68
1:D:137:LYS:N	1:D:137:LYS:HD3	2.08	0.68
1:C:109:GLU:O	1:C:137:LYS:HE2	1.93	0.67
1:A:98:HIS:CD2	1:A:101:ARG:HH21	2.06	0.67
1:D:75:ARG:NH2	5:D:1529:HOH:O	2.26	0.67
1:C:40:GLU:CG	1:C:43:GLU:HG3	2.25	0.67
1:D:319:ARG:HH21	1:D:323:GLN:HE21	1.40	0.67
1:D:37:GLN:HE21	1:D:37:GLN:H	1.42	0.67
1:B:319:ARG:HH21	1:B:323:GLN:HE21	1.42	0.67
1:B:40:GLU:CG	1:B:43:GLU:HG3	2.25	0.67
3:B:2510:AMP:H5'1	3:B:2510:AMP:H8	1.61	0.66
1:D:282:ARG:HD2	5:D:1355:HOH:O	1.94	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:245:ILE:O	1:A:248:ARG:HG2	1.96	0.65
1:C:314:TRP:O	1:C:319:ARG:HD3	1.96	0.65
1:C:351:ARG:HD3	1:D:191:PRO:HB2	1.78	0.65
1:B:93:PRO:HG3	5:B:1181:HOH:O	1.97	0.65
5:A:1694:HOH:O	1:B:345:ASN:HB2	1.96	0.65
1:A:318:GLU:HG2	1:C:293:GLU:HB3	1.78	0.64
1:B:293:GLU:CB	1:D:318:GLU:HG2	2.28	0.64
5:A:1272:HOH:O	1:B:191:PRO:HG2	1.98	0.64
1:D:314:TRP:O	1:D:319:ARG:HD3	1.97	0.63
1:B:151:LYS:HD2	1:B:151:LYS:N	2.01	0.63
1:A:318:GLU:HG2	1:C:293:GLU:CB	2.29	0.63
1:C:12:ASP:OD1	1:C:150:ARG:NH2	2.31	0.63
1:B:37:GLN:H	1:B:37:GLN:HE21	1.45	0.63
1:D:395:ARG:NH1	1:D:395:ARG:HG3	2.12	0.63
1:C:35:ILE:H	1:C:37:GLN:HE22	1.43	0.62
1:C:119:GLY:HA3	4:C:3520:AS1:OD1	1.99	0.62
1:D:126:GLU:HG2	1:D:138:VAL:HG11	1.82	0.62
1:D:35:ILE:H	1:D:37:GLN:HE22	1.47	0.62
1:D:197:LYS:HG3	5:D:1447:HOH:O	1.98	0.62
1:C:122:GLN:HE21	1:C:143:ARG:HH11	1.46	0.62
1:B:293:GLU:HB2	1:D:318:GLU:CG	2.30	0.62
1:A:214:GLU:OE1	1:A:249:HIS:HE1	1.83	0.62
1:B:202:MET:HE2	5:B:1479:HOH:O	2.00	0.62
1:A:87:GLY:HA2	1:A:179:LEU:HD13	1.82	0.61
1:D:87:GLY:HA2	1:D:179:LEU:HD13	1.82	0.61
1:B:120:ASN:H	4:B:2520:AS1:CD	2.13	0.61
1:B:122:GLN:HE21	1:B:143:ARG:HH11	1.46	0.61
1:C:395:ARG:HG3	1:C:395:ARG:HH11	1.64	0.61
1:B:35:ILE:N	1:B:37:GLN:NE2	2.40	0.61
1:B:318:GLU:CG	1:D:293:GLU:CB	2.79	0.61
1:D:120:ASN:N	4:D:4520:AS1:OD2	2.35	0.60
1:C:88:THR:HG21	4:C:3520:AS1:H12	1.84	0.60
1:A:119:GLY:HA3	4:A:520:AS1:OD2	2.02	0.60
3:B:2510:AMP:C8	3:B:2510:AMP:H5'1	2.38	0.59
1:D:29:ILE:HD12	1:D:29:ILE:N	2.17	0.59
1:B:150:ARG:O	1:B:154:ILE:HD12	2.02	0.59
4:D:4520:AS1:HA	4:D:4520:AS1:CB	2.19	0.59
1:B:282:ARG:NH2	5:B:1160:HOH:O	2.36	0.59
1:C:248:ARG:HG3	1:C:249:HIS:CD2	2.37	0.59
1:C:35:ILE:N	1:C:37:GLN:NE2	2.48	0.59
1:D:214:GLU:OE1	1:D:249:HIS:HE1	1.86	0.59
1:B:293:GLU:CB	1:D:318:GLU:CG	2.81	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:318:GLU:HG3	1:C:293:GLU:HB2	1.84	0.58
1:A:374:LYS:NZ	1:D:374:LYS:HE3	2.18	0.58
1:C:98:HIS:HD2	1:C:101:ARG:NH1	2.02	0.58
1:A:27:GLU:HG2	1:A:54:SER:OG	2.02	0.58
5:C:1151:HOH:O	1:D:191:PRO:HG2	2.04	0.57
1:D:1:MET:O	1:D:26:ALA:HB1	2.04	0.57
1:C:151:LYS:HD2	1:C:151:LYS:H	1.68	0.57
1:C:41:VAL:HG12	1:C:58:ALA:HB1	1.85	0.57
1:A:88:THR:HG21	4:A:520:AS1:H12	1.87	0.57
1:D:394:GLU:O	1:D:395:ARG:C	2.43	0.57
1:B:88:THR:HG21	4:B:2520:AS1:H12	1.86	0.57
1:D:67:ARG:HD2	1:D:68:ASP:OD1	2.04	0.57
1:A:109:GLU:O	1:A:137:LYS:HE2	2.05	0.57
1:D:119:GLY:HA3	4:D:4520:AS1:OD2	2.04	0.57
1:C:120:ASN:N	4:C:3520:AS1:OD1	2.38	0.56
1:B:293:GLU:HB2	1:D:318:GLU:HG3	1.87	0.56
1:B:122:GLN:NE2	1:B:143:ARG:HH11	2.03	0.56
1:B:214:GLU:OE1	1:B:249:HIS:HE1	1.88	0.56
1:C:98:HIS:CD2	1:C:101:ARG:NH1	2.74	0.56
1:D:74:MET:HE2	5:D:1833:HOH:O	2.05	0.56
1:C:164:VAL:CG1	1:C:165:PRO:HD2	2.34	0.56
1:B:314:TRP:CE3	1:B:319:ARG:HD2	2.41	0.56
1:D:136:ILE:C	1:D:137:LYS:HD3	2.25	0.56
1:C:343:LYS:HD2	1:D:343:LYS:CD	2.35	0.56
1:D:122:GLN:HE21	1:D:143:ARG:HH11	1.52	0.56
1:C:162:ILE:O	1:C:164:VAL:HG23	2.07	0.55
1:C:319:ARG:HH21	1:C:323:GLN:NE2	2.02	0.55
1:A:318:GLU:CG	1:C:293:GLU:CB	2.84	0.55
1:A:90:ILE:HD11	1:A:179:LEU:HD12	1.88	0.55
1:C:21:LYS:HD2	1:C:26:ALA:O	2.06	0.55
1:A:120:ASN:N	4:A:520:AS1:OD2	2.39	0.55
1:C:133:LYS:HD3	1:C:136:ILE:HD13	1.86	0.55
1:C:122:GLN:NE2	1:C:143:ARG:HH11	2.05	0.55
1:C:33:ALA:HB1	1:C:61:LEU:HD12	1.88	0.55
1:D:7:TYR:CZ	1:D:9:GLY:HA2	2.42	0.54
1:B:87:GLY:HA2	1:B:179:LEU:HD13	1.88	0.54
1:C:343:LYS:CD	1:D:343:LYS:HD2	2.36	0.54
1:C:2:LYS:HB2	1:C:108:ALA:HA	1.90	0.54
1:D:120:ASN:H	4:D:4520:AS1:CD	2.21	0.54
1:B:374:LYS:O	1:B:374:LYS:HG2	2.07	0.54
1:A:319:ARG:NH1	1:A:323:GLN:HE21	2.04	0.53
1:C:164:VAL:HG12	1:C:165:PRO:CD	2.36	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:189:GLU:O	1:C:191:PRO:HD3	2.08	0.53
1:B:70:VAL:O	1:B:74:MET:HG3	2.09	0.53
1:C:40:GLU:HG2	1:C:43:GLU:HG3	1.91	0.53
1:D:88:THR:HG21	4:D:4520:AS1:H12	1.90	0.53
1:B:318:GLU:CG	1:D:293:GLU:HB3	2.37	0.53
1:C:126:GLU:HG2	1:C:138:VAL:HG11	1.91	0.53
1:A:120:ASN:H	4:A:520:AS1:CD	2.22	0.53
1:D:351:ARG:CD	1:D:351:ARG:N	2.67	0.52
1:B:40:GLU:HB3	1:B:43:GLU:HG3	1.91	0.52
1:C:21:LYS:O	1:C:25:ARG:HA	2.08	0.52
1:D:146:SER:O	1:D:148:GLN:HG3	2.10	0.52
1:A:395:ARG:NH1	5:A:1722:HOH:O	2.41	0.52
1:C:32:THR:CG2	1:C:41:VAL:HG13	2.39	0.52
1:C:374:LYS:HE2	5:C:1586:HOH:O	2.08	0.52
1:C:151:LYS:HD2	1:C:151:LYS:N	2.24	0.52
1:D:345:ASN:HB3	1:D:347:TYR:CE1	2.44	0.52
1:B:372:ASP:OD2	1:B:372:ASP:C	2.48	0.52
1:A:343:LYS:HD2	1:B:343:LYS:HD3	1.90	0.52
1:D:98:HIS:CD2	1:D:101:ARG:HH21	2.23	0.52
1:C:373:GLN:HG3	1:D:117:GLY:HA3	1.92	0.52
1:D:18:LYS:O	1:D:22:GLU:HG3	2.09	0.52
1:B:35:ILE:N	1:B:37:GLN:HE22	2.04	0.52
1:B:1:MET:O	1:B:26:ALA:HB1	2.10	0.52
1:A:395:ARG:HH22	1:D:355:LYS:HE3	1.75	0.51
1:C:37:GLN:N	1:C:37:GLN:HE21	2.03	0.51
1:C:18:LYS:HE3	1:C:160:HIS:HB3	1.92	0.51
1:D:350:GLY:C	1:D:351:ARG:HD2	2.29	0.51
1:C:209:ALA:O	1:C:343:LYS:NZ	2.42	0.51
1:A:374:LYS:HZ2	1:D:374:LYS:HE3	1.75	0.51
1:B:359:ARG:HB3	1:B:361:ASP:HB2	1.93	0.51
1:B:318:GLU:HG3	1:D:293:GLU:HB3	1.93	0.51
1:B:179:LEU:O	1:B:272:THR:HG23	2.11	0.51
1:D:393:VAL:O	1:D:395:ARG:HG3	2.11	0.50
1:A:93:PRO:HG3	5:A:1256:HOH:O	2.10	0.50
1:C:180:HIS:ND1	1:C:180:HIS:C	2.64	0.50
1:C:227:ALA:HB2	1:C:232:ARG:HD2	1.92	0.50
1:D:93:PRO:HG3	5:D:1293:HOH:O	2.11	0.50
1:A:351:ARG:HD3	1:B:191:PRO:HB2	1.93	0.50
1:A:302:SER:HB2	1:A:303:PRO:HD3	1.92	0.50
1:C:35:ILE:N	1:C:37:GLN:HE22	2.06	0.50
1:A:180:HIS:C	1:A:180:HIS:ND1	2.64	0.50
1:A:119:GLY:HA3	4:A:520:AS1:CD	2.41	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:249:HIS:HD2	5:A:1159:HOH:O	1.94	0.50
1:A:12:ASP:OD1	1:A:150:ARG:NH2	2.33	0.50
1:C:73:MET:CE	1:C:90:ILE:HG12	2.42	0.49
1:C:154:ILE:HG23	1:C:164:VAL:CG1	2.30	0.49
1:C:90:ILE:HD11	1:C:179:LEU:HD12	1.94	0.49
1:C:164:VAL:CG1	1:C:165:PRO:CD	2.89	0.49
1:B:249:HIS:HD2	5:B:1825:HOH:O	1.95	0.49
1:B:34:ASP:HA	1:B:37:GLN:HE22	1.76	0.49
1:B:122:GLN:HE22	1:B:143:ARG:HD3	1.78	0.49
1:D:12:ASP:OD1	1:D:150:ARG:NH2	2.37	0.49
1:B:318:GLU:CG	1:D:293:GLU:HB2	2.35	0.49
1:C:179:LEU:O	1:C:272:THR:HG23	2.12	0.49
1:C:93:PRO:HG3	5:C:1406:HOH:O	2.12	0.48
1:B:202:MET:CE	5:B:1479:HOH:O	2.60	0.48
1:A:6:ALA:HB1	3:A:510:AMP:N3	2.29	0.48
1:C:6:ALA:HB2	1:C:99:LEU:HD11	1.96	0.48
1:C:204:GLN:HG2	5:C:1667:HOH:O	2.13	0.48
1:C:149:GLY:HA3	1:C:151:LYS:HE3	1.95	0.48
1:A:117:GLY:H	1:B:373:GLN:NE2	2.11	0.48
1:B:296:HIS:O	1:B:300:MET:HE2	2.14	0.48
1:C:302:SER:HB2	1:C:303:PRO:HD3	1.94	0.47
1:B:121:ASP:N	4:B:2520:AS1:OD2	2.28	0.47
1:C:151:LYS:CD	1:C:151:LYS:N	2.76	0.47
1:B:328:HIS:CE1	1:B:331:ARG:HH22	2.32	0.47
1:D:16:ILE:HG23	1:D:141:PRO:HG2	1.97	0.47
1:D:11:LEU:O	1:D:15:ILE:HG13	2.14	0.47
1:C:258:GLU:OE2	4:C:3520:AS1:N4	2.48	0.47
1:B:87:GLY:HA3	1:B:180:HIS:CD2	2.49	0.47
1:A:148:GLN:HB2	1:A:152:GLU:OE2	2.14	0.47
1:B:180:HIS:ND1	1:B:180:HIS:C	2.66	0.47
1:D:176:ALA:CB	1:D:181:ILE:HG12	2.45	0.47
1:A:24:TYR:OH	1:A:144:GLU:OE2	2.28	0.47
1:D:122:GLN:NE2	1:D:143:ARG:HH11	2.13	0.47
1:A:137:LYS:N	1:A:137:LYS:HD3	2.30	0.46
1:A:189:GLU:OE2	1:B:360:GLN:OE1	2.33	0.46
1:A:171:PRO:HB2	1:A:187:VAL:HG13	1.96	0.46
1:C:119:GLY:HA3	4:C:3520:AS1:CD	2.44	0.46
1:B:96:ALA:HB3	1:B:128:THR:HG21	1.97	0.46
1:A:71:PHE:O	1:A:75:ARG:HG3	2.14	0.46
1:C:71:PHE:O	1:C:75:ARG:HG3	2.14	0.46
1:B:321:ALA:HA	1:D:328:HIS:CD2	2.50	0.46
1:C:154:ILE:CG2	1:C:164:VAL:HG11	2.31	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:90:ILE:HD11	1:B:179:LEU:HD12	1.97	0.46
1:D:302:SER:HB2	1:D:303:PRO:HD3	1.97	0.46
1:B:164:VAL:HG13	1:B:165:PRO:HD2	1.96	0.46
1:A:359:ARG:NH1	1:A:361:ASP:OD2	2.48	0.46
1:D:180:HIS:C	1:D:180:HIS:ND1	2.69	0.46
1:D:2:LYS:HB2	1:D:108:ALA:HA	1.98	0.46
1:A:395:ARG:NH1	5:A:1382:HOH:O	2.49	0.46
3:C:3510:AMP:H5'1	3:C:3510:AMP:H8	1.79	0.46
1:A:328:HIS:CD2	1:C:321:ALA:HA	2.51	0.46
4:B:2520:AS1:CB	4:B:2520:AS1:C	2.94	0.45
1:B:180:HIS:CD2	5:B:1074:HOH:O	2.69	0.45
1:D:150:ARG:O	1:D:154:ILE:HD12	2.16	0.45
1:A:370:GLY:N	5:A:1286:HOH:O	2.49	0.45
1:B:6:ALA:HB1	3:B:2510:AMP:N3	2.31	0.45
1:D:90:ILE:HD11	1:D:179:LEU:HD12	1.98	0.45
1:D:262:VAL:HG13	1:D:262:VAL:O	2.16	0.45
1:D:96:ALA:HB3	1:D:128:THR:HG21	1.98	0.45
1:A:282:ARG:HD2	5:A:1351:HOH:O	2.17	0.45
4:B:2520:AS1:CB	4:B:2520:AS1:HA	2.21	0.45
1:B:151:LYS:CD	1:B:151:LYS:N	2.71	0.45
1:C:392:LEU:O	1:C:395:ARG:N	2.48	0.45
1:D:174:MET:HA	1:D:182:SER:O	2.16	0.45
1:A:118:LYS:HD3	1:B:365:PHE:CE2	2.51	0.45
1:C:393:VAL:O	1:C:395:ARG:NH1	2.49	0.45
1:B:40:GLU:CB	1:B:43:GLU:HG3	2.46	0.45
1:A:6:ALA:HB2	1:A:99:LEU:HD11	1.99	0.45
1:B:71:PHE:O	1:B:75:ARG:HG3	2.15	0.45
1:B:355:LYS:NZ	1:C:394:GLU:OE1	2.49	0.45
1:C:40:GLU:HG3	1:C:43:GLU:HG3	1.99	0.45
1:A:73:MET:CE	1:A:90:ILE:HG12	2.46	0.45
1:C:395:ARG:HG3	1:C:395:ARG:NH1	2.32	0.45
1:D:66:VAL:HA	1:D:70:VAL:HB	1.97	0.45
1:C:393:VAL:C	1:C:395:ARG:H	2.19	0.45
1:A:66:VAL:HA	1:A:70:VAL:HB	1.98	0.45
1:A:298:ARG:C	1:A:298:ARG:HD2	2.37	0.45
1:C:34:ASP:HB3	1:C:60:ASP:HA	1.99	0.45
1:D:66:VAL:HG21	1:D:236:ALA:HA	1.99	0.44
1:A:84:TYR:CZ	1:A:86:LEU:HA	2.52	0.44
1:B:319:ARG:HH21	1:B:323:GLN:NE2	2.11	0.44
1:B:180:HIS:HD2	5:B:1074:HOH:O	2.00	0.44
3:A:510:AMP:H8	3:A:510:AMP:H5'1	1.82	0.44
1:B:393:VAL:C	1:B:395:ARG:H	2.20	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:101:ARG:HD3	5:C:1261:HOH:O	2.17	0.44
1:D:117:GLY:HA2	1:D:122:GLN:NE2	2.32	0.44
1:B:393:VAL:C	1:B:395:ARG:N	2.69	0.44
1:C:17:LEU:O	1:C:17:LEU:HD23	2.17	0.44
1:C:120:ASN:H	4:C:3520:AS1:CD	2.30	0.44
3:D:4510:AMP:O2P	4:D:4520:AS1:C	2.66	0.44
1:A:238:LEU:C	1:A:238:LEU:HD23	2.38	0.44
1:A:117:GLY:HA3	1:B:373:GLN:NE2	2.32	0.44
1:D:92:ARG:NH2	3:D:4510:AMP:H5'2	2.33	0.44
4:C:3520:AS1:C	4:C:3520:AS1:CB	2.95	0.44
1:C:241:ARG:NH2	5:C:1653:HOH:O	2.47	0.44
1:B:32:THR:O	1:B:58:ALA:HA	2.17	0.44
1:B:318:GLU:HG2	1:D:293:GLU:HB3	1.99	0.43
1:D:146:SER:O	1:D:147:PHE:C	2.56	0.43
1:A:345:ASN:HB3	1:A:347:TYR:CE1	2.53	0.43
1:B:117:GLY:HA2	1:B:122:GLN:NE2	2.33	0.43
1:D:34:ASP:HA	1:D:37:GLN:HE22	1.83	0.43
1:C:238:LEU:C	1:C:238:LEU:HD23	2.38	0.43
1:B:157:ALA:HB1	1:B:162:ILE:HB	2.00	0.43
1:A:260:ARG:HD2	1:A:264:MET:SD	2.58	0.43
1:A:104:GLU:OE1	1:A:133:LYS:NZ	2.51	0.43
1:D:339:LEU:N	1:D:339:LEU:HD12	2.33	0.43
4:A:520:AS1:C	4:A:520:AS1:CB	2.97	0.43
1:B:298:ARG:O	1:B:298:ARG:HD2	2.18	0.43
1:B:37:GLN:HE21	1:B:37:GLN:N	2.14	0.43
1:B:180:HIS:HB2	1:B:271:GLU:O	2.19	0.43
1:A:282:ARG:NH2	5:A:1218:HOH:O	2.51	0.43
1:B:288:LEU:HG	1:B:351:ARG:HG3	1.99	0.43
1:D:35:ILE:N	1:D:37:GLN:NE2	2.56	0.43
1:A:171:PRO:O	1:A:186:GLY:HA3	2.18	0.43
1:A:86:LEU:HB3	1:A:89:SER:HB2	2.01	0.43
1:C:17:LEU:HD11	1:C:53:ALA:HB2	2.01	0.43
1:A:390:ARG:O	1:A:394:GLU:HG2	2.19	0.42
1:B:98:HIS:HD2	1:B:101:ARG:NH2	1.90	0.42
1:A:69:PHE:C	1:A:72:PRO:HD2	2.39	0.42
1:D:353:ALA:HA	1:D:354:PRO:HD3	1.89	0.42
1:A:180:HIS:HB2	1:A:271:GLU:O	2.20	0.42
1:C:29:ILE:HD12	1:C:55:LYS:HB3	2.02	0.42
1:A:171:PRO:HG2	1:A:172:TYR:CD2	2.54	0.42
1:D:314:TRP:CE3	1:D:319:ARG:HD2	2.55	0.42
1:D:117:GLY:HA2	1:D:122:GLN:HE22	1.85	0.42
1:B:50:ARG:NH1	5:B:1836:HOH:O	2.53	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:248:ARG:HG3	1:A:249:HIS:ND1	2.35	0.41
1:D:7:TYR:CE2	1:D:9:GLY:HA2	2.55	0.41
1:A:136:ILE:HG13	1:A:137:LYS:N	2.35	0.41
1:A:7:TYR:CE2	1:A:48:ALA:HB2	2.56	0.41
1:A:118:LYS:HD3	1:B:365:PHE:CZ	2.56	0.41
1:A:359:ARG:HG2	5:A:1609:HOH:O	2.20	0.41
1:D:395:ARG:NH1	1:D:395:ARG:CG	2.80	0.41
1:C:122:GLN:HE21	1:C:143:ARG:NH1	2.17	0.41
1:B:328:HIS:CD2	1:D:321:ALA:HA	2.56	0.41
1:A:298:ARG:O	1:A:298:ARG:HD2	2.21	0.41
1:A:262:VAL:HG13	5:A:1759:HOH:O	2.20	0.41
1:D:392:LEU:O	1:D:395:ARG:N	2.54	0.41
1:C:393:VAL:C	1:C:395:ARG:N	2.74	0.41
1:D:362:LEU:CD2	1:D:370:GLY:HA3	2.50	0.41
1:D:319:ARG:HH21	1:D:323:GLN:NE2	2.11	0.41
1:D:197:LYS:HB3	1:D:197:LYS:HE2	1.92	0.41
1:C:123:VAL:HG12	1:C:127:LEU:CD1	2.51	0.41
4:D:4520:AS1:C	4:D:4520:AS1:CB	2.98	0.41
1:C:158:GLU:HG3	1:C:164:VAL:HG21	2.03	0.41
1:C:314:TRP:HA	1:C:319:ARG:HD2	2.03	0.41
1:C:314:TRP:O	1:C:319:ARG:CD	2.67	0.41
1:B:345:ASN:HB3	1:B:347:TYR:CE1	2.55	0.41
1:C:71:PHE:CZ	1:C:225:PRO:HG2	2.56	0.41
1:A:181:ILE:HG22	1:A:183:TYR:CE1	2.56	0.41
1:B:355:LYS:CE	1:C:394:GLU:OE1	2.69	0.41
1:D:172:TYR:CZ	1:D:199:MET:HG3	2.57	0.40
1:A:32:THR:O	1:A:58:ALA:HA	2.21	0.40
1:C:101:ARG:NH2	5:C:1507:HOH:O	2.49	0.40
1:C:282:ARG:O	1:C:286:GLU:HG3	2.22	0.40
1:B:203:THR:HB	1:B:250:GLY:HA2	2.03	0.40
1:C:66:VAL:HA	1:C:70:VAL:HB	2.03	0.40
1:C:392:LEU:O	1:C:395:ARG:HA	2.22	0.40
1:A:394:GLU:O	1:A:395:ARG:HB2	2.22	0.40
1:C:204:GLN:HG3	1:C:205:ASP:N	2.36	0.40
1:D:298:ARG:HD2	1:D:298:ARG:O	2.21	0.40
1:B:64:GLU:OE1	1:B:98:HIS:HE1	2.05	0.40
1:B:7:TYR:CE2	1:B:9:GLY:HA2	2.55	0.40
1:C:242:LEU:HA	1:C:242:LEU:HD23	1.92	0.40
1:C:87:GLY:HA3	1:C:180:HIS:CD2	2.56	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	380/400 (95%)	373 (98%)	6 (2%)	1 (0%)	50	44
1	B	380/400 (95%)	369 (97%)	11 (3%)	0	100	100
1	C	380/400 (95%)	372 (98%)	8 (2%)	0	100	100
1	D	380/400 (95%)	372 (98%)	7 (2%)	1 (0%)	50	44
All	All	1520/1600 (95%)	1486 (98%)	32 (2%)	2 (0%)	59	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	343	LYS
1	D	343	LYS

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	310/319 (97%)	293 (94%)	17 (6%)	30	23
1	B	310/319 (97%)	292 (94%)	18 (6%)	28	21
1	C	310/319 (97%)	295 (95%)	15 (5%)	35	28
1	D	310/319 (97%)	290 (94%)	20 (6%)	24	17
All	All	1240/1276 (97%)	1170 (94%)	70 (6%)	30	22

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	LYS

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Mol	Chain	Res	Type
1	A	37	GLN
1	A	90	ILE
1	A	94	LEU
1	A	118	LYS
1	A	135	ASP
1	A	137	LYS
1	A	145	TRP
1	A	151	LYS
1	A	179	LEU
1	A	180	HIS
1	A	182	SER
1	A	188	LEU
1	A	239	LEU
1	A	267	ARG
1	A	351	ARG
1	A	361	ASP
1	B	37	GLN
1	B	62	LYS
1	B	90	ILE
1	B	94	LEU
1	B	145	TRP
1	B	146	SER
1	B	151	LYS
1	B	154	ILE
1	B	179	LEU
1	B	180	HIS
1	B	182	SER
1	B	239	LEU
1	B	267	ARG
1	B	318	GLU
1	B	319	ARG
1	B	351	ARG
1	B	361	ASP
1	B	372	ASP
1	C	25	ARG
1	C	37	GLN
1	C	90	ILE
1	C	145	TRP
1	C	151	LYS
1	C	178	LEU
1	C	179	LEU
1	C	180	HIS

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Mol	Chain	Res	Type
1	C	182	SER
1	C	239	LEU
1	C	267	ARG
1	C	319	ARG
1	C	345	ASN
1	C	351	ARG
1	C	373	GLN
1	D	37	GLN
1	D	62	LYS
1	D	90	ILE
1	D	94	LEU
1	D	145	TRP
1	D	174	MET
1	D	178	LEU
1	D	179	LEU
1	D	180	HIS
1	D	188	LEU
1	D	239	LEU
1	D	262	VAL
1	D	264	MET
1	D	267	ARG
1	D	319	ARG
1	D	351	ARG
1	D	364	SER
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 (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	HIS
1	A	180	HIS
1	A	249	HIS
1	A	323	GLN
1	A	360	GLN
1	A	373	GLN
1	B	37	GLN
1	B	98	HIS
1	B	122	GLN
1	B	148	GLN
1	B	180	HIS

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Mol	Chain	Res	Type
1	B	249	HIS
1	B	323	GLN
1	B	360	GLN
1	B	373	GLN
1	C	37	GLN
1	C	98	HIS
1	C	122	GLN
1	C	180	HIS
1	C	323	GLN
1	C	345	ASN
1	C	360	GLN
1	C	373	GLN
1	D	37	GLN
1	D	98	HIS
1	D	122	GLN
1	D	180	HIS
1	D	249	HIS
1	D	323	GLN
1	D	360	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

12 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected

value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	AMP	A	510	-	25,25,25	2.88	15 (60%)	38,38,38	2.42	12 (31%)
4	AS1	A	520	-	19,19,19	3.92	5 (26%)	24,24,24	7.16	13 (54%)
2	SO4	A	540	-	4,4,4	2.70	2 (50%)	6,6,6	1.02	0
3	AMP	B	2510	-	25,25,25	2.99	15 (60%)	38,38,38	2.48	13 (34%)
4	AS1	B	2520	-	19,19,19	3.91	4 (21%)	24,24,24	7.16	13 (54%)
2	SO4	B	2540	-	4,4,4	3.15	2 (50%)	6,6,6	0.96	0
3	AMP	C	3510	-	25,25,25	2.69	14 (56%)	38,38,38	2.45	14 (36%)
4	AS1	C	3520	-	19,19,19	3.86	5 (26%)	24,24,24	7.09	13 (54%)
2	SO4	C	3540	-	4,4,4	3.08	2 (50%)	6,6,6	0.84	0
3	AMP	D	4510	-	25,25,25	2.70	15 (60%)	38,38,38	2.45	14 (36%)
4	AS1	D	4520	-	19,19,19	3.93	5 (26%)	24,24,24	7.39	13 (54%)
2	SO4	D	4540	-	4,4,4	2.97	2 (50%)	6,6,6	0.92	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
3	AMP	A	510	-	-	0/10/26/26	0/1/3/3
4	AS1	A	520	-	2/2/6/8	1/23/23/23	0/0/0/0
2	SO4	A	540	-	-	0/0/0/0	0/0/0/0
3	AMP	B	2510	-	-	0/10/26/26	0/1/3/3
4	AS1	B	2520	-	2/2/6/8	1/23/23/23	0/0/0/0
2	SO4	B	2540	-	-	0/0/0/0	0/0/0/0
3	AMP	C	3510	-	-	0/10/26/26	0/1/3/3
4	AS1	C	3520	-	2/2/6/8	1/23/23/23	0/0/0/0
2	SO4	C	3540	-	-	0/0/0/0	0/0/0/0
3	AMP	D	4510	-	-	0/10/26/26	0/1/3/3
4	AS1	D	4520	-	2/2/6/8	0/23/23/23	0/0/0/0
2	SO4	D	4540	-	-	0/0/0/0	0/0/0/0

All (86) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	4520	AS1	CB-CA	12.71	1.81	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	2520	AS1	CB-CA	12.34	1.80	1.53
4	A	520	AS1	CB-CA	12.34	1.80	1.53
4	C	3520	AS1	CB-CA	12.23	1.80	1.53
4	B	2520	AS1	C1-N2	7.23	1.63	1.46
4	A	520	AS1	C1-N2	6.99	1.62	1.46
4	C	3520	AS1	C-N3	6.81	1.43	1.33
4	C	3520	AS1	C1-N2	6.58	1.61	1.46
4	D	4520	AS1	C1-N2	6.53	1.61	1.46
4	B	2520	AS1	C-N3	6.28	1.42	1.33
4	A	520	AS1	C-N3	6.16	1.42	1.33
4	D	4520	AS1	C-N3	6.01	1.42	1.33
4	A	520	AS1	CA-N1	-5.89	1.32	1.45
4	D	4520	AS1	CA-N1	-5.66	1.33	1.45
4	B	2520	AS1	CA-N1	-5.53	1.33	1.45
4	C	3520	AS1	CA-N1	-5.26	1.34	1.45
3	B	2510	AMP	C2'-C1'	5.17	1.60	1.53
2	B	2540	SO4	O1-S	5.15	1.63	1.47
3	B	2510	AMP	O4'-C1'	5.08	1.49	1.41
3	D	4510	AMP	C8-N9	5.03	1.44	1.36
3	A	510	AMP	C8-N9	4.99	1.44	1.36
3	A	510	AMP	C2'-C1'	4.97	1.60	1.53
3	A	510	AMP	O4'-C1'	4.76	1.48	1.41
3	C	3510	AMP	C8-N9	4.70	1.43	1.36
3	B	2510	AMP	P-O3P	-4.68	1.37	1.54
2	C	3540	SO4	O1-S	4.68	1.62	1.47
2	D	4540	SO4	O1-S	4.66	1.62	1.47
3	A	510	AMP	C3'-C4'	4.63	1.65	1.53
3	B	2510	AMP	C8-N9	4.54	1.43	1.36
3	B	2510	AMP	C3'-C4'	4.33	1.64	1.53
3	A	510	AMP	P-O3P	-4.33	1.38	1.54
3	A	510	AMP	P-O1P	4.32	1.65	1.51
3	C	3510	AMP	C3'-C4'	4.32	1.64	1.53
3	C	3510	AMP	O4'-C1'	4.30	1.48	1.41
3	D	4510	AMP	P-O3P	-4.29	1.39	1.54
3	D	4510	AMP	C3'-C4'	4.27	1.64	1.53
3	C	3510	AMP	C2'-C1'	4.26	1.59	1.53
3	D	4510	AMP	P-O1P	4.23	1.65	1.51
3	C	3510	AMP	P-O1P	4.17	1.65	1.51
3	B	2510	AMP	P-O1P	4.06	1.64	1.51
3	D	4510	AMP	C2'-C1'	4.05	1.59	1.53
2	A	540	SO4	O1-S	4.02	1.60	1.47
2	C	3540	SO4	O3-S	-3.97	1.33	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	2510	AMP	P-O2P	3.93	1.69	1.54
3	A	510	AMP	P-O5'	-3.82	1.46	1.60
2	D	4540	SO4	O3-S	-3.67	1.34	1.47
3	C	3510	AMP	P-O3P	-3.64	1.41	1.54
2	A	540	SO4	O3-S	-3.55	1.35	1.47
3	B	2510	AMP	P-O5'	-3.55	1.47	1.60
2	B	2540	SO4	O3-S	-3.50	1.35	1.47
3	D	4510	AMP	P-O5'	-3.35	1.48	1.60
3	D	4510	AMP	O4'-C1'	3.18	1.46	1.41
3	D	4510	AMP	O3'-C3'	3.10	1.50	1.43
3	D	4510	AMP	O5'-C5'	3.08	1.57	1.44
3	A	510	AMP	O3'-C3'	3.07	1.50	1.43
3	B	2510	AMP	O5'-C5'	3.06	1.57	1.44
3	B	2510	AMP	C2-N1	3.02	1.39	1.33
3	C	3510	AMP	P-O5'	-2.97	1.49	1.60
3	C	3510	AMP	C2-N3	2.93	1.38	1.32
3	C	3510	AMP	P-O2P	2.86	1.65	1.54
3	C	3510	AMP	C2-N1	2.86	1.39	1.33
3	C	3510	AMP	O5'-C5'	2.83	1.56	1.44
3	A	510	AMP	C1'-N9	2.83	1.57	1.48
3	B	2510	AMP	C2-N3	2.80	1.37	1.32
3	A	510	AMP	C2-N1	2.70	1.39	1.33
3	D	4510	AMP	C2-N1	2.55	1.38	1.33
3	A	510	AMP	C2-N3	2.55	1.37	1.32
3	B	2510	AMP	O3'-C3'	2.45	1.48	1.43
3	D	4510	AMP	C2-N3	2.45	1.37	1.32
3	A	510	AMP	O5'-C5'	2.44	1.55	1.44
4	D	4520	AS1	OG1-CG	-2.41	1.21	1.30
3	D	4510	AMP	C1'-N9	2.41	1.56	1.48
3	B	2510	AMP	C5-C4	2.40	1.45	1.40
3	A	510	AMP	C5-C4	2.39	1.45	1.40
4	C	3520	AS1	OG1-CG	-2.39	1.21	1.30
3	C	3510	AMP	C4-N3	2.31	1.39	1.35
3	D	4510	AMP	P-O2P	2.29	1.63	1.54
3	C	3510	AMP	O3'-C3'	2.25	1.48	1.43
3	B	2510	AMP	C2'-C3'	2.25	1.59	1.53
3	B	2510	AMP	C1'-N9	2.23	1.55	1.48
3	A	510	AMP	C2'-C3'	2.23	1.59	1.53
4	A	520	AS1	OG1-CG	-2.22	1.22	1.30
3	C	3510	AMP	C1'-N9	2.21	1.55	1.48
3	D	4510	AMP	C5-C4	2.15	1.45	1.40
3	D	4510	AMP	C2'-C3'	2.14	1.59	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	510	AMP	P-O2P	2.11	1.62	1.54

All (105) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	4520	AS1	C3-C2-C1	23.42	171.94	111.94
4	A	520	AS1	C3-C2-C1	22.86	170.51	111.94
4	B	2520	AS1	C3-C2-C1	22.77	170.28	111.94
4	C	3520	AS1	C3-C2-C1	22.14	168.67	111.94
4	D	4520	AS1	C2-C1-N2	17.58	165.14	112.21
4	A	520	AS1	C2-C1-N2	15.74	159.61	112.21
4	B	2520	AS1	C2-C1-N2	15.59	159.17	112.21
4	C	3520	AS1	C2-C1-N2	15.58	159.12	112.21
4	B	2520	AS1	C5-C4-N4	11.73	128.78	109.36
4	C	3520	AS1	C5-C4-N4	11.62	128.60	109.36
4	D	4520	AS1	C5-C4-N4	11.39	128.22	109.36
4	A	520	AS1	C5-C4-N4	11.30	128.07	109.36
4	C	3520	AS1	N2-C-N1	11.08	137.27	119.13
4	B	2520	AS1	N2-C-N1	11.07	137.25	119.13
4	D	4520	AS1	N2-C-N1	10.85	136.88	119.13
4	A	520	AS1	N2-C-N1	10.54	136.39	119.13
4	A	520	AS1	C1-N2-C	8.15	138.34	124.03
4	B	2520	AS1	C1-N2-C	7.20	136.68	124.03
4	C	3520	AS1	C1-N2-C	7.09	136.49	124.03
4	D	4520	AS1	C1-N2-C	7.07	136.46	124.03
4	B	2520	AS1	N2-C-N3	-7.06	106.92	120.23
4	C	3520	AS1	N2-C-N3	-6.85	107.32	120.23
4	D	4520	AS1	N2-C-N3	-6.68	107.64	120.23
4	A	520	AS1	N2-C-N3	-6.45	108.08	120.23
3	B	2510	AMP	C8-N9-C4	-6.38	102.03	106.90
3	C	3510	AMP	N3-C2-N1	-6.37	123.38	128.71
3	D	4510	AMP	N3-C2-N1	-5.98	123.71	128.71
3	A	510	AMP	N3-C2-N1	-5.94	123.74	128.71
3	A	510	AMP	C8-N9-C4	-5.86	102.43	106.90
3	B	2510	AMP	N3-C2-N1	-5.62	124.01	128.71
3	C	3510	AMP	C4-C5-N7	5.59	114.31	109.52
3	D	4510	AMP	C8-N9-C4	-5.57	102.65	106.90
3	C	3510	AMP	C8-N9-C4	-5.39	102.78	106.90
3	B	2510	AMP	C4-C5-N7	5.15	113.93	109.52
3	A	510	AMP	C4-C5-N7	4.87	113.69	109.52
3	D	4510	AMP	C4-C5-N7	4.82	113.65	109.52
3	B	2510	AMP	P-O5'-C5'	4.81	132.10	118.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4510	AMP	O4'-C1'-N9	4.56	112.68	108.44
4	D	4520	AS1	OG1-CG-OG2	-4.53	113.83	124.07
4	C	3520	AS1	OG1-CG-OG2	-4.44	114.04	124.07
4	A	520	AS1	CA-CB-CD	4.42	127.13	112.94
3	C	3510	AMP	O4'-C1'-N9	4.39	112.52	108.44
4	C	3520	AS1	CA-CB-CD	4.34	126.87	112.94
3	A	510	AMP	P-O5'-C5'	4.33	130.72	118.19
4	B	2520	AS1	OG1-CG-OG2	-4.31	114.32	124.07
4	D	4520	AS1	CA-CB-CD	4.29	126.72	112.94
4	A	520	AS1	OG1-CG-OG2	-4.27	114.42	124.07
4	B	2520	AS1	CA-CB-CD	4.22	126.48	112.94
3	D	4510	AMP	P-O5'-C5'	4.12	130.11	118.19
4	A	520	AS1	CB-CA-CG	3.97	118.41	111.19
4	C	3520	AS1	N1-C-N3	-3.94	113.09	120.14
4	C	3520	AS1	CB-CA-N1	3.92	118.66	110.61
3	D	4510	AMP	O2P-P-O5'	3.84	117.25	106.65
3	A	510	AMP	O4'-C1'-N9	3.80	111.97	108.44
3	B	2510	AMP	O2P-P-O5'	3.77	117.06	106.65
3	A	510	AMP	O3P-P-O5'	3.75	116.98	106.65
4	B	2520	AS1	CG-CA-N1	3.67	119.98	110.53
4	D	4520	AS1	CB-CA-N1	3.65	118.11	110.61
4	B	2520	AS1	N1-C-N3	-3.64	113.63	120.14
4	C	3520	AS1	CB-CA-CG	3.63	117.79	111.19
4	A	520	AS1	CB-CA-N1	3.61	118.03	110.61
4	D	4520	AS1	CB-CA-CG	3.50	117.56	111.19
4	B	2520	AS1	CB-CA-CG	3.50	117.56	111.19
3	A	510	AMP	O2P-P-O5'	3.49	116.29	106.65
3	B	2510	AMP	O3P-P-O5'	3.49	116.29	106.65
4	D	4520	AS1	N1-C-N3	-3.43	114.01	120.14
3	C	3510	AMP	O2P-P-O5'	3.42	116.08	106.65
4	C	3520	AS1	CG-CA-N1	3.42	119.34	110.53
4	A	520	AS1	C3-C4-N4	3.38	118.40	110.14
4	D	4520	AS1	CG-CA-N1	3.37	119.22	110.53
4	B	2520	AS1	CB-CA-N1	3.34	117.47	110.61
4	C	3520	AS1	C3-C4-N4	3.33	118.28	110.14
4	A	520	AS1	N1-C-N3	-3.31	114.23	120.14
3	C	3510	AMP	N6-C6-N1	3.29	125.83	119.36
4	A	520	AS1	CG-CA-N1	3.28	119.00	110.53
3	C	3510	AMP	O3P-P-O5'	3.28	115.69	106.65
3	B	2510	AMP	C3'-C2'-C1'	3.23	105.97	100.91
4	D	4520	AS1	C3-C4-N4	3.23	118.03	110.14
4	B	2520	AS1	C3-C4-N4	3.13	117.80	110.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4510	AMP	O3P-P-O5'	3.10	115.21	106.65
3	C	3510	AMP	O2P-P-O1P	-3.07	100.39	110.44
3	B	2510	AMP	O2P-P-O1P	-3.00	100.64	110.44
3	A	510	AMP	O2P-P-O1P	-2.99	100.66	110.44
3	C	3510	AMP	P-O5'-C5'	2.90	126.58	118.19
3	B	2510	AMP	N6-C6-N1	2.90	125.06	119.36
3	D	4510	AMP	O2P-P-O1P	-2.89	100.98	110.44
3	A	510	AMP	C3'-C2'-C1'	2.88	105.41	100.91
3	D	4510	AMP	C3'-C2'-C1'	2.84	105.36	100.91
3	B	2510	AMP	O3'-C3'-C4'	-2.84	102.72	111.08
3	D	4510	AMP	O3'-C3'-C4'	-2.80	102.83	111.08
3	B	2510	AMP	O5'-C5'-C4'	2.80	119.21	108.94
3	C	3510	AMP	O3'-C3'-C4'	-2.72	103.07	111.08
3	A	510	AMP	N6-C6-N1	2.71	124.68	119.36
3	D	4510	AMP	N6-C6-N1	2.55	124.38	119.36
3	B	2510	AMP	O4'-C4'-C5'	-2.55	100.25	109.36
3	D	4510	AMP	C2-N1-C6	2.54	123.35	118.77
3	A	510	AMP	C2-N1-C6	2.51	123.31	118.77
3	C	3510	AMP	C3'-C2'-C1'	2.40	104.66	100.91
3	D	4510	AMP	C8-N9-C1'	2.37	131.05	126.38
3	A	510	AMP	O3'-C3'-C4'	-2.34	104.19	111.08
3	B	2510	AMP	C2-N1-C6	2.33	122.98	118.77
3	C	3510	AMP	C2-N1-C6	2.32	122.95	118.77
3	C	3510	AMP	O5'-C5'-C4'	2.30	117.39	108.94
3	D	4510	AMP	C2'-C3'-C4'	-2.20	98.28	102.65
3	C	3510	AMP	O3'-C3'-C2'	2.03	118.44	111.83

All (8) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	D	4520	AS1	C4
4	D	4520	AS1	CA
4	C	3520	AS1	C4
4	C	3520	AS1	CA
4	B	2520	AS1	C4
4	B	2520	AS1	CA
4	A	520	AS1	C4
4	A	520	AS1	CA

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	2520	AS1	C2-C1-N2-C
4	C	3520	AS1	C2-C1-N2-C
4	A	520	AS1	C2-C1-N2-C

There are no ring outliers.

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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	386/400 (96%)	-0.27	9 (2%) 57 57	20, 27, 49, 78	0
1	B	386/400 (96%)	-0.31	7 (1%) 65 66	20, 27, 47, 65	0
1	C	386/400 (96%)	-0.13	8 (2%) 60 61	21, 29, 55, 72	0
1	D	386/400 (96%)	-0.19	10 (2%) 53 53	21, 30, 52, 72	0
All	All	1544/1600 (96%)	-0.22	34 (2%) 57 59	20, 29, 51, 78	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	165	PRO	5.4
1	D	165	PRO	4.7
1	A	165	PRO	4.6
1	B	370	GLY	4.4
1	A	370	GLY	4.3
1	A	365	PHE	4.1
1	B	165	PRO	4.1
1	C	370	GLY	4.1
1	B	365	PHE	3.1
1	A	197	LYS	3.0
1	D	370	GLY	2.9
1	D	148	GLN	2.9
1	D	39	GLU	2.8
1	C	395	ARG	2.7
1	D	159	ALA	2.7
1	C	40	GLU	2.6
1	C	365	PHE	2.6
1	C	25	ARG	2.6
1	D	371	TYR	2.6
1	D	395	ARG	2.5
1	A	148	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	395	ARG	2.4
1	C	43	GLU	2.4
1	A	38	GLY	2.3
1	C	164	VAL	2.3
1	D	201	ARG	2.3
1	D	164	VAL	2.3
1	D	365	PHE	2.3
1	A	364	SER	2.3
1	A	164	VAL	2.2
1	B	40	GLU	2.1
1	B	148	GLN	2.1
1	A	361	ASP	2.1
1	B	371	TYR	2.1

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	AS1	C	3520	20/20	0.27	8.80	23,35,54,54	0
4	AS1	A	520	20/20	0.24	7.91	24,36,53,54	0
4	AS1	D	4520	20/20	0.26	7.44	23,39,56,58	0
3	AMP	A	510	23/23	0.28	6.92	36,45,68,68	23
3	AMP	C	3510	23/23	0.25	5.67	41,47,69,70	23
4	AS1	B	2520	20/20	0.23	5.03	22,36,59,61	0
3	AMP	D	4510	23/23	0.28	4.81	40,47,69,70	23
3	AMP	B	2510	23/23	0.25	4.54	31,41,60,63	23

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	SO4	B	2540	5/5	0.10	-0.13	37,40,41,42	0
2	SO4	A	540	5/5	0.09	-0.32	41,43,44,44	0
2	SO4	C	3540	5/5	0.08	-1.32	43,44,45,45	0
2	SO4	D	4540	5/5	0.07	-1.95	44,45,46,47	0

6.5 Other polymers ⓘ

There are no such residues in this entry.