



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:22 AM GMT

PDB ID : 2A2D
Title : X-ray structure of human n-acetyl galactosamine kinase complexed with Mn-AMPPNP and n-acetyl galactosamine
Authors : Thoden, J.B.; Holden, H.M.
Deposited on : 2005-06-22
Resolution : 2.20 Å(reported)

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

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

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

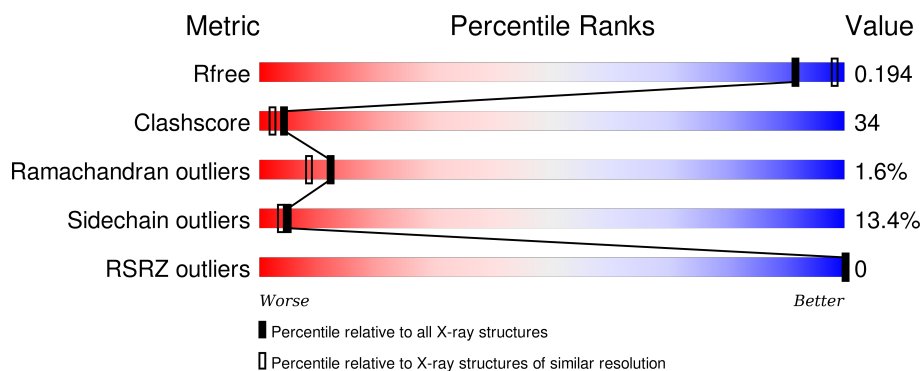
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

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



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

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	478	

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
4	ANP	A	461	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3765 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.

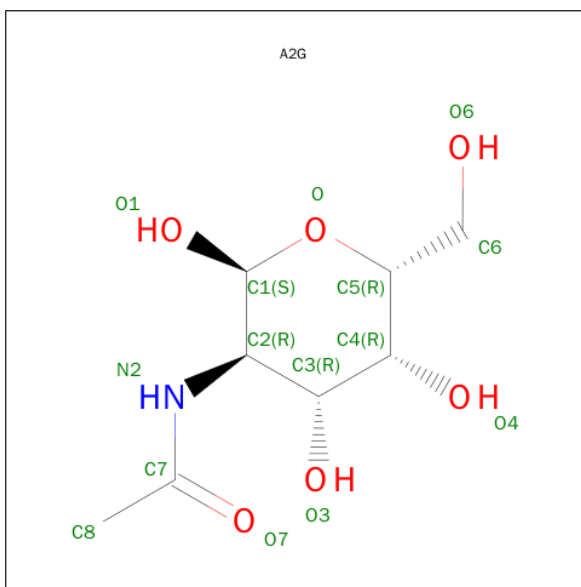
- Molecule 1 is a protein called N-acetylgalactosamine kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	448	3462	2198	587	649	28	0	3	0

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	CLONING ARTIFACT	UNP Q01415
A	-18	GLY	-	CLONING ARTIFACT	UNP Q01415
A	-17	SER	-	CLONING ARTIFACT	UNP Q01415
A	-16	SER	-	CLONING ARTIFACT	UNP Q01415
A	-15	HIS	-	EXPRESSION TAG	UNP Q01415
A	-14	HIS	-	EXPRESSION TAG	UNP Q01415
A	-13	HIS	-	EXPRESSION TAG	UNP Q01415
A	-12	HIS	-	EXPRESSION TAG	UNP Q01415
A	-11	HIS	-	EXPRESSION TAG	UNP Q01415
A	-10	HIS	-	EXPRESSION TAG	UNP Q01415
A	-9	SER	-	CLONING ARTIFACT	UNP Q01415
A	-8	SER	-	CLONING ARTIFACT	UNP Q01415
A	-7	GLU	-	CLONING ARTIFACT	UNP Q01415
A	-6	ASN	-	CLONING ARTIFACT	UNP Q01415
A	-5	LEU	-	CLONING ARTIFACT	UNP Q01415
A	-4	TYR	-	CLONING ARTIFACT	UNP Q01415
A	-3	PHE	-	CLONING ARTIFACT	UNP Q01415
A	-2	GLN	-	CLONING ARTIFACT	UNP Q01415
A	-1	GLY	-	CLONING ARTIFACT	UNP Q01415
A	0	HIS	-	CLONING ARTIFACT	UNP Q01415

- Molecule 2 is SUGAR (N-ACETYL-D-GALACTOSAMINE) (three-letter code: A2G) (formula: C₈H₁₅NO₆).

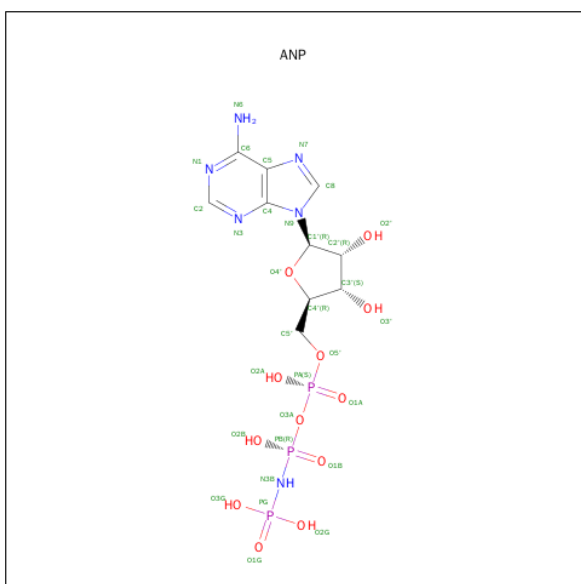


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			15	8	1	6		

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mn	0	0
			1	1		

- Molecule 4 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



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

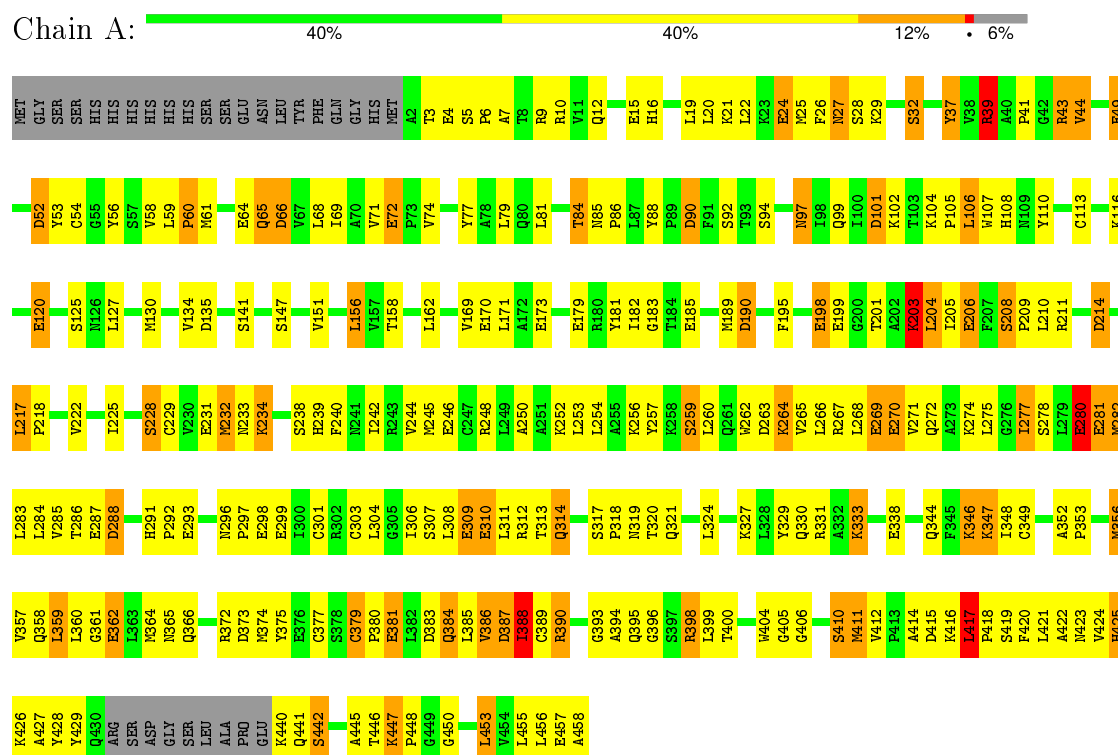
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	256	Total	O	0	0
			256	256		

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: N-acetylgalactosamine kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	119.10 Å 119.10 Å 65.90 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.20 29.78 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.9 (30.00-2.20) 98.9 (29.78-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.20 Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.203 , 0.247 0.199 , 0.194	Depositor DCC
R_{free} test set	2691 reflections (11.14%)	DCC
Wilson B-factor (Å ²)	2.6	Xtriage
Anisotropy	0.914	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 260.0	EDS
Estimated twinning fraction	0.236 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.29$, $\langle L^2 \rangle = 0.12$	Xtriage
Outliers	0 of 26904 reflections	Xtriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	3765	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.65% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MN, ANP, A2G

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	1.01	27/3538 (0.8%)	1.43	34/4786 (0.7%)

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	338	GLU	CD-OE2	7.20	1.33	1.25
1	A	309	GLU	CD-OE2	6.99	1.33	1.25
1	A	185	GLU	CD-OE2	6.63	1.32	1.25
1	A	24	GLU	CD-OE2	6.57	1.32	1.25
1	A	280	GLU	CD-OE2	6.28	1.32	1.25
1	A	270	GLU	CD-OE2	6.23	1.32	1.25
1	A	287	GLU	CD-OE2	6.22	1.32	1.25
1	A	362	GLU	CD-OE2	6.12	1.32	1.25
1	A	64	GLU	CD-OE2	6.02	1.32	1.25
1	A	120	GLU	CD-OE2	5.90	1.32	1.25
1	A	4	GLU	CD-OE2	5.71	1.31	1.25
1	A	72	GLU	CD-OE2	5.68	1.31	1.25
1	A	246	GLU	CD-OE2	5.68	1.31	1.25
1	A	298	GLU	CD-OE2	5.65	1.31	1.25
1	A	206	GLU	CD-OE2	5.55	1.31	1.25
1	A	299	GLU	CD-OE2	5.50	1.31	1.25
1	A	231	GLU	CD-OE2	5.50	1.31	1.25
1	A	269	GLU	CD-OE2	5.48	1.31	1.25
1	A	281	GLU	CD-OE2	5.46	1.31	1.25
1	A	381	GLU	CD-OE2	5.34	1.31	1.25
1	A	199	GLU	CD-OE2	5.28	1.31	1.25
1	A	457	GLU	CD-OE2	5.21	1.31	1.25
1	A	310	GLU	CD-OE2	5.19	1.31	1.25
1	A	173	GLU	CD-OE2	5.19	1.31	1.25
1	A	198	GLU	CD-OE2	5.14	1.31	1.25
1	A	49	GLU	CD-OE2	5.12	1.31	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	179	GLU	CD-OE2	5.06	1.31	1.25

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	387	ASP	CB-CG-OD1	8.85	126.27	118.30
1	A	390	ARG	NE-CZ-NH2	-8.71	115.95	120.30
1	A	415	ASP	CB-CA-C	-8.71	92.98	110.40
1	A	415	ASP	CB-CG-OD2	-8.26	110.86	118.30
1	A	329	TYR	CB-CG-CD2	7.79	125.67	121.00
1	A	288	ASP	CB-CG-OD2	-7.76	111.32	118.30
1	A	387	ASP	CB-CG-OD2	-7.67	111.40	118.30
1	A	263	ASP	CB-CG-OD2	-7.43	111.61	118.30
1	A	453	LEU	CB-CA-C	-7.22	96.48	110.20
1	A	214	ASP	CB-CG-OD1	7.20	124.78	118.30
1	A	90	ASP	CB-CG-OD2	-7.07	111.94	118.30
1	A	390	ARG	NE-CZ-NH1	6.98	123.79	120.30
1	A	263	ASP	CB-CG-OD1	6.83	124.44	118.30
1	A	101	ASP	CB-CG-OD2	-6.70	112.27	118.30
1	A	383	ASP	CB-CG-OD2	-6.47	112.47	118.30
1	A	288	ASP	CB-CG-OD1	6.31	123.98	118.30
1	A	372	ARG	NE-CZ-NH2	-6.22	117.19	120.30
1	A	39	ARG	NE-CZ-NH1	6.16	123.38	120.30
1	A	214	ASP	CB-CG-OD2	-6.12	112.79	118.30
1	A	329	TYR	CB-CG-CD1	-6.11	117.33	121.00
1	A	52	ASP	CB-CG-OD2	-5.89	113.00	118.30
1	A	398	ARG	NE-CZ-NH2	-5.88	117.36	120.30
1	A	101	ASP	CB-CG-OD1	5.76	123.48	118.30
1	A	373	ASP	CB-CG-OD2	-5.74	113.14	118.30
1	A	415	ASP	CB-CG-OD1	5.73	123.45	118.30
1	A	151	VAL	CA-CB-CG1	-5.68	102.38	110.90
1	A	190	ASP	CB-CG-OD2	-5.47	113.38	118.30
1	A	135	ASP	CB-CG-OD1	5.34	123.11	118.30
1	A	66	ASP	CB-CG-OD1	5.33	123.10	118.30
1	A	208	SER	C-N-CD	-5.30	108.95	120.60
1	A	84	THR	CA-CB-CG2	-5.23	105.07	112.40
1	A	405	GLY	C-N-CA	-5.20	111.38	122.30
1	A	379	CYS	N-CA-CB	5.09	119.77	110.60
1	A	203	LYS	N-CA-CB	5.02	119.64	110.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	3462	0	3503	236	0
2	A	15	0	14	0	0
3	A	1	0	0	0	0
4	A	31	0	13	9	0
5	A	256	0	0	13	0
All	All	3765	0	3530	236	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 34.

All (236) 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:A:357:VAL:HG12	1:A:395:GLN:HG2	1.25	1.14
1:A:390:ARG:HA	1:A:394:ALA:HB3	1.32	1.09
1:A:447:LYS:HG3	1:A:448:PRO:HD2	1.40	1.02
1:A:65:GLN:H	1:A:65:GLN:HE21	1.12	0.97
1:A:218:PRO:HG2	1:A:411:MET:CE	1.96	0.94
1:A:264:LYS:HD3	1:A:274:LYS:NZ	1.88	0.88
1:A:417:LEU:HB3	1:A:418:PRO:HD3	1.56	0.88
1:A:301:CYS:SG	1:A:308:LEU:HD23	2.16	0.85
1:A:417:LEU:HD12	1:A:421:LEU:HG	1.58	0.83
1:A:447:LYS:HG3	1:A:448:PRO:CD	2.08	0.83
1:A:234:LYS:H	1:A:234:LYS:HD3	1.45	0.82
1:A:271:VAL:HG12	1:A:282:MET:CE	2.09	0.81
1:A:7:ALA:HB3	1:A:453:LEU:CD2	2.10	0.80
1:A:352:ALA:HA	1:A:356:MET:SD	2.23	0.77
1:A:390:ARG:CA	1:A:394:ALA:HB3	2.11	0.77
1:A:278:SER:OG	1:A:281:GLU:HG3	1.83	0.77
1:A:264:LYS:HD3	1:A:274:LYS:HZ3	1.49	0.77
1:A:271:VAL:HG12	1:A:282:MET:HE3	1.67	0.76
1:A:318:PRO:HA	1:A:321:GLN:OE1	1.86	0.76
1:A:317:SER:O	1:A:321:GLN:HG3	1.88	0.74
1:A:306:ILE:CG2	1:A:310:GLU:HB2	2.18	0.74
1:A:357:VAL:CG1	1:A:395:GLN:HG2	2.12	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:418:PRO:HD3	5:A:532:HOH:O	1.88	0.73
1:A:330:GLN:HG3	1:A:374:MET:HE1	1.70	0.72
1:A:379:CYS:HB2	1:A:380:PRO:HD2	1.72	0.71
1:A:445:ALA:HB3	5:A:542:HOH:O	1.91	0.71
1:A:234:LYS:N	1:A:234:LYS:HD3	2.05	0.70
1:A:7:ALA:HB3	1:A:453:LEU:HD21	1.73	0.70
1:A:357:VAL:HG12	1:A:395:GLN:CG	2.16	0.70
1:A:44:VAL:HG11	1:A:225:ILE:HD11	1.74	0.70
1:A:393:GLY:HA2	1:A:416:LYS:HD3	1.75	0.68
1:A:218:PRO:HG2	1:A:411:MET:HE2	1.76	0.67
1:A:393:GLY:CA	1:A:416:LYS:HD3	2.24	0.66
1:A:422:ALA:O	1:A:426:LYS:HB2	1.94	0.66
1:A:346:LYS:O	1:A:349:CYS:N	2.28	0.66
1:A:198:GLU:HB2	1:A:203:LYS:HE3	1.78	0.66
1:A:233:ASN:ND2	4:A:461:ANP:O1A	2.30	0.65
1:A:44:VAL:CG1	1:A:225:ILE:HD11	2.27	0.64
1:A:387:ASP:O	1:A:390:ARG:N	2.31	0.64
1:A:352:ALA:N	1:A:353:PRO:HD3	2.12	0.64
1:A:420:PHE:O	1:A:424:VAL:HG23	1.97	0.64
1:A:271:VAL:HG12	1:A:282:MET:HE1	1.80	0.64
1:A:307:SER:N	1:A:310:GLU:OE1	2.31	0.64
1:A:9:ARG:HB3	1:A:455:LEU:HD23	1.80	0.64
1:A:275:LEU:HD12	1:A:282:MET:CE	2.28	0.64
1:A:275:LEU:HD12	1:A:282:MET:HE2	1.80	0.64
1:A:20:LEU:O	1:A:24:GLU:HG3	1.98	0.64
1:A:362:GLU:O	1:A:366:GLN:HG3	1.98	0.63
1:A:307:SER:OG	1:A:310:GLU:HG3	1.99	0.63
1:A:232:MET:SD	1:A:234:LYS:HD2	2.39	0.63
1:A:417:LEU:HD11	1:A:421:LEU:HD11	1.80	0.62
1:A:234:LYS:HE3	4:A:461:ANP:PG	2.39	0.62
1:A:420:PHE:HD2	1:A:421:LEU:HD23	1.65	0.62
1:A:104:LYS:O	1:A:106:LEU:HD12	2.01	0.61
1:A:440:LYS:N	5:A:541:HOH:O	2.32	0.61
1:A:365:ASN:OD1	1:A:390:ARG:NE	2.30	0.61
1:A:417:LEU:O	1:A:421:LEU:HG	2.01	0.61
1:A:222:VAL:HG21	1:A:417:LEU:HD22	1.82	0.61
1:A:65:GLN:H	1:A:65:GLN:NE2	1.92	0.60
1:A:414:ALA:HB3	5:A:710:HOH:O	2.01	0.60
1:A:272:GLN:HG2	1:A:272:GLN:O	2.02	0.60
1:A:361:GLY:O	1:A:365:ASN:ND2	2.34	0.60
1:A:310:GLU:HA	1:A:314:GLN:HG3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:LYS:HD3	1:A:274:LYS:HZ1	1.65	0.59
1:A:44:VAL:HG23	1:A:61:MET:HG2	1.84	0.59
1:A:16:HIS:HB2	1:A:19:LEU:HB2	1.85	0.59
1:A:218:PRO:HG2	1:A:411:MET:HE1	1.83	0.58
1:A:425:HIS:O	1:A:429:TYR:HB2	2.02	0.58
1:A:417:LEU:HB3	5:A:532:HOH:O	2.02	0.58
1:A:189:MET:HE3	1:A:190:ASP:OD1	2.03	0.58
1:A:417:LEU:HB3	1:A:418:PRO:CD	2.32	0.58
1:A:379:CYS:HB2	1:A:380:PRO:CD	2.33	0.58
1:A:25:MET:HE3	5:A:674:HOH:O	2.04	0.58
1:A:306:ILE:HG22	1:A:307:SER:H	1.69	0.57
1:A:102:LYS:HE2	1:A:183:GLY:HA2	1.87	0.57
1:A:419:SER:O	1:A:423:ASN:N	2.28	0.56
1:A:421:LEU:O	1:A:425:HIS:HB2	2.05	0.56
1:A:107:TRP:CE3	4:A:461:ANP:H2	2.39	0.56
1:A:84:THR:HG23	1:A:134:VAL:O	2.06	0.56
1:A:306:ILE:HG22	1:A:307:SER:N	2.20	0.56
1:A:417:LEU:HD12	1:A:421:LEU:CG	2.34	0.56
1:A:306:ILE:HG23	1:A:310:GLU:HB2	1.86	0.56
1:A:386:VAL:O	1:A:390:ARG:HG3	2.06	0.56
1:A:234:LYS:HE3	4:A:461:ANP:N3B	2.21	0.56
1:A:12:GLN:NE2	1:A:458:ALA:O	2.36	0.56
1:A:417:LEU:O	1:A:420:PHE:HB3	2.05	0.55
1:A:353:PRO:HD2	1:A:356:MET:HG3	1.89	0.55
1:A:304:LEU:HB2	1:A:306:ILE:HD12	1.89	0.55
1:A:441:GLN:HG2	5:A:486:HOH:O	2.05	0.55
1:A:359:LEU:O	1:A:362:GLU:HB2	2.06	0.55
1:A:97:ASN:HA	5:A:701:HOH:O	2.05	0.55
1:A:381:GLU:N	1:A:381:GLU:OE1	2.30	0.55
1:A:107:TRP:HB2	4:A:461:ANP:N3	2.22	0.54
1:A:447:LYS:CG	1:A:448:PRO:HD2	2.28	0.54
1:A:304:LEU:HB2	1:A:306:ILE:CD1	2.38	0.54
1:A:240:PHE:O	1:A:244:VAL:HG23	2.07	0.54
1:A:267:ARG:O	1:A:270:GLU:HB2	2.07	0.54
1:A:65:GLN:HE21	1:A:65:GLN:N	1.94	0.54
1:A:39:ARG:HD3	1:A:39:ARG:C	2.27	0.54
1:A:346:LYS:O	1:A:347:LYS:C	2.46	0.54
1:A:206:GLU:O	1:A:210:LEU:HD12	2.07	0.54
1:A:358:GLN:HA	1:A:395:GLN:HE21	1.72	0.54
1:A:380:PRO:O	1:A:384:GLN:HG2	2.08	0.53
1:A:265:VAL:HA	1:A:270:GLU:OE1	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:310:GLU:O	1:A:314:GLN:HG3	2.09	0.53
1:A:54:CYS:HB3	1:A:268:LEU:HB2	1.90	0.53
1:A:110:TYR:CE1	1:A:182:ILE:HG12	2.44	0.52
1:A:416:LYS:O	1:A:417:LEU:C	2.47	0.52
1:A:446:THR:HG22	1:A:447:LYS:N	2.24	0.52
1:A:313:THR:HG22	1:A:314:GLN:N	2.25	0.51
1:A:25:MET:HE3	1:A:84:THR:HG22	1.91	0.51
1:A:85:ASN:HB3	1:A:88:TYR:HD1	1.75	0.51
1:A:116:LYS:O	1:A:120:GLU:HG3	2.11	0.51
1:A:280:GLU:CD	1:A:280:GLU:H	2.13	0.51
1:A:359:LEU:O	1:A:362:GLU:N	2.44	0.51
1:A:21:LYS:O	1:A:25:MET:HG3	2.11	0.50
1:A:56:TYR:CE1	1:A:269:GLU:HG2	2.46	0.50
1:A:102:LYS:CE	1:A:183:GLY:HA2	2.41	0.50
1:A:238:SER:O	1:A:239:HIS:C	2.49	0.50
1:A:385:LEU:HD13	1:A:428:TYR:CG	2.45	0.50
1:A:399:LEU:HD12	1:A:399:LEU:C	2.31	0.50
1:A:419:SER:O	1:A:420:PHE:C	2.49	0.50
1:A:310:GLU:CA	1:A:314:GLN:HG3	2.41	0.50
1:A:423:ASN:O	1:A:427:ALA:N	2.40	0.50
1:A:41:PRO:HA	1:A:66:ASP:HB3	1.94	0.50
1:A:388:ILE:HG22	1:A:389:CYS:N	2.27	0.49
1:A:281:GLU:HA	1:A:284:LEU:HD12	1.94	0.49
1:A:306:ILE:HG22	1:A:310:GLU:HB2	1.94	0.49
1:A:393:GLY:HA3	1:A:416:LYS:HD3	1.93	0.49
1:A:306:ILE:CG2	1:A:307:SER:H	2.25	0.49
1:A:238:SER:O	1:A:242:ILE:HG13	2.13	0.49
1:A:250:ALA:O	1:A:254:LEU:HD12	2.11	0.49
1:A:308:LEU:O	1:A:311:LEU:HB3	2.13	0.49
1:A:69:ILE:HG12	1:A:134:VAL:HG22	1.94	0.49
1:A:27:ASN:HA	1:A:32:SER:O	2.12	0.49
1:A:204:LEU:HD13	1:A:346:LYS:CD	2.43	0.48
1:A:37:TYR:N	1:A:37:TYR:CD1	2.80	0.48
1:A:16:HIS:CD2	1:A:16:HIS:N	2.82	0.48
1:A:296:ASN:HB2	1:A:297:PRO:HD2	1.95	0.48
1:A:357:VAL:O	1:A:358:GLN:C	2.51	0.48
1:A:256:LYS:HD3	1:A:262:TRP:CD1	2.47	0.48
1:A:225:ILE:HD12	1:A:446:THR:OG1	2.13	0.48
1:A:234:LYS:HE2	4:A:461:ANP:O3A	2.13	0.48
1:A:348:ILE:HG22	1:A:349:CYS:N	2.29	0.48
1:A:364:MET:HE2	1:A:396:GLY:HA3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:257:TYR:C	1:A:259:SER:H	2.17	0.48
1:A:5:SER:HB3	1:A:6:PRO:HD2	1.95	0.48
1:A:97:ASN:C	1:A:97:ASN:HD22	2.16	0.48
1:A:275:LEU:HB3	1:A:277:ILE:HG12	1.94	0.47
1:A:238:SER:HB3	1:A:319:ASN:HD22	1.79	0.47
1:A:256:LYS:HE3	1:A:303:CYS:O	2.14	0.47
1:A:208:SER:OG	1:A:208:SER:O	2.30	0.47
1:A:85:ASN:OD1	1:A:86:PRO:HD2	2.15	0.47
1:A:309:GLU:O	1:A:313:THR:HB	2.15	0.47
1:A:447:LYS:HE3	1:A:448:PRO:HD2	1.97	0.47
1:A:284:LEU:HB3	5:A:607:HOH:O	2.13	0.47
1:A:44:VAL:CG2	1:A:61:MET:HG2	2.45	0.47
1:A:218:PRO:HG3	1:A:360:LEU:CD1	2.44	0.47
1:A:278:SER:HG	1:A:281:GLU:HG3	1.78	0.47
1:A:204:LEU:HD13	1:A:346:LYS:HD2	1.95	0.47
1:A:104:LYS:O	1:A:106:LEU:N	2.47	0.47
1:A:171:LEU:HB3	1:A:195:PHE:CZ	2.50	0.47
1:A:26:PHE:HZ	1:A:72:GLU:HB3	1.78	0.47
1:A:113:CYS:SG	1:A:181:TYR:HB2	2.55	0.47
1:A:217:LEU:HD13	1:A:217:LEU:N	2.29	0.47
1:A:49:GLU:HB2	1:A:398:ARG:NH2	2.29	0.46
1:A:29:LYS:NZ	1:A:90:ASP:OD1	2.40	0.46
1:A:377:CYS:HA	1:A:404:TRP:CE2	2.51	0.46
1:A:10:ARG:HB3	1:A:458:ALA:HB3	1.98	0.46
1:A:217:LEU:N	1:A:217:LEU:CD1	2.79	0.46
1:A:245:MET:HE2	1:A:245:MET:CA	2.45	0.46
1:A:352:ALA:N	1:A:353:PRO:CD	2.78	0.46
1:A:266:LEU:HD23	1:A:266:LEU:HA	1.68	0.46
1:A:97:ASN:ND2	5:A:684:HOH:O	2.39	0.45
1:A:447:LYS:HG3	1:A:448:PRO:N	2.32	0.45
1:A:422:ALA:O	1:A:425:HIS:HB2	2.16	0.45
1:A:58:VAL:HG21	1:A:189:MET:HB2	1.98	0.45
1:A:228:SER:O	1:A:229:CYS:HB2	2.16	0.45
1:A:234:LYS:CE	4:A:461:ANP:N3B	2.80	0.45
1:A:208:SER:HA	1:A:209:PRO:HA	1.79	0.45
1:A:201:THR:O	1:A:448:PRO:HB3	2.17	0.44
1:A:43:ARG:HD2	1:A:43:ARG:C	2.37	0.44
1:A:25:MET:CE	1:A:84:THR:HG22	2.48	0.44
1:A:182:ILE:HA	1:A:182:ILE:HD12	1.72	0.44
1:A:22:LEU:CD2	1:A:68:LEU:HG	2.48	0.44
1:A:56:TYR:CE1	1:A:269:GLU:CG	3.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:399:LEU:O	1:A:399:LEU:HD12	2.16	0.44
1:A:248:ARG:O	1:A:252:LYS:HG3	2.17	0.44
1:A:447:LYS:HE3	5:A:641:HOH:O	2.18	0.44
1:A:156:LEU:HD12	1:A:156:LEU:O	2.18	0.44
1:A:446:THR:CG2	1:A:447:LYS:N	2.80	0.44
1:A:291:HIS:ND1	1:A:293:GLU:O	2.49	0.44
1:A:387:ASP:O	1:A:390:ARG:HB2	2.17	0.43
1:A:238:SER:HA	1:A:319:ASN:HD21	1.83	0.43
1:A:285:VAL:O	1:A:288:ASP:N	2.50	0.43
1:A:286:THR:HG21	1:A:333:LYS:HD3	2.00	0.43
1:A:417:LEU:CD1	1:A:421:LEU:HD11	2.45	0.43
1:A:71:VAL:HG12	1:A:72:GLU:N	2.32	0.43
1:A:266:LEU:N	1:A:270:GLU:OE1	2.42	0.43
1:A:310:GLU:O	1:A:314:GLN:N	2.51	0.43
1:A:74:VAL:HG22	1:A:130:MET:N	2.33	0.43
1:A:52:ASP:OD1	1:A:53:TYR:N	2.51	0.43
1:A:238:SER:HA	5:A:508:HOH:O	2.18	0.43
1:A:256:LYS:HD3	1:A:262:TRP:NE1	2.33	0.43
1:A:306:ILE:CG2	1:A:307:SER:N	2.80	0.43
1:A:41:PRO:HG3	1:A:450:GLY:O	2.17	0.43
1:A:327:LYS:HB3	1:A:330:GLN:OE1	2.18	0.43
1:A:412:VAL:HG21	1:A:420:PHE:CG	2.54	0.43
1:A:396:GLY:O	1:A:410:SER:HB3	2.19	0.43
1:A:292:PRO:O	1:A:293:GLU:C	2.54	0.43
1:A:364:MET:HE2	1:A:396:GLY:CA	2.49	0.43
1:A:286:THR:CG2	1:A:333:LYS:HD3	2.49	0.42
1:A:60:PRO:HD2	1:A:205:ILE:HD12	2.01	0.42
1:A:198:GLU:O	1:A:201:THR:HG22	2.20	0.42
1:A:79:LEU:HD11	1:A:81:LEU:HD21	2.01	0.42
1:A:387:ASP:O	1:A:388:ILE:C	2.57	0.42
1:A:280:GLU:O	1:A:283:LEU:HB2	2.20	0.42
1:A:77:TYR:O	1:A:127:LEU:HD23	2.20	0.41
1:A:277:ILE:HD12	1:A:282:MET:HG2	2.01	0.41
1:A:301:CYS:SG	1:A:308:LEU:HA	2.60	0.41
1:A:158:THR:O	1:A:162:LEU:HG	2.20	0.41
1:A:429:TYR:CE2	1:A:442:SER:HB2	2.56	0.41
1:A:234:LYS:HE2	4:A:461:ANP:PA	2.60	0.41
1:A:238:SER:HB3	1:A:319:ASN:ND2	2.36	0.41
1:A:228:SER:HB3	1:A:406:GLY:O	2.20	0.41
1:A:169:VAL:HG12	1:A:170:GLU:N	2.36	0.41
1:A:456:LEU:HD23	1:A:456:LEU:HA	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:PRO:HB3	5:A:477:HOH:O	2.20	0.41
1:A:77:TYR:O	1:A:94:SER:HA	2.21	0.40
1:A:147:SER:HB3	4:A:461:ANP:O5'	2.21	0.40
1:A:330:GLN:HA	1:A:374:MET:HE3	2.03	0.40
1:A:79:LEU:O	1:A:92:SER:HA	2.21	0.40
1:A:210:LEU:HA	1:A:210:LEU:HD12	1.84	0.40
1:A:277:ILE:CD1	1:A:282:MET:HG2	2.51	0.40
1:A:81:LEU:HD12	1:A:108:HIS:HB2	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	447/478 (94%)	399 (89%)	41 (9%)	7 (2%)	12 8

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	359	LEU
1	A	346	LYS
1	A	60	PRO
1	A	347	LYS
1	A	105	PRO
1	A	417	LEU
1	A	388	ILE

5.3.2 Protein sidechains [i](#)

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

resolution.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	383/407 (94%)	332 (87%)	51 (13%)	5 4

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	THR
1	A	15	GLU
1	A	27	ASN
1	A	28	SER
1	A	32	SER
1	A	37	TYR
1	A	39	ARG
1	A	43	ARG
1	A	44	VAL
1	A	59	LEU
1	A	65	GLN
1	A	97	ASN
1	A	99	GLN
1	A	101	ASP
1	A	106	LEU
1	A	125	SER
1	A	141	SER
1	A	156	LEU
1	A	203	LYS
1	A	204	LEU
1	A	211	ARG
1	A	214	ASP
1	A	217	LEU
1	A	228	SER
1	A	232	MET
1	A	234	LYS
1	A	253	LEU
1	A	259	SER
1	A	260	LEU
1	A	264	LYS
1	A	277	ILE
1	A	280	GLU
1	A	282	MET
1	A	312	ARG

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Mol	Chain	Res	Type
1	A	314	GLN
1	A	324	LEU
1	A	331	ARG
1	A	333	LYS
1	A	344	GLN
1	A	356	MET
1	A	375	TYR
1	A	384	GLN
1	A	386	VAL
1	A	388	ILE
1	A	400	THR
1	A	410	SER
1	A	411	MET
1	A	417	LEU
1	A	425	HIS
1	A	442	SER
1	A	447	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	HIS
1	A	27	ASN
1	A	65	GLN
1	A	97	ASN
1	A	233	ASN
1	A	366	GLN
1	A	395	GLN
1	A	423	ASN
1	A	425	HIS

5.3.3 RNA ⓘ

There are no RNA molecules 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 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	A2G	A	459	-	15,15,15	1.02	1 (6%)	17,21,21	1.97	4 (23%)
4	ANP	A	461	3	27,33,33	1.51	4 (14%)	30,52,52	2.19	7 (23%)

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	A2G	A	459	-	-	0/6/26/26	0/1/1/1
4	ANP	A	461	3	-	0/12/38/38	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	461	ANP	C6-N6	-3.03	1.25	1.34
4	A	461	ANP	O4'-C1'	-2.11	1.38	1.41
2	A	459	A2G	C1-C2	2.76	1.56	1.53
4	A	461	ANP	PG-O1G	3.14	1.49	1.46
4	A	461	ANP	PB-O1B	3.44	1.50	1.46

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	461	ANP	N3-C2-N1	-6.31	124.06	128.89
4	A	461	ANP	O1B-PB-N3B	-3.74	106.17	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	459	A2G	O3-C3-C2	-3.47	102.58	109.66
4	A	461	ANP	C2'-C1'-N9	-2.59	110.33	114.29
4	A	461	ANP	O3G-PG-O1G	-2.26	107.47	113.49
2	A	459	A2G	O7-C7-C8	-2.24	117.95	122.06
2	A	459	A2G	C8-C7-N2	2.05	120.03	116.11
4	A	461	ANP	O3A-PB-N3B	4.02	117.49	106.44
4	A	461	ANP	C2-N1-C6	4.32	126.49	118.77
4	A	461	ANP	O3A-PA-O5'	4.37	114.53	102.94
2	A	459	A2G	O-C5-C4	5.87	120.70	109.68

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	461	ANP	9	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	448/478 (93%)	-1.13	0 100 100	8, 26, 70, 94	0

There are no RSRZ outliers to report.

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	ANP	A	461	31/31	0.99	0.06	-0.32	6,26,43,67	0
2	A2G	A	459	15/15	0.99	0.05	-0.41	3,13,22,29	0
3	MN	A	460	1/1	1.00	0.03	-1.13	27,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.