



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 04:07 PM GMT

PDB ID : 2J91  
Title : CRYSTAL STRUCTURE OF HUMAN ADENYLOSUCCINATE LYASE IN COMPLEX WITH AMP  
Authors : Stenmark, P.; Moche, M.; Arrowsmith, C.; Berglund, H.; Busam, R.; Collins, R.; Edwards, A.; Ericsson, U.B.; Flodin, S.; Flores, A.; Graslund, S.; Hammarstrom, M.; Hallberg, B.M.; Holmberg Schiavone, L.; Hogbom, M.; Johansson, I.; Karlberg, T.; Kosinska, U.; Kotenyova, T.; Magnusdottir, A.; Nilsson, M.E.; Nilsson-Ehle, P.; Nyman, T.; Ogg, D.; Persson, C.; Sagemark, J.; Sundstrom, M.; Uppenberg, J.; Uppsten, M.; Thorsell, A.G.; Van Den Berg, S.; Wallden, K.; Weigelt, J.; Nordlund, P.  
Deposited on : 2006-11-01  
Resolution : 1.80 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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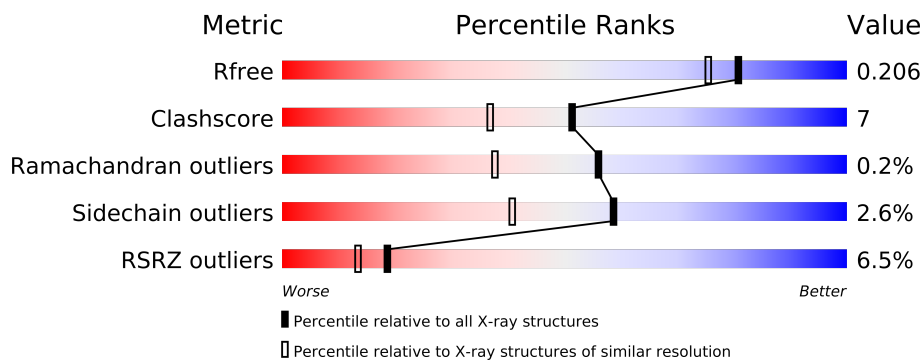
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 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 1.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	66092	3513 (1.80-1.80)
Clashscore	79885	4461 (1.80-1.80)
Ramachandran outliers	78287	4404 (1.80-1.80)
Sidechain outliers	78261	4403 (1.80-1.80)
RSRZ outliers	66119	3515 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	503	
1	B	503	
1	C	503	
1	D	503	

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

Mol	Type	Chain	Res	Geometry	Electron density
2	AMP	C	1000	-	X
4	GOL	A	1475	-	X
4	GOL	D	1474	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 16310 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 ADENYLOSUCCINATE LYASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	457	Total	C	N	O	S	0	2	0
			3664	2308	652	679	25			
1	B	463	Total	C	N	O	S	0	1	0
			3702	2331	658	687	26			
1	C	456	Total	C	N	O	S	0	3	0
			3662	2306	655	675	26			
1	D	456	Total	C	N	O	S	0	4	0
			3667	2313	654	675	25			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	63	ARG	GLN	ENGINEERED MUTATION	UNP P30566
B	63	ARG	GLN	ENGINEERED MUTATION	UNP P30566
C	63	ARG	GLN	ENGINEERED MUTATION	UNP P30566
D	63	ARG	GLN	ENGINEERED MUTATION	UNP P30566

- Molecule 2 is ADENOSINE MONOPHOSPHATE (three-letter code: AMP) (formula: C<sub>10</sub>H<sub>14</sub>N<sub>5</sub>O<sub>7</sub>P).



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

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

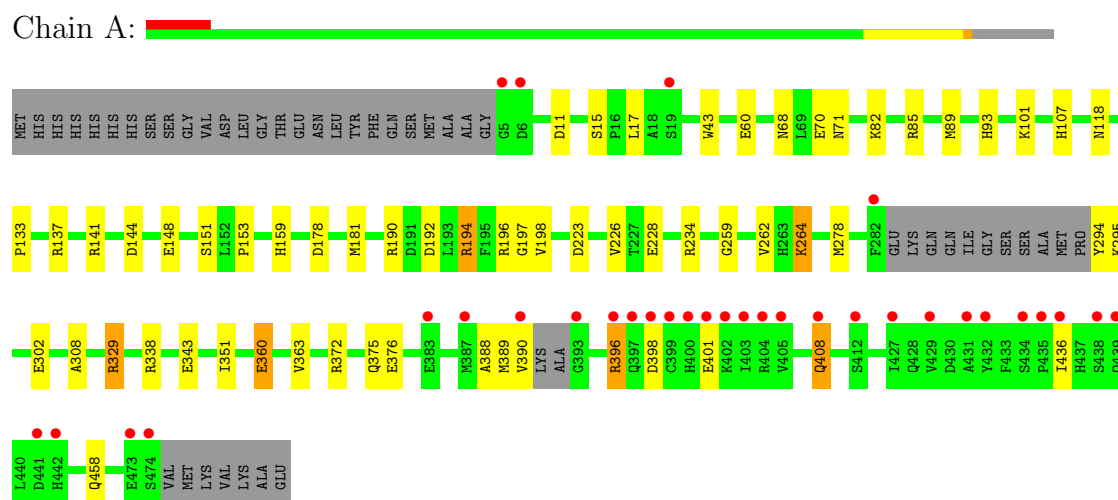
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	334	Total	O	0	0
			334	334		
5	B	377	Total	O	0	0
			377	377		
5	C	373	Total	O	0	0
			373	373		
5	D	411	Total	O	0	0
			411	411		

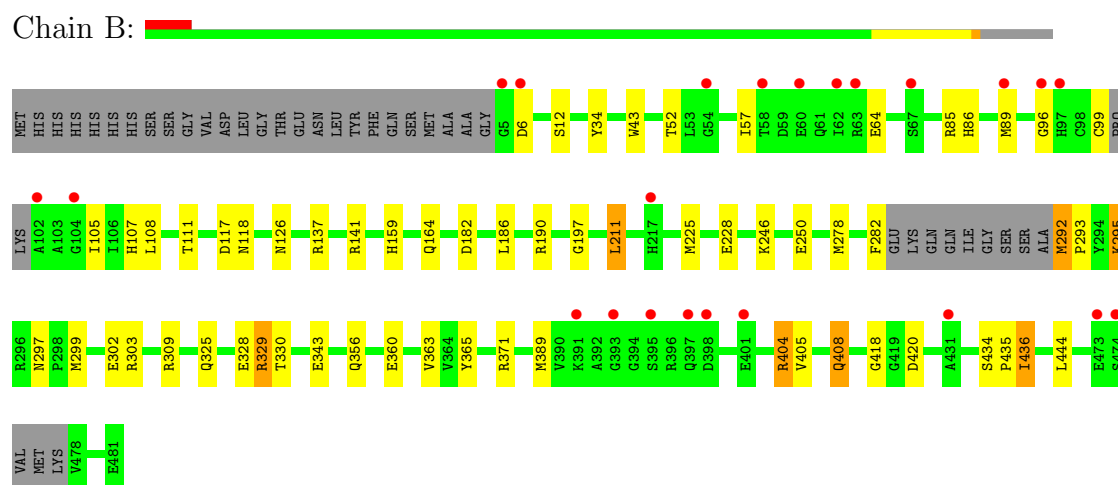
### 3 Residue-property plots

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.

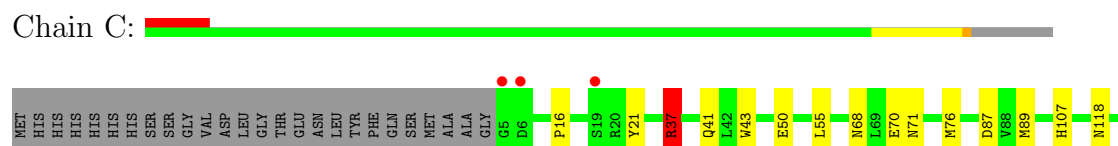
#### • Molecule 1: ADENYLOSUCCINATE LYASE

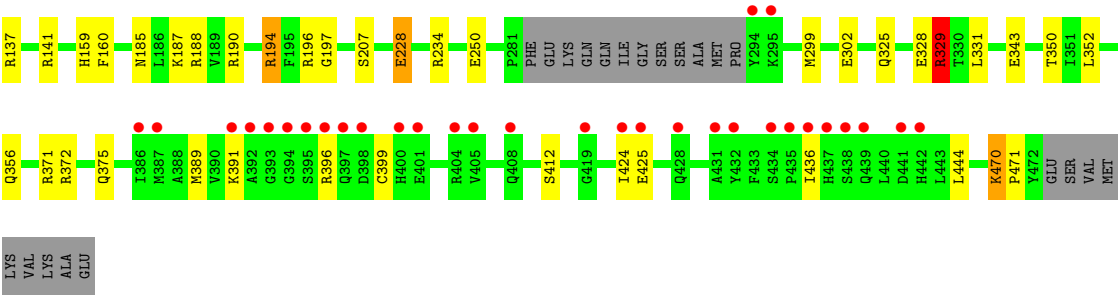


#### • Molecule 1: ADENYLOSUCCINATE LYASE



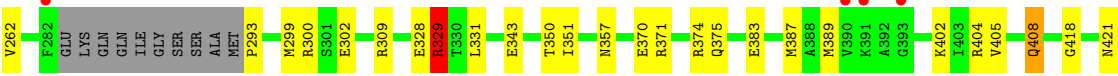
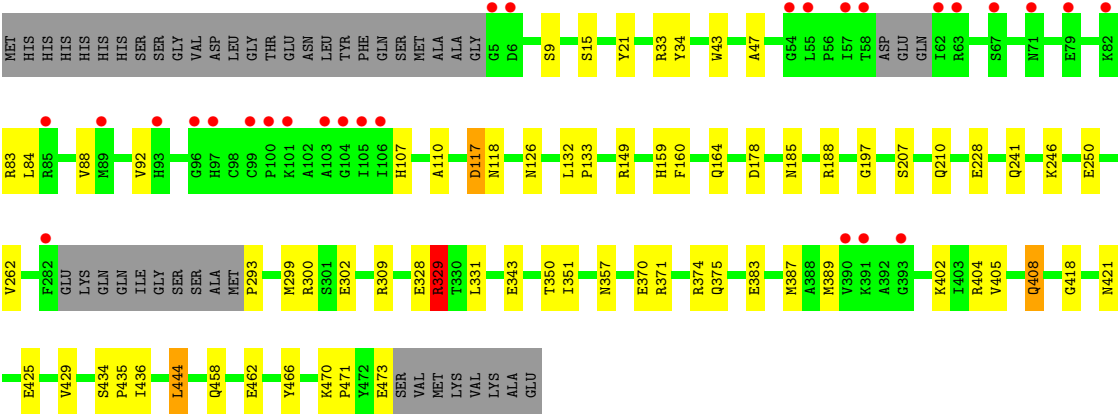
#### • Molecule 1: ADENYLOSUCCINATE LYASE





● Molecule 1: ADENYLOSUCCINATE LYASE

Chain D:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.37Å 104.34Å 213.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	106.60 – 1.80 14.98 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (106.60-1.80) 100.0 (14.98-1.80)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.79 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.157 , 0.199 0.168 , 0.206	Depositor DCC
$R_{free}$ test set	8809 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	18.7	Xtriage
Anisotropy	0.203	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 58.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 176155 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	16310	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 4.51% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities 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: GOL, CL, AMP

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	0.68	0/3737	0.76	6/5044 (0.1%)
1	B	0.70	1/3771 (0.0%)	0.74	2/5086 (0.0%)
1	C	0.71	0/3735	0.79	6/5041 (0.1%)
1	D	0.68	0/3744	0.76	5/5052 (0.1%)
All	All	0.69	1/14987 (0.0%)	0.76	19/20223 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	295	LYS	CE-NZ	6.33	1.64	1.49

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	194	ARG	NE-CZ-NH2	-10.51	115.04	120.30
1	A	194	ARG	NE-CZ-NH2	-7.88	116.36	120.30
1	C	194	ARG	NE-CZ-NH1	7.48	124.04	120.30
1	B	309	ARG	NE-CZ-NH2	-6.23	117.19	120.30
1	A	137	ARG	NE-CZ-NH2	-6.23	117.19	120.30
1	A	194	ARG	NE-CZ-NH1	6.09	123.35	120.30
1	A	137	ARG	NE-CZ-NH1	5.99	123.30	120.30
1	D	149	ARG	NE-CZ-NH2	-5.93	117.33	120.30
1	D	444	LEU	CA-CB-CG	-5.84	101.86	115.30
1	C	87	ASP	CB-CG-OD2	-5.77	113.11	118.30
1	C	87	ASP	CB-CG-OD1	5.61	123.35	118.30
1	C	37	ARG	NE-CZ-NH2	-5.53	117.53	120.30
1	D	178	ASP	CB-CG-OD1	5.48	123.23	118.30
1	A	338	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	D	309	ARG	NE-CZ-NH2	-5.45	117.58	120.30
1	C	37	ARG	CG-CD-NE	-5.40	100.47	111.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	178	ASP	CB-CG-OD2	-5.21	113.61	118.30
1	A	295	LYS	CD-CE-NZ	5.10	123.43	111.70
1	B	295	LYS	CD-CE-NZ	5.08	123.39	111.70

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	3664	0	3674	47	0
1	B	3702	0	3712	64	0
1	C	3662	0	3686	47	0
1	D	3667	0	3692	72	0
2	A	23	0	12	1	0
2	B	23	0	12	1	0
2	C	23	0	12	2	0
2	D	23	0	12	3	0
3	A	1	0	0	0	0
3	B	1	0	0	1	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	6	0	8	0	0
4	B	6	0	8	0	0
4	C	6	0	8	0	0
4	D	6	0	8	0	0
5	A	334	0	0	11	1
5	B	377	0	0	17	1
5	C	373	0	0	12	0
5	D	411	0	0	17	0
All	All	16310	0	14844	217	1

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 7.

All (217) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:B:2094:HOH:O	1:C:331:LEU:HD23	1.29	1.29
5:A:2101:HOH:O	1:D:331:LEU:HD23	1.21	1.26
1:A:388:ALA:HB1	1:A:436:ILE:HD11	1.37	1.04
1:B:389:MET:HE3	1:B:436:ILE:HD13	1.38	1.02
1:D:331:LEU:HD13	5:D:2265:HOH:O	1.59	1.00
1:A:178:ASP:HA	1:A:181:MET:HE3	1.46	0.98
1:B:389:MET:CE	1:B:436:ILE:HD13	1.94	0.96
1:D:34:TYR:OH	1:D:126:ASN:ND2	1.99	0.95
1:D:389:MET:HE2	1:D:436:ILE:HG23	1.47	0.93
1:A:178:ASP:HA	1:A:181:MET:CE	1.99	0.92
1:A:389:MET:HE3	1:A:436:ILE:HD13	1.58	0.85
1:D:375:GLN:HE22	1:D:421:ASN:H	1.22	0.85
1:A:388:ALA:CB	1:A:436:ILE:HD11	2.06	0.85
1:D:389:MET:HE3	1:D:436:ILE:HG13	1.60	0.84
5:B:2094:HOH:O	1:C:331:LEU:CD2	1.97	0.82
1:D:293:PRO:N	5:D:2234:HOH:O	2.13	0.82
1:D:389:MET:CE	1:D:436:ILE:HG13	2.12	0.78
1:B:299:MET:HE1	5:C:2119:HOH:O	1.84	0.78
1:A:264:LYS:NZ	5:A:2222:HOH:O	2.07	0.77
1:A:294:TYR:N	5:A:2236:HOH:O	2.16	0.77
1:B:389:MET:HE3	1:B:436:ILE:CD1	2.14	0.77
1:D:299:MET:HE2	5:D:2253:HOH:O	1.86	0.76
1:D:299:MET:CE	5:D:2253:HOH:O	2.33	0.75
1:A:70:GLU:OE1	5:A:2062:HOH:O	2.05	0.74
1:D:458:GLN:HG3	5:D:2387:HOH:O	1.86	0.74
1:B:111:THR:HG22	3:B:1001:CL:CL	2.25	0.73
1:B:111:THR:CG2	5:B:2098:HOH:O	2.36	0.73
1:D:164:GLN:OE1	5:D:2142:HOH:O	2.05	0.72
1:B:137:ARG:HH11	1:B:356:GLN:HE22	1.36	0.72
1:B:299:MET:SD	5:C:2043:HOH:O	2.47	0.72
1:C:299[A]:MET:HE3	2:C:1000:AMP:C8	2.25	0.71
1:B:126:ASN:HB3	5:B:2107:HOH:O	1.89	0.71
1:B:404:ARG:O	1:B:408:GLN:HG2	1.91	0.70
1:B:282:PHE:C	5:B:2221:HOH:O	2.30	0.70
1:D:389:MET:HE2	1:D:436:ILE:CG2	2.20	0.70
1:D:262:VAL:HG11	1:D:351:ILE:CG2	2.21	0.70
1:C:228:GLU:HG2	5:C:2246:HOH:O	1.91	0.70
1:D:408:GLN:NE2	5:D:2337:HOH:O	2.25	0.69
1:D:370:GLU:OE1	5:D:2298:HOH:O	2.10	0.69
1:D:375:GLN:NE2	1:D:421:ASN:H	1.91	0.68
5:A:2222:HOH:O	1:B:325:GLN:O	2.10	0.68
1:B:389:MET:HE3	1:B:436:ILE:HB	1.76	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:164:GLN:OE1	5:B:2139:HOH:O	2.12	0.67
1:C:43:TRP:HE1	1:C:107:HIS:HD2	1.40	0.67
1:B:57:ILE:HG12	1:B:105:ILE:HD12	1.77	0.67
1:C:43:TRP:HE1	1:C:107:HIS:CD2	2.13	0.67
1:B:299:MET:CE	5:C:2119:HOH:O	2.42	0.66
1:D:389:MET:HE1	1:D:435:PRO:HD2	1.77	0.66
1:C:250:GLU:CD	5:C:2266:HOH:O	2.34	0.66
1:B:111:THR:HG23	5:B:2098:HOH:O	1.97	0.65
1:A:343:GLU:OE2	1:D:343:GLU:OE2	2.15	0.65
1:C:68:ASN:ND2	1:C:71:ASN:HD22	1.95	0.65
1:C:372:ARG:HA	1:C:375:GLN:CG	2.27	0.65
1:B:302:GLU:OE1	1:D:159:HIS:HD2	1.80	0.65
1:A:264:LYS:CE	5:A:2222:HOH:O	2.46	0.64
1:C:207:SER:OG	5:C:2220:HOH:O	1.88	0.64
1:A:198:VAL:HG13	1:A:226:VAL:HG21	1.79	0.63
1:D:408:GLN:HA	1:D:408:GLN:HE21	1.61	0.63
1:B:365:TYR:OH	5:B:2275:HOH:O	2.14	0.63
1:D:408:GLN:NE2	5:D:2336:HOH:O	2.30	0.63
1:B:141:ARG:CZ	1:B:360:GLU:OE2	2.48	0.62
1:C:89:MET:HG3	5:C:2016:HOH:O	1.99	0.62
1:A:302:GLU:OE1	1:C:159:HIS:HD2	1.82	0.62
1:B:389:MET:CE	1:B:436:ILE:HB	2.29	0.62
1:A:144:ASP:O	1:A:148:GLU:HG3	1.99	0.62
1:A:118:ASN:HD21	1:A:197:GLY:H	1.46	0.61
1:B:118:ASN:HD21	1:B:197:GLY:H	1.47	0.61
1:C:352:LEU:O	1:C:356:GLN:HG3	1.99	0.61
1:A:43:TRP:HE1	1:A:107:HIS:CD2	2.18	0.61
1:C:21:TYR:HB3	1:C:350:THR:HG21	1.81	0.61
1:C:118:ASN:HD21	1:C:197:GLY:H	1.48	0.61
1:C:372:ARG:O	1:C:375:GLN:HG3	2.00	0.61
1:A:82:LYS:HZ1	1:A:85:ARG:CZ	2.14	0.61
1:A:181:MET:CE	1:B:246:LYS:HD3	2.31	0.60
1:A:43:TRP:HE1	1:A:107:HIS:HD2	1.50	0.60
1:D:389:MET:CE	1:D:436:ILE:CG1	2.78	0.60
1:D:118:ASN:HD21	1:D:197:GLY:H	1.47	0.59
1:D:300:ARG:HH11	1:D:357:ASN:ND2	2.00	0.59
1:B:159:HIS:HD2	1:D:302:GLU:OE1	1.84	0.59
1:A:17:LEU:CD1	1:A:343:GLU:HG2	2.32	0.59
1:D:300:ARG:HH11	1:D:357:ASN:HD21	1.51	0.59
1:B:43:TRP:HE1	1:B:107:HIS:CD2	2.20	0.59
1:C:68:ASN:HD22	1:C:71:ASN:HD22	1.51	0.57
1:B:117:ASP:OD1	5:B:2101:HOH:O	2.18	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:37:ARG:HH12	1:C:41:GLN:HE21	1.52	0.56
1:D:246:LYS:O	1:D:250:GLU:HG2	2.05	0.56
1:D:389:MET:HE3	1:D:436:ILE:CG1	2.34	0.56
1:A:398:ASP:O	1:A:401:GLU:HG2	2.05	0.56
1:A:181:MET:HE1	1:B:246:LYS:HD3	1.87	0.56
1:D:470:LYS:HB3	1:D:471:PRO:HD3	1.88	0.55
1:A:372:ARG:HA	1:A:375:GLN:HG2	1.88	0.55
1:B:389:MET:HE2	1:B:436:ILE:HD13	1.81	0.55
1:C:185:ASN:HD22	1:C:188[A]:ARG:HH12	1.55	0.55
1:D:370:GLU:HG3	5:D:2293:HOH:O	2.07	0.54
1:B:328:GLU:O	1:B:329:ARG:HB2	2.07	0.54
1:D:389:MET:HE2	1:D:436:ILE:CG1	2.37	0.54
1:A:375:GLN:HG3	1:A:376:GLU:HG2	1.90	0.54
1:D:83:ARG:NH2	1:D:84:LEU:HD21	2.22	0.54
1:D:262:VAL:HG11	1:D:351:ILE:HG21	1.87	0.54
1:B:299:MET:HE3	2:B:1000:AMP:H2'	1.90	0.53
1:C:372:ARG:HA	1:C:375:GLN:HG2	1.91	0.53
1:C:16:PRO:HB2	1:C:21:TYR:CE2	2.44	0.53
1:C:137:ARG:HH11	1:C:356:GLN:HE22	1.55	0.52
1:B:420:ASP:OD1	1:D:371:ARG:NH2	2.40	0.52
5:A:2222:HOH:O	1:B:325:GLN:C	2.46	0.52
1:B:303:ARG:NH2	5:B:2240:HOH:O	2.42	0.52
1:B:96:GLY:O	1:B:99:CYS:O	2.28	0.51
1:C:389:MET:SD	1:C:436:ILE:HD13	2.50	0.51
1:B:43:TRP:HE1	1:B:107:HIS:HD2	1.59	0.51
1:A:194:ARG:HD3	1:A:234:ARG:CD	2.41	0.51
1:D:466:TYR:OH	5:D:2396:HOH:O	2.19	0.50
1:B:389:MET:HE3	1:B:436:ILE:CG1	2.42	0.50
1:C:375:GLN:HB3	5:C:2341:HOH:O	2.12	0.49
1:C:371:ARG:O	1:C:375:GLN:HG2	2.12	0.49
1:B:295:LYS:HE3	1:B:297:ASN:OD1	2.13	0.49
1:B:34:TYR:OH	1:B:126:ASN:ND2	2.29	0.49
1:C:187:LYS:HE2	5:C:2207:HOH:O	2.11	0.49
1:A:11:ASP:OD1	1:D:33:ARG:HD2	2.13	0.49
1:D:389:MET:HE2	1:D:436:ILE:HG13	1.94	0.48
1:D:241:GLN:HB3	1:D:329:ARG:HG2	1.95	0.48
1:A:178:ASP:HA	1:A:181:MET:HE2	1.93	0.48
1:B:137:ARG:HD2	1:B:356:GLN:HE21	1.79	0.48
1:B:292:MET:O	1:B:292:MET:HG2	2.12	0.48
1:C:470:LYS:HB3	1:C:471:PRO:HD3	1.95	0.48
1:A:262:VAL:HG11	1:A:351:ILE:HG23	1.95	0.48
1:B:52:THR:HB	1:B:225:MET:HE1	1.96	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:107:HIS:HE1	5:C:2146:HOH:O	1.97	0.47
1:C:141:ARG:NH2	1:C:356:GLN:HB3	2.30	0.47
1:B:389:MET:HE3	1:B:436:ILE:CB	2.43	0.47
1:B:137:ARG:HD2	1:B:356:GLN:NE2	2.30	0.47
1:A:159:HIS:HD2	1:C:302:GLU:OE1	1.98	0.47
1:C:194:ARG:HD3	1:C:234:ARG:CD	2.45	0.47
1:B:299:MET:HB3	5:B:2094:HOH:O	2.14	0.47
1:A:408:GLN:HE21	1:A:408:GLN:N	2.13	0.47
1:A:390:VAL:HG21	1:A:396:ARG:HH11	1.79	0.47
1:B:328:GLU:O	1:B:329:ARG:CB	2.61	0.47
1:D:331:LEU:CD1	5:D:2265:HOH:O	2.36	0.47
1:D:389:MET:HE1	1:D:435:PRO:CD	2.43	0.47
1:A:141:ARG:CZ	1:A:360:GLU:OE2	2.62	0.47
1:A:264:LYS:HE3	5:A:2146:HOH:O	2.15	0.47
1:A:198:VAL:HG22	1:A:223:ASP:HA	1.97	0.47
1:B:141:ARG:NH1	1:B:360:GLU:OE2	2.48	0.47
1:A:262:VAL:HG11	1:A:351:ILE:CG2	2.45	0.47
1:D:375:GLN:HE22	1:D:421:ASN:N	2.01	0.46
1:C:424:ILE:HG13	1:C:444:LEU:HD13	1.98	0.46
1:A:15:SER:HB2	1:D:15:SER:HB2	1.97	0.46
1:D:262:VAL:HG11	1:D:351:ILE:HG23	1.94	0.46
1:D:43:TRP:HE1	1:D:107:HIS:CD2	2.34	0.46
1:B:329:ARG:HG2	1:B:330:THR:N	2.30	0.46
5:B:2373:HOH:O	1:D:159:HIS:HE1	1.99	0.46
1:C:325:GLN:O	5:C:2299:HOH:O	2.20	0.45
1:B:371:ARG:HD2	1:D:418:GLY:O	2.17	0.45
1:B:64:GLU:OE2	1:B:99:CYS:C	2.55	0.45
1:D:132:LEU:HB2	1:D:133:PRO:HD3	1.98	0.45
1:B:434:SER:N	1:B:435:PRO:CD	2.80	0.45
1:B:389:MET:HE2	1:B:435:PRO:HB2	1.98	0.45
1:B:418:GLY:O	1:D:371:ARG:HD2	2.17	0.45
1:C:50:GLU:HB3	1:C:55:LEU:HD12	1.99	0.45
1:A:17:LEU:HD12	1:A:343:GLU:HG2	1.98	0.45
1:A:194:ARG:HD3	1:A:234:ARG:HD3	1.99	0.45
1:B:182:ASP:O	1:B:186:LEU:HD13	2.15	0.45
1:C:21:TYR:HB3	1:C:350:THR:CG2	2.47	0.45
1:A:141:ARG:CZ	5:A:2123:HOH:O	2.64	0.45
1:B:86:HIS:CE1	1:B:89:MET:HG2	2.52	0.45
1:D:299:MET:HE1	2:D:1000:AMP:H2'	1.99	0.44
1:B:292:MET:N	1:B:293:PRO:CD	2.81	0.44
1:C:389:MET:HG2	1:C:399:CYS:SG	2.58	0.44
1:A:89:MET:O	1:A:93:HIS:HD2	2.00	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:88:VAL:O	1:D:92:VAL:HG23	2.18	0.44
1:D:185:ASN:HD22	1:D:188[A]:ARG:HH12	1.65	0.44
1:B:246:LYS:O	1:B:250:GLU:HG2	2.18	0.44
1:C:389:MET:HA	1:C:436:ILE:HD11	2.00	0.44
1:A:192:ASP:OD2	5:A:2160:HOH:O	2.21	0.44
1:A:375:GLN:NE2	5:A:2288:HOH:O	2.51	0.44
1:A:151:SER:O	1:A:153:PRO:HD3	2.18	0.44
1:C:372:ARG:HA	1:C:375:GLN:CD	2.38	0.43
1:B:108:LEU:HG	1:B:211:LEU:HD11	1.99	0.43
1:C:424:ILE:CD1	1:C:444:LEU:HD11	2.48	0.43
1:D:374:ARG:NH2	5:D:2305:HOH:O	2.44	0.43
5:B:2373:HOH:O	1:D:159:HIS:CE1	2.71	0.43
1:D:107:HIS:CD2	1:D:110:ALA:HB3	2.53	0.43
1:C:37:ARG:HH12	1:C:41:GLN:NE2	2.15	0.43
1:B:137:ARG:HH11	1:B:356:GLN:NE2	2.09	0.43
1:B:85:ARG:NH2	5:B:2076:HOH:O	2.51	0.42
2:A:1000:AMP:C8	1:D:331:LEU:HD22	2.54	0.42
1:A:68:ASN:ND2	1:A:71:ASN:HD22	2.17	0.42
1:C:76:MET:HE3	5:C:2096:HOH:O	2.19	0.42
1:B:343:GLU:OE1	1:C:343:GLU:OE2	2.38	0.42
1:D:434:SER:N	1:D:435:PRO:CD	2.82	0.42
1:D:371:ARG:HG3	5:D:2308:HOH:O	2.18	0.42
1:D:207:SER:OG	5:D:2185:HOH:O	2.14	0.42
1:A:278:MET:HA	1:A:363:VAL:O	2.19	0.42
1:D:425:GLU:O	1:D:429:VAL:HG13	2.20	0.42
1:D:34:TYR:HH	1:D:126:ASN:ND2	2.12	0.42
1:D:458:GLN:O	1:D:462:GLU:HG3	2.20	0.42
1:B:107:HIS:HE1	5:B:2099:HOH:O	2.01	0.42
1:A:118:ASN:ND2	1:A:196:ARG:HB3	2.34	0.42
1:D:299:MET:CE	2:D:1000:AMP:H2'	2.50	0.41
1:D:299:MET:HE1	2:D:1000:AMP:O2'	2.20	0.41
1:C:328:GLU:O	1:C:329:ARG:HB3	2.21	0.41
1:C:118:ASN:ND2	1:C:196:ARG:HB3	2.36	0.41
1:D:118:ASN:HD21	1:D:197:GLY:N	2.16	0.41
1:D:117:ASP:OD1	5:D:2084:HOH:O	2.20	0.41
1:C:37:ARG:NE	1:C:70:GLU:OE2	2.53	0.41
1:D:383:GLU:OE2	1:D:387:MET:HE2	2.21	0.41
1:B:12:SER:OG	5:B:2006:HOH:O	2.21	0.41
1:B:126:ASN:CB	5:B:2107:HOH:O	2.56	0.41
1:D:408:GLN:HA	1:D:408:GLN:NE2	2.30	0.41
1:C:194:ARG:HD3	1:C:234:ARG:HD3	2.02	0.41
1:A:259:GLY:O	1:A:308:ALA:HB1	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:328:GLU:O	1:D:329:ARG:HB3	2.21	0.41
1:D:383:GLU:HG3	1:D:387:MET:HE3	2.03	0.41
1:A:60:GLU:OE1	1:A:101:LYS:NZ	2.40	0.40
1:C:299[A]:MET:CE	2:C:1000:AMP:C8	3.02	0.40
1:B:278:MET:HA	1:B:363:VAL:O	2.21	0.40
1:D:47:ALA:HB2	5:D:2078:HOH:O	2.21	0.40
1:D:21:TYR:HB3	1:D:350:THR:HG21	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:A:2091:HOH:O	5:B:2124:HOH:O[3.555]	2.06	0.14

## 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	453/503 (90%)	447 (99%)	5 (1%)	1 (0%)	56	38
1	B	456/503 (91%)	450 (99%)	5 (1%)	1 (0%)	56	38
1	C	455/503 (90%)	449 (99%)	5 (1%)	1 (0%)	56	38
1	D	454/503 (90%)	447 (98%)	6 (1%)	1 (0%)	56	38
All	All	1818/2012 (90%)	1793 (99%)	21 (1%)	4 (0%)	56	38

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	329	ARG
1	B	329	ARG
1	C	329	ARG
1	D	329	ARG



### 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	403/438 (92%)	393 (98%)	10 (2%)	60	42
1	B	406/438 (93%)	396 (98%)	10 (2%)	60	42
1	C	402/438 (92%)	392 (98%)	10 (2%)	60	42
1	D	403/438 (92%)	390 (97%)	13 (3%)	51	32
All	All	1614/1752 (92%)	1571 (97%)	43 (3%)	59	39

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	133	PRO
1	A	190	ARG
1	A	228	GLU
1	A	264	LYS
1	A	329	ARG
1	A	360	GLU
1	A	396	ARG
1	A	408	GLN
1	A	458[A]	GLN
1	A	458[B]	GLN
1	B	6	ASP
1	B	190	ARG
1	B	211	LEU
1	B	228	GLU
1	B	292	MET
1	B	404	ARG
1	B	405	VAL
1	B	408	GLN
1	B	436	ILE
1	B	444	LEU
1	C	37	ARG
1	C	160	PHE
1	C	190	ARG
1	C	228	GLU
1	C	329	ARG
1	C	391	LYS

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Mol	Chain	Res	Type
1	C	396	ARG
1	C	412	SER
1	C	425	GLU
1	C	470	LYS
1	D	9	SER
1	D	117	ASP
1	D	160	PHE
1	D	210	GLN
1	D	228[A]	GLU
1	D	228[B]	GLU
1	D	329	ARG
1	D	402	LYS
1	D	404	ARG
1	D	405	VAL
1	D	408	GLN
1	D	444	LEU
1	D	473	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	ASN
1	A	93	HIS
1	A	107	HIS
1	A	118	ASN
1	A	126	ASN
1	A	159	HIS
1	A	177	GLN
1	A	185	ASN
1	A	408	GLN
1	A	439	GLN
1	B	68	ASN
1	B	107	HIS
1	B	118	ASN
1	B	159	HIS
1	B	177	GLN
1	B	185	ASN
1	B	356	GLN
1	B	408	GLN
1	C	7	HIS
1	C	41	GLN
1	C	68	ASN

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Mol	Chain	Res	Type
1	C	107	HIS
1	C	118	ASN
1	C	126	ASN
1	C	159	HIS
1	C	177	GLN
1	C	185	ASN
1	C	210	GLN
1	C	356	GLN
1	C	458	GLN
1	D	68	ASN
1	D	71	ASN
1	D	107	HIS
1	D	118	ASN
1	D	126	ASN
1	D	159	HIS
1	D	177	GLN
1	D	185	ASN
1	D	357	ASN
1	D	375	GLN
1	D	408	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 ⓘ

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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
2	AMP	A	1000	-	25,25,25	1.21	2 (8%)	38,38,38	1.99	9 (23%)
4	GOL	A	1475	-	5,5,5	0.32	0	5,5,5	0.59	0
2	AMP	B	1000	-	25,25,25	1.19	3 (12%)	38,38,38	1.96	8 (21%)
4	GOL	B	1482	-	5,5,5	0.34	0	5,5,5	0.66	0
2	AMP	C	1000	-	25,25,25	1.29	3 (12%)	38,38,38	1.96	10 (26%)
4	GOL	C	1473	-	5,5,5	0.45	0	5,5,5	0.36	0
2	AMP	D	1000	-	25,25,25	1.21	3 (12%)	38,38,38	1.85	10 (26%)
4	GOL	D	1474	-	5,5,5	0.63	0	5,5,5	0.54	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	AMP	A	1000	-	-	1/10/26/26	0/1/3/3
4	GOL	A	1475	-	-	0/4/4/4	0/0/0/0
2	AMP	B	1000	-	-	0/10/26/26	0/1/3/3
4	GOL	B	1482	-	-	0/4/4/4	0/0/0/0
2	AMP	C	1000	-	-	0/10/26/26	0/1/3/3
4	GOL	C	1473	-	-	0/4/4/4	0/0/0/0
2	AMP	D	1000	-	-	0/10/26/26	0/1/3/3
4	GOL	D	1474	-	-	0/4/4/4	0/0/0/0

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1000	AMP	C4-N9	-3.62	1.32	1.37
2	C	1000	AMP	C5-C4	3.56	1.48	1.40
2	D	1000	AMP	C4-N9	-3.42	1.32	1.37
2	A	1000	AMP	C5-C4	3.09	1.47	1.40
2	A	1000	AMP	C4-N9	-2.96	1.33	1.37
2	C	1000	AMP	C4-N9	-2.49	1.34	1.37
2	C	1000	AMP	O4'-C1'	2.48	1.45	1.41
2	D	1000	AMP	C2-N3	2.47	1.37	1.32
2	B	1000	AMP	C5-C4	2.32	1.45	1.40
2	B	1000	AMP	O4'-C1'	2.27	1.44	1.41
2	D	1000	AMP	C8-N9	-2.21	1.33	1.36

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1000	AMP	N3-C2-N1	-6.53	123.25	128.71
2	A	1000	AMP	N3-C2-N1	-6.41	123.35	128.71
2	B	1000	AMP	N3-C4-N9	5.72	135.76	125.43
2	B	1000	AMP	N3-C2-N1	-5.50	124.11	128.71
2	A	1000	AMP	N3-C4-N9	5.46	135.30	125.43
2	C	1000	AMP	N3-C4-N9	5.44	135.26	125.43
2	D	1000	AMP	N3-C4-N9	5.33	135.06	125.43
2	D	1000	AMP	C8-N9-C4	4.33	110.20	106.90
2	B	1000	AMP	C8-N9-C4	4.25	110.14	106.90
2	D	1000	AMP	N3-C2-N1	-3.84	125.50	128.71
2	A	1000	AMP	C8-N9-C4	3.66	109.69	106.90
2	B	1000	AMP	O3P-P-O5'	-3.47	97.07	106.65
2	C	1000	AMP	C8-N9-C4	3.36	109.46	106.90
2	D	1000	AMP	N6-C6-N1	3.29	125.83	119.36
2	D	1000	AMP	O4'-C1'-C2'	-3.10	102.02	106.77
2	C	1000	AMP	C4-C5-N7	-2.95	107.00	109.52
2	A	1000	AMP	C5-C4-N3	-2.94	119.30	125.70
2	C	1000	AMP	C5-C4-N3	-2.94	119.30	125.70
2	A	1000	AMP	O2P-P-O1P	2.89	119.89	110.44
2	D	1000	AMP	C1'-N9-C4	-2.82	121.77	126.64
2	B	1000	AMP	C5-C4-N3	-2.81	119.59	125.70
2	B	1000	AMP	O4'-C1'-C2'	-2.74	102.56	106.77
2	C	1000	AMP	O3P-P-O5'	-2.52	99.69	106.65
2	A	1000	AMP	C4-C5-N7	-2.45	107.42	109.52
2	B	1000	AMP	C1'-N9-C4	-2.40	122.49	126.64
2	A	1000	AMP	C1'-N9-C4	-2.36	122.55	126.64
2	A	1000	AMP	C2-N3-C4	2.28	120.49	114.01
2	D	1000	AMP	C5-C4-N3	-2.26	120.78	125.70
2	C	1000	AMP	O4'-C1'-N9	2.21	110.50	108.44
2	C	1000	AMP	C2-N3-C4	2.17	120.18	114.01
2	A	1000	AMP	C4'-O4'-C1'	-2.16	107.40	109.75
2	D	1000	AMP	C5-C6-N6	-2.15	115.86	120.72
2	C	1000	AMP	C1'-N9-C4	-2.11	122.99	126.64
2	D	1000	AMP	C5-C4-N9	-2.09	104.15	107.16
2	D	1000	AMP	O3P-P-O2P	2.05	115.61	107.61
2	C	1000	AMP	C2-N1-C6	2.05	122.47	118.77
2	B	1000	AMP	C2-N3-C4	2.03	119.80	114.01

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1000	AMP	C2'-C1'-N9-C8

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, 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	457/503 (90%)	0.23	33 (7%)	15 11	22, 28, 51, 72	0
1	B	463/503 (92%)	0.30	23 (4%)	28 22	23, 31, 49, 59	0
1	C	456/503 (90%)	0.31	34 (7%)	14 10	23, 28, 49, 62	0
1	D	456/503 (90%)	0.38	28 (6%)	21 15	23, 31, 53, 65	0
All	All	1832/2012 (91%)	0.30	118 (6%)	18 14	22, 29, 51, 72	0

All (118) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	63	ARG	5.3
1	D	100	PRO	5.3
1	A	5	GLY	5.0
1	C	397	GLN	5.0
1	C	431	ALA	4.9
1	D	93	HIS	4.9
1	D	82	LYS	4.8
1	C	294	TYR	4.8
1	A	404	ARG	4.7
1	C	396	ARG	4.7
1	C	438	SER	4.7
1	A	431	ALA	4.6
1	A	438	SER	4.6
1	D	85	ARG	4.6
1	A	429	VAL	4.5
1	D	5	GLY	4.4
1	A	432	TYR	4.4
1	D	101	LYS	4.2
1	C	295	LYS	4.2
1	D	62	ILE	4.1
1	A	387	MET	4.1

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Mol	Chain	Res	Type	RSRZ
1	C	391	LYS	4.0
1	A	401	GLU	4.0
1	C	392	ALA	4.0
1	A	439	GLN	4.0
1	C	5	GLY	4.0
1	B	474	SER	4.0
1	C	393	GLY	4.0
1	A	408	GLN	3.7
1	B	97	HIS	3.7
1	C	441	ASP	3.7
1	B	104	GLY	3.7
1	B	63	ARG	3.7
1	B	393	GLY	3.6
1	A	19	SER	3.6
1	A	398	ASP	3.6
1	D	54	GLY	3.6
1	C	432	TYR	3.5
1	A	441	ASP	3.5
1	D	393	GLY	3.5
1	D	89	MET	3.5
1	A	397	GLN	3.4
1	D	105	ILE	3.4
1	A	396	ARG	3.4
1	A	390	VAL	3.3
1	C	436	ILE	3.3
1	D	58	THR	3.3
1	D	6	ASP	3.3
1	B	5	GLY	3.3
1	D	99	CYS	3.3
1	A	436	ILE	3.2
1	B	391	LYS	3.2
1	D	103	ALA	3.2
1	B	60	GLU	3.2
1	C	398	ASP	3.1
1	C	395	SER	3.0
1	A	434	SER	3.0
1	D	97	HIS	3.0
1	B	58	THR	3.0
1	A	6	ASP	3.0
1	D	57	ILE	3.0
1	B	67	SER	3.0
1	A	442	HIS	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	67	SER	2.9
1	A	282	PHE	2.8
1	D	96	GLY	2.8
1	A	427	ILE	2.8
1	C	408	GLN	2.7
1	A	435	PRO	2.7
1	C	404	ARG	2.7
1	C	442	HIS	2.7
1	A	474	SER	2.7
1	D	79	GLU	2.7
1	B	431	ALA	2.7
1	D	104	GLY	2.7
1	C	400	HIS	2.7
1	B	401	GLU	2.7
1	B	89	MET	2.6
1	B	395	SER	2.6
1	B	473	GLU	2.6
1	C	425	GLU	2.6
1	A	393	GLY	2.6
1	C	387	MET	2.6
1	C	437	HIS	2.6
1	B	102	ALA	2.5
1	A	405	VAL	2.5
1	C	19	SER	2.5
1	C	434	SER	2.5
1	C	6	ASP	2.4
1	D	282	PHE	2.4
1	C	428	GLN	2.4
1	C	439	GLN	2.4
1	A	383	GLU	2.3
1	D	55	LEU	2.3
1	A	473	GLU	2.3
1	B	96	GLY	2.3
1	B	6	ASP	2.3
1	B	62	ILE	2.3
1	C	424	ILE	2.3
1	D	71	ASN	2.3
1	C	405	VAL	2.2
1	A	400	HIS	2.2
1	B	397	GLN	2.2
1	A	412	SER	2.2
1	C	394	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	390	VAL	2.2
1	C	386	ILE	2.2
1	C	435	PRO	2.2
1	A	399	CYS	2.2
1	B	398	ASP	2.2
1	B	54	GLY	2.1
1	D	391	LYS	2.1
1	D	106	ILE	2.1
1	C	401	GLU	2.1
1	A	402	LYS	2.1
1	B	217	HIS	2.1
1	C	419	GLY	2.0
1	A	403	ILE	2.0

## 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, 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	GOL	A	1475	6/6	0.14	3.18	21,23,26,27	0
2	AMP	C	1000	23/23	0.26	2.91	35,46,47,49	0
4	GOL	D	1474	6/6	0.16	2.44	21,25,28,28	0
4	GOL	B	1482	6/6	0.14	1.92	35,36,37,38	0
2	AMP	A	1000	23/23	0.15	0.19	26,35,38,40	0
3	CL	A	1001	1/1	0.08	-0.72	17,17,17,17	0
4	GOL	C	1473	6/6	0.09	-1.20	17,23,24,25	0
2	AMP	D	1000	23/23	0.07	-1.40	13,16,20,21	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	AMP	B	1000	23/23	0.08	-1.61	12,14,18,20	0
3	CL	D	1001	1/1	0.06	-1.68	29,29,29,29	0
3	CL	C	1001	1/1	0.06	-2.94	13,13,13,13	0
3	CL	B	1001	1/1	0.03	-5.50	27,27,27,27	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.