



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:57 PM GMT

PDB ID : 3VN9
Title : Refined Crystal structure of non-phosphorylated MAP2K6 in a putative auto-inhibition state
Authors : Kinoshita, T.; Matsuzaka, H.; Nakai, R.; Kirii, Y.; Yokota, K.; Tada, T.; Matsumoto, T.
Deposited on : 2012-01-05
Resolution : 2.60 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

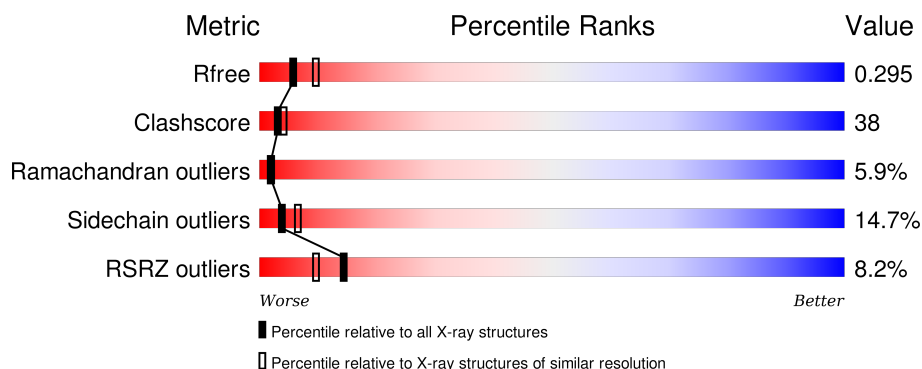
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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	340	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2322 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.

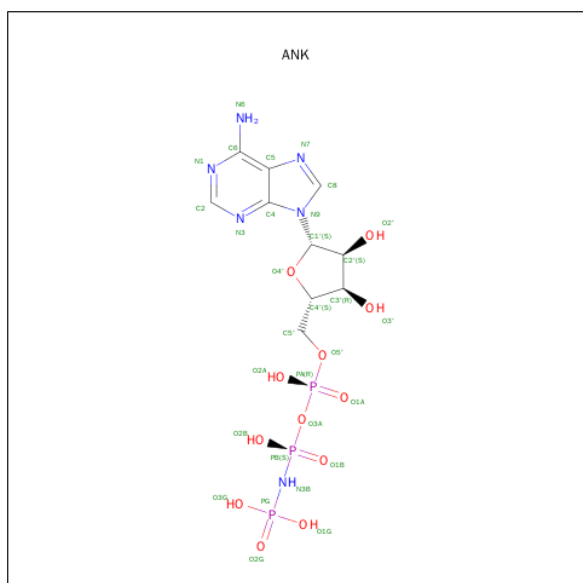
- Molecule 1 is a protein called Dual specificity mitogen-activated protein kinase kinase 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	291	2272	1460	375	421	16	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	335	HIS	-	EXPRESSION TAG	UNP P52564
A	336	HIS	-	EXPRESSION TAG	UNP P52564
A	337	HIS	-	EXPRESSION TAG	UNP P52564
A	338	HIS	-	EXPRESSION TAG	UNP P52564
A	339	HIS	-	EXPRESSION TAG	UNP P52564
A	340	HIS	-	EXPRESSION TAG	UNP P52564

- Molecule 2 is 9-{5-O-[(R)-HYDROXY{[(S)-HYDROXY(PHOSPHONOAMINO)PHOSPHORYL]OXY}PHOSPHORYL]-BETA-L-RIBOFURANOSYL}-9H-PURIN-6-AMINE (three-letter code: ANK) (formula: C₁₀H₁₇N₆O₁₂P₃).



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

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

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

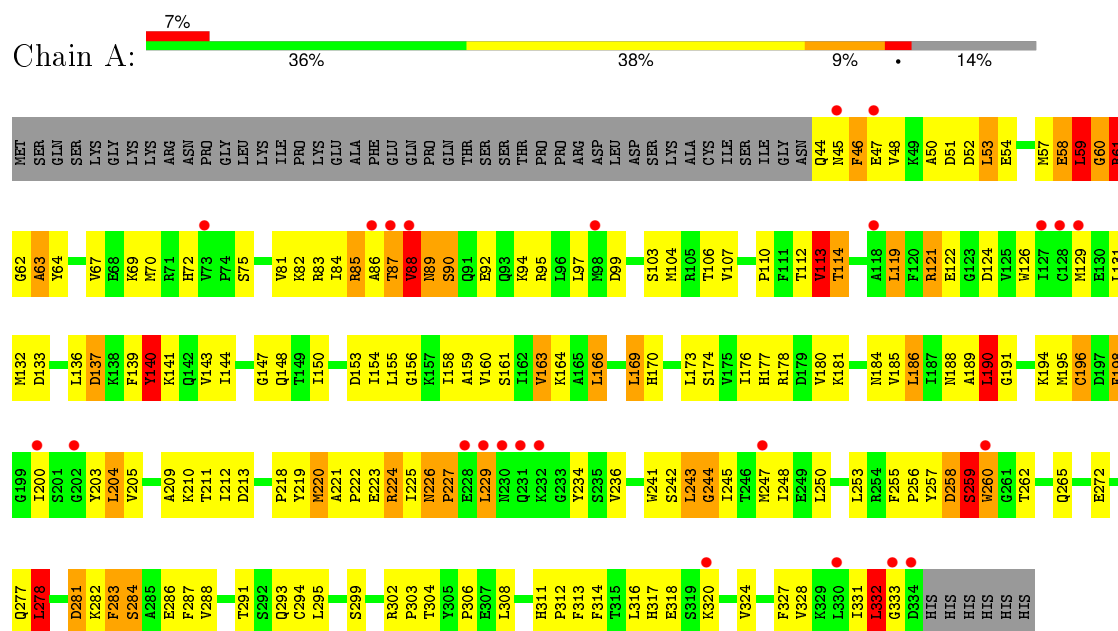
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	18	Total	O	0	0
			18	18		

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: Dual specificity mitogen-activated protein kinase kinase 6



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	83.46 Å 83.46 Å 101.15 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.60 19.66 – 2.60	Depositor EDS
% Data completeness (in resolution range)	93.6 (20.00-2.60) 93.9 (19.66-2.60)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 2.59 Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.264 , 0.280 0.283 , 0.295	Depositor DCC
R_{free} test set	1207 reflections (11.06%)	DCC
Wilson B-factor (Å ²)	76.2	Xtriage
Anisotropy	0.339	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 68.2	EDS
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	5 of 12124 reflections (0.041%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2322	wwPDB-VP
Average B, all atoms (Å ²)	98.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.43% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ANK

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	0.75	2/2321 (0.1%)	0.97	13/3142 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	59	LEU	C-O	-5.60	1.12	1.23
1	A	63	ALA	N-CA	-5.02	1.36	1.46

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	62	GLY	N-CA-C	-8.30	92.35	113.10
1	A	59	LEU	CA-CB-CG	8.26	134.30	115.30
1	A	57	MET	N-CA-C	-7.71	90.17	111.00
1	A	63	ALA	C-N-CA	7.57	140.62	121.70
1	A	226	ASN	C-N-CD	-6.50	106.31	120.60
1	A	58	GLU	N-CA-C	-6.13	94.44	111.00
1	A	63	ALA	CA-C-N	-6.07	103.84	117.20
1	A	60	GLY	N-CA-C	-6.00	98.11	113.10
1	A	61	ARG	N-CA-C	-5.95	94.93	111.00
1	A	332	LEU	CA-CB-CG	5.94	128.96	115.30
1	A	278	LEU	CA-CB-CG	5.92	128.92	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	59	LEU	CB-CG-CD1	5.69	120.67	111.00
1	A	190	LEU	CA-CB-CG	5.01	126.82	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	259	SER	Peptide

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	2272	0	2275	175	1
2	A	31	0	12	4	0
3	A	1	0	0	0	0
4	A	18	0	0	11	0
All	All	2322	0	2287	175	1

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 38.

All (175) 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:52:ASP:C	1:A:53:LEU:HD23	1.63	1.16
1:A:50:ALA:CB	1:A:51:ASP:HA	1.75	1.13
1:A:262:THR:HG23	1:A:265:GLN:HB2	1.26	1.10
1:A:259:SER:HB3	1:A:260:TRP:HA	1.09	1.08
1:A:53:LEU:HD23	1:A:53:LEU:N	1.50	1.08
1:A:46:PHE:HA	1:A:47:GLU:HG2	1.35	1.05
1:A:143:VAL:HG22	1:A:331:ILE:HD13	1.33	1.05
1:A:259:SER:CB	1:A:260:TRP:HA	1.90	1.00
1:A:50:ALA:HB1	1:A:51:ASP:CA	1.92	0.99
1:A:50:ALA:HB1	1:A:51:ASP:HA	1.00	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:SER:HB3	1:A:260:TRP:CA	1.93	0.97
1:A:224:ARG:O	1:A:227:PRO:HD3	1.65	0.95
1:A:53:LEU:CD2	1:A:53:LEU:N	2.30	0.94
1:A:104:MET:HA	4:A:514:HOH:O	1.67	0.94
1:A:225:ILE:C	1:A:227:PRO:HD2	1.89	0.92
1:A:159:ALA:HA	1:A:247:MET:CE	2.02	0.90
1:A:143:VAL:HG22	1:A:331:ILE:CD1	2.05	0.86
1:A:46:PHE:CA	1:A:47:GLU:HG2	2.06	0.85
1:A:304:THR:OG1	1:A:306:PRO:HD2	1.77	0.85
1:A:90:SER:HB2	1:A:92:GLU:OE1	1.75	0.85
1:A:218:PRO:HB2	1:A:220:MET:CE	2.09	0.83
1:A:159:ALA:HA	1:A:247:MET:HE3	1.60	0.81
1:A:224:ARG:O	1:A:227:PRO:CD	2.28	0.81
1:A:288:VAL:HG21	4:A:512:HOH:O	1.81	0.80
1:A:173:LEU:CD1	4:A:514:HOH:O	2.31	0.79
1:A:218:PRO:HB2	1:A:220:MET:HE3	1.64	0.79
1:A:189:ALA:C	1:A:191:GLY:H	1.85	0.78
1:A:53:LEU:HB2	1:A:70:MET:HE1	1.67	0.76
1:A:112:THR:O	1:A:113:VAL:HB	1.85	0.76
1:A:328:VAL:O	1:A:332:LEU:HD22	1.89	0.73
1:A:132:MET:HG2	1:A:186:LEU:HB3	1.69	0.72
1:A:63:ALA:HB1	4:A:504:HOH:O	1.90	0.71
1:A:225:ILE:O	1:A:227:PRO:HD2	1.89	0.71
1:A:85:ARG:O	1:A:87:THR:N	2.24	0.70
1:A:278:LEU:HD21	1:A:283:PHE:HB2	1.74	0.70
1:A:159:ALA:HA	1:A:247:MET:HE1	1.73	0.70
1:A:158:ILE:HG22	1:A:247:MET:HE1	1.72	0.70
1:A:176:ILE:HD13	1:A:234:TYR:HB3	1.73	0.69
1:A:85:ARG:C	1:A:87:THR:H	1.95	0.69
1:A:200:ILE:HG22	1:A:204:LEU:HD22	1.75	0.68
1:A:205:VAL:O	1:A:210:LYS:HE3	1.94	0.67
1:A:177:HIS:HB2	1:A:198:PHE:CE1	2.30	0.67
1:A:156:GLY:O	1:A:160:VAL:HG12	1.95	0.67
1:A:154:ILE:HG23	1:A:324:VAL:HG23	1.77	0.66
1:A:200:ILE:H	1:A:200:ILE:HD12	1.59	0.66
1:A:58:GLU:HG2	1:A:59:LEU:N	2.11	0.66
1:A:262:THR:CG2	1:A:265:GLN:HB2	2.15	0.65
1:A:58:GLU:OE1	1:A:61:ARG:NH1	2.29	0.65
1:A:248:ILE:HG22	1:A:256:PRO:HG3	1.77	0.65
1:A:184:ASN:ND2	2:A:401:ANK:O2B	2.29	0.65
1:A:64:TYR:N	4:A:504:HOH:O	2.29	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:ALA:C	1:A:191:GLY:N	2.51	0.64
1:A:52:ASP:C	1:A:53:LEU:CD2	2.55	0.64
1:A:53:LEU:HB2	1:A:70:MET:CE	2.28	0.63
1:A:60:GLY:HA3	1:A:67:VAL:CG1	2.27	0.63
1:A:67:VAL:HA	1:A:81:VAL:O	1.99	0.63
1:A:218:PRO:HB2	1:A:220:MET:HE1	1.81	0.62
1:A:133:ASP:HB2	1:A:188:ASN:HA	1.81	0.62
1:A:223:GLU:OE1	1:A:223:GLU:N	2.30	0.62
1:A:58:GLU:CG	1:A:59:LEU:N	2.63	0.62
1:A:139:PHE:O	1:A:141:LYS:N	2.33	0.62
1:A:170:HIS:NE2	1:A:236:VAL:HG21	2.15	0.62
1:A:60:GLY:HA3	1:A:67:VAL:HG12	1.83	0.61
1:A:218:PRO:CB	1:A:220:MET:HE3	2.30	0.61
1:A:186:LEU:HD22	1:A:196:CYS:HB3	1.81	0.61
1:A:318:GLU:OE1	1:A:318:GLU:N	2.34	0.61
1:A:272:GLU:HG2	4:A:517:HOH:O	2.02	0.60
1:A:137:ASP:HA	1:A:140:TYR:HD2	1.64	0.60
1:A:303:PRO:HB2	1:A:308:LEU:HD12	1.83	0.60
1:A:110:PRO:O	1:A:194:LYS:HE3	2.01	0.59
1:A:46:PHE:HA	1:A:47:GLU:CG	2.23	0.59
1:A:188:ASN:OD1	1:A:189:ALA:O	2.19	0.59
1:A:50:ALA:CB	1:A:51:ASP:CA	2.59	0.58
1:A:85:ARG:C	1:A:87:THR:N	2.58	0.57
1:A:177:HIS:CD2	1:A:198:PHE:HA	2.39	0.57
1:A:137:ASP:HA	1:A:140:TYR:CD2	2.39	0.57
1:A:218:PRO:CB	1:A:220:MET:CE	2.82	0.57
1:A:60:GLY:O	1:A:67:VAL:N	2.37	0.56
1:A:131:LEU:HD23	1:A:132:MET:N	2.21	0.56
1:A:316:LEU:O	1:A:320:LYS:HB2	2.05	0.56
1:A:90:SER:CB	1:A:92:GLU:OE1	2.52	0.56
1:A:173:LEU:HD11	4:A:514:HOH:O	2.00	0.55
1:A:176:ILE:CD1	1:A:234:TYR:HB3	2.36	0.55
1:A:159:ALA:CA	1:A:247:MET:HE1	2.37	0.55
1:A:226:ASN:N	1:A:227:PRO:HD2	2.21	0.54
1:A:112:THR:O	1:A:195:MET:O	2.24	0.54
1:A:90:SER:O	1:A:94:LYS:HB2	2.08	0.54
1:A:72:HIS:HD2	1:A:75:SER:OG	1.91	0.54
1:A:137:ASP:O	1:A:141:LYS:HE3	2.08	0.53
1:A:106:THR:O	1:A:107:VAL:C	2.46	0.53
1:A:262:THR:HG23	1:A:265:GLN:CB	2.18	0.53
1:A:245:ILE:CD1	4:A:509:HOH:O	2.57	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:ALA:CA	1:A:247:MET:CE	2.82	0.52
1:A:119:LEU:HD13	1:A:126:TRP:HB2	1.92	0.52
1:A:87:THR:O	1:A:88:VAL:HG13	2.09	0.52
1:A:281:ASP:OD1	1:A:282:LYS:HG2	2.10	0.51
1:A:169:LEU:HD13	1:A:198:PHE:CZ	2.45	0.51
1:A:243:LEU:HD12	1:A:243:LEU:C	2.30	0.51
1:A:170:HIS:O	1:A:174:SER:HA	2.11	0.51
1:A:262:THR:HG22	1:A:265:GLN:NE2	2.26	0.51
1:A:150:ILE:HG23	1:A:327:PHE:CE2	2.47	0.50
1:A:222:PRO:O	1:A:226:ASN:OD1	2.30	0.50
1:A:200:ILE:CG2	1:A:204:LEU:HD22	2.41	0.50
1:A:327:PHE:CZ	1:A:331:ILE:HD11	2.46	0.50
1:A:46:PHE:HD1	1:A:48:VAL:HA	1.75	0.50
1:A:132:MET:O	2:A:401:ANK:H2	2.12	0.50
1:A:286:GLU:HB3	1:A:312:PRO:HB2	1.94	0.49
1:A:241:TRP:CD1	1:A:241:TRP:C	2.85	0.49
1:A:200:ILE:HG22	1:A:204:LEU:CD2	2.41	0.49
1:A:119:LEU:N	1:A:119:LEU:HD12	2.27	0.49
1:A:282:LYS:HB2	1:A:283:PHE:CE2	2.49	0.48
1:A:163:VAL:HG12	1:A:164:LYS:N	2.28	0.48
1:A:177:HIS:ND1	1:A:178:ARG:O	2.46	0.48
1:A:139:PHE:CZ	1:A:331:ILE:HD12	2.49	0.48
1:A:287:PHE:HB2	1:A:313:PHE:CE1	2.49	0.48
1:A:82:LYS:NZ	2:A:401:ANK:O1G	2.45	0.47
1:A:219:TYR:O	1:A:242:SER:OG	2.30	0.47
1:A:46:PHE:N	1:A:47:GLU:HG2	2.30	0.47
1:A:163:VAL:HG11	1:A:314:PHE:CE1	2.50	0.46
1:A:259:SER:CB	1:A:260:TRP:CA	2.67	0.46
1:A:136:LEU:HD21	1:A:158:ILE:HG12	1.96	0.46
1:A:258:ASP:O	1:A:259:SER:O	2.33	0.46
1:A:278:LEU:C	1:A:278:LEU:HD23	2.36	0.46
1:A:282:LYS:HB2	1:A:283:PHE:CD2	2.51	0.46
1:A:67:VAL:HG11	2:A:401:ANK:O4'	2.16	0.45
1:A:139:PHE:C	1:A:141:LYS:N	2.70	0.45
1:A:169:LEU:HD13	1:A:198:PHE:CE2	2.52	0.45
1:A:221:ALA:HB1	1:A:222:PRO:HD2	1.98	0.45
1:A:139:PHE:C	1:A:141:LYS:H	2.20	0.44
1:A:294:CYS:O	1:A:302:ARG:HD3	2.17	0.44
1:A:89:ASN:ND2	1:A:89:ASN:H	2.15	0.44
1:A:150:ILE:HG22	1:A:154:ILE:HB	1.99	0.44
1:A:181:LYS:O	1:A:185:VAL:HG23	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:ARG:O	1:A:227:PRO:HD2	2.12	0.44
1:A:287:PHE:O	1:A:288:VAL:C	2.56	0.44
1:A:160:VAL:HG13	1:A:161:SER:H	1.83	0.44
1:A:136:LEU:HA	1:A:139:PHE:HB3	2.00	0.43
1:A:176:ILE:HD12	1:A:178:ARG:HG2	2.00	0.43
1:A:284:SER:O	1:A:287:PHE:HB3	2.19	0.43
1:A:166:LEU:HD21	1:A:180:VAL:HG21	1.99	0.43
1:A:97:LEU:HD21	1:A:203:TYR:HB3	2.00	0.43
1:A:200:ILE:N	1:A:200:ILE:HD12	2.30	0.43
1:A:121:ARG:HB3	1:A:122:GLU:H	1.60	0.43
1:A:205:VAL:O	1:A:205:VAL:HG12	2.18	0.43
1:A:212:ILE:HD12	4:A:516:HOH:O	2.19	0.43
1:A:153:ASP:HA	1:A:317:HIS:CE1	2.54	0.42
1:A:245:ILE:HD12	4:A:509:HOH:O	2.17	0.42
1:A:209:ALA:O	1:A:213:ASP:HB2	2.19	0.42
1:A:84:ILE:CG2	1:A:85:ARG:N	2.82	0.42
1:A:226:ASN:O	1:A:227:PRO:HB2	2.19	0.42
1:A:122:GLU:C	1:A:124:ASP:H	2.22	0.42
1:A:44:GLN:HG3	1:A:45:ASN:H	1.84	0.42
1:A:150:ILE:HB	4:A:515:HOH:O	2.18	0.42
1:A:243:LEU:O	1:A:244:GLY:C	2.56	0.42
1:A:311:HIS:HA	1:A:312:PRO:HD2	1.75	0.42
1:A:204:LEU:C	1:A:205:VAL:HG23	2.40	0.42
1:A:53:LEU:CB	1:A:70:MET:CE	2.96	0.41
1:A:324:VAL:O	1:A:328:VAL:HG23	2.20	0.41
1:A:220:MET:HE1	1:A:220:MET:H	1.85	0.41
1:A:58:GLU:O	1:A:59:LEU:HB3	2.20	0.41
1:A:245:ILE:HG13	1:A:295:LEU:HD11	2.03	0.41
1:A:204:LEU:HD12	1:A:204:LEU:HA	1.73	0.41
1:A:211:THR:O	1:A:212:ILE:C	2.59	0.41
1:A:242:SER:O	1:A:243:LEU:C	2.58	0.41
1:A:177:HIS:HB2	1:A:198:PHE:CD1	2.56	0.41
1:A:158:ILE:O	1:A:159:ALA:C	2.59	0.41
1:A:291:THR:HG23	1:A:295:LEU:HD23	2.03	0.41
1:A:166:LEU:HD12	1:A:166:LEU:HA	1.87	0.41
1:A:293:GLN:HB3	1:A:303:PRO:CG	2.51	0.41
1:A:144:ILE:HD11	1:A:253:LEU:O	2.21	0.41
1:A:95:ARG:O	1:A:99:ASP:HB2	2.20	0.40
1:A:69:LYS:HG3	1:A:131:LEU:CD1	2.51	0.40
1:A:255:PHE:CE2	1:A:257:TYR:HB2	2.56	0.40
1:A:170:HIS:CD2	1:A:236:VAL:CG2	3.05	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:VAL:O	1:A:114:THR:CB	2.69	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	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:GLU:OE1	1:A:236:VAL:CG2[2_445]	2.00	0.20

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	289/340 (85%)	227 (78%)	45 (16%)	17 (6%)	2 2

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	46	PHE
1	A	86	ALA
1	A	88	VAL
1	A	113	VAL
1	A	227	PRO
1	A	229	LEU
1	A	259	SER
1	A	281	ASP
1	A	147	GLY
1	A	333	GLY
1	A	59	LEU
1	A	87	THR
1	A	140	TYR
1	A	190	LEU
1	A	85	ARG

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Mol	Chain	Res	Type
1	A	114	THR
1	A	244	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	251/303 (83%)	214 (85%)	37 (15%)	4 6

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	LEU
1	A	59	LEU
1	A	61	ARG
1	A	83	ARG
1	A	88	VAL
1	A	89	ASN
1	A	90	SER
1	A	103	SER
1	A	113	VAL
1	A	119	LEU
1	A	121	ARG
1	A	129	MET
1	A	137	ASP
1	A	140	TYR
1	A	148	GLN
1	A	155	LEU
1	A	163	VAL
1	A	166	LEU
1	A	169	LEU
1	A	186	LEU
1	A	190	LEU
1	A	196	CYS
1	A	198	PHE
1	A	204	LEU

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Mol	Chain	Res	Type
1	A	220	MET
1	A	224	ARG
1	A	229	LEU
1	A	243	LEU
1	A	250	LEU
1	A	258	ASP
1	A	260	TRP
1	A	277	GLN
1	A	278	LEU
1	A	283	PHE
1	A	284	SER
1	A	299	SER
1	A	332	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	GLN
1	A	72	HIS
1	A	89	ASN
1	A	148	GLN
1	A	184	ASN
1	A	265	GLN
1	A	317	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 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	ANK	A	401	3	27,33,33	1.81	5 (18%)	30,52,52	4.12	9 (30%)

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	ANK	A	401	3	-	1/12/38/38	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	ANK	PB-O1B	2.17	1.48	1.46
2	A	401	ANK	C5-C4	3.53	1.48	1.40
2	A	401	ANK	PG-N3B	4.10	1.74	1.63
2	A	401	ANK	PG-O2G	4.30	1.51	1.46
2	A	401	ANK	PB-N3B	4.61	1.75	1.63

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	ANK	C4'-O4'-C1'	-16.80	91.26	109.72
2	A	401	ANK	N3-C2-N1	-4.76	125.25	128.89
2	A	401	ANK	O2G-PG-N3B	-3.12	107.11	111.90
2	A	401	ANK	C4-C5-N7	-2.30	107.36	109.48
2	A	401	ANK	O2A-PA-O1A	2.29	124.93	112.53
2	A	401	ANK	O4'-C1'-N9	3.68	115.81	108.10
2	A	401	ANK	O3G-PG-O1G	4.04	119.55	107.58
2	A	401	ANK	C1'-N9-C4	4.33	133.47	126.94
2	A	401	ANK	C2'-C1'-N9	10.23	129.92	114.29

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	ANK	O2G-PG-N3B-PB

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	ANK	4	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	291/340 (85%)	0.47	24 (8%) 14 10	30, 95, 134, 175	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	230	ASN	8.5
1	A	229	LEU	4.8
1	A	231	GLN	4.7
1	A	202	GLY	4.4
1	A	88	VAL	4.1
1	A	334	ASP	4.1
1	A	86	ALA	3.8
1	A	260	TRP	3.6
1	A	320	LYS	3.4
1	A	228	GLU	3.3
1	A	333	GLY	3.0
1	A	200	ILE	3.0
1	A	98	MET	2.9
1	A	45	ASN	2.8
1	A	128	CYS	2.7
1	A	232	LYS	2.5
1	A	247	MET	2.5
1	A	118	ALA	2.4
1	A	129	MET	2.2
1	A	47	GLU	2.1
1	A	73	VAL	2.1
1	A	330	LEU	2.1
1	A	127	ILE	2.0
1	A	87	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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(\AA^2)	Q<0.9
2	ANK	A	401	31/31	0.96	0.17	-0.55	74,86,98,105	0
3	MG	A	402	1/1	0.96	0.09	-	98,98,98,98	0

6.5 Other polymers [i](#)

There are no such residues in this entry.