

Full wwPDB X-ray Structure Validation Report (i)

Nov 2, 2021 - 04:51 AM EDT

PDB ID	:	2YU1
Title	:	Crystal structure of hJHDM1A complexed with a-ketoglutarate
Authors	:	Han, Z.
Deposited on	:	2007-04-05
Resolution	:	2.70 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 https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.70 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.



Motria	Whole archive	Similar resolution
wietric	$(\# { m Entries})$	$(\# { m Entries}, { m resolution} { m range}({ m \AA}))$
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain					
1	А	451	39%	40%	6%	14%		



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3255 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 JmjC domain-containing histone demethylation protein 1A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	А	386	Total 3170	C 2040	N 527	O 581	S 22	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	252	SER	LYS	engineered mutation	UNP Q9Y2K7

• Molecule 2 is FE (II) ION (three-letter code: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	Total Fe 1 1	0	0

• Molecule 3 is 2-OXOGLUTARIC ACID (three-letter code: AKG) (formula: C₅H₆O₅).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	А	1	Total 10	$\begin{array}{c} \mathrm{C} \\ \mathrm{5} \end{array}$	O 5	0	0

• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	74	Total O 74 74	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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.

Note EDS was not executed.

• Molecule 1: JmjC domain-containing histone demethylation protein 1A





4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source	
Space group	P 31 2 1	Depositor	
Cell constants	81.25Å 81.25Å 124.01Å	Depositor	
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor	
Resolution (Å)	20.00 - 2.70	Depositor	
% Data completeness	99.9 (20.00-2.70)	Depositor	
(in resolution range)		Depositor	
R_{merge}	0.06	Depositor	
R_{sym}	(Not available)	Depositor	
Refinement program	CNS 1.1	Depositor	
R, R_{free}	0.238 , 0.273	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	3255	wwPDB-VP	
Average B, all atoms $(Å^2)$	51.0	wwPDB-VP	



5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: AKG, FE2

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).

Mal	Chain	Bond	lengths	Bond angles		
IVIOI	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.43	0/3256	0.69	2/4418~(0.0%)	

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	121	MET	N-CA-C	5.38	125.54	111.00
1	А	117	LYS	N-CA-C	5.24	125.14	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	А	3170	0	3063	230	0
2	А	1	0	0	0	0
3	А	10	0	4	3	0
4	А	74	0	0	11	0
All	All	3255	0	3067	230	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 37.



A 1 -		Interatomic	Clash overlap (Å)	
Atom-1	Atom-2	distance (Å)		
1:A:122:THR:HG23	1:A:125:GLN:HE22	1.13	1.08	
1:A:281:GLY:HA2	1:A:311:ILE:HD11	1.39	1.04	
1:A:100:MET:HG3	1:A:101:CYS:H	1.18	1.03	
1:A:102:VAL:HA	1:A:151:THR:HG21	1.06	1.02	
1:A:168:TRP:HB3	1:A:332:MET:HE3	1.38	1.02	
1:A:227:GLY:HA3	1:A:292:THR:HG22	1.42	1.00	
1:A:102:VAL:HA	1:A:151:THR:CG2	1.95	0.95	
1:A:102:VAL:CA	1:A:151:THR:HG21	1.97	0.94	
1:A:174:PRO:HG2	1:A:177:LEU:HD12	1.51	0.92	
1:A:172:MET:HG2	1:A:336:VAL:HG22	1.54	0.89	
1:A:228:GLY:H	1:A:292:THR:HG21	1.37	0.88	
1:A:228:GLY:H	1:A:292:THR:CG2	1.85	0.88	
1:A:168:TRP:HB3	1:A:332:MET:CE	2.09	0.83	
1:A:100:MET:HG3	1:A:101:CYS:N	1.92	0.83	
1:A:64:VAL:HG11	1:A:512:PRO:HD3	1.61	0.82	
1:A:122:THR:HG23	1:A:125:GLN:NE2	1.94	0.82	
1:A:279:PRO:HG2	1:A:282:TRP:CD1	2.14	0.82	
1:A:231:PHE:HB2	1:A:268:ILE:HG22	1.64	0.80	
1:A:237:THR:HG22	1:A:240:ASN:H	1.46	0.79	
1:A:115:THR:HG23	1:A:116:GLN:H	1.47	0.78	
1:A:121:MET:HE1	1:A:126:TRP:HA	1.65	0.77	
1:A:102:VAL:HG11	1:A:146:LEU:HD21	1.67	0.77	
1:A:290:THR:O	1:A:292:THR:HG23	1.85	0.77	
1:A:81:SER:HB2	1:A:86:ILE:HD11	1.68	0.76	
1:A:50:ASN:HB3	4:A:763:HOH:O	1.86	0.76	
1:A:111:MET:HG3	1:A:117:LYS:H	1.51	0.76	
1:A:256:ILE:HD13	1:A:261:ARG:NH1	2.01	0.75	
1:A:91:PRO:HA	1:A:225:HIS:CE1	2.20	0.75	
1:A:98:VAL:O	1:A:102:VAL:HG22	1.86	0.74	
1:A:168:TRP:O	1:A:172:MET:HB2	1.88	0.73	
1:A:107:MET:SD	1:A:122:THR:HG22	2.28	0.73	
1:A:512:PRO:HG2	1:A:515:GLN:NE2	2.04	0.72	
1:A:205:ARG:HG3	1:A:289:PRO:O	1.89	0.72	
1:A:253:GLN:HG2	4:A:717:HOH:O	1.90	0.72	
1:A:228:GLY:O	1:A:292:THR:HG21	1.93	0.69	
1:A:109:ASP:HA	1:A:119:ILE:HG22	1.75	0.68	
1:A:113:VAL:HG11	1:A:207:CYS:HB2	1.75	0.68	
1:A:112:ASP:O	1:A:114:ASN:N	2.27	0.68	
1:A:197:GLN:HG2	4:A:746:HOH:O	1.93	0.68	
1:A:338:GLU:HG2	1:A:349:HIS:HB2	1.75	0.67	

All (230) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.



	lo uo pugo	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:279:PRO:HG2	1:A:282:TRP:NE1	2.10	0.66
1:A:288:THR:HG22	1:A:290:THR:O	1.97	0.65
1:A:144:ILE:CA	1:A:201:LEU:HD12	2.26	0.65
1:A:495:LEU:HD23	1:A:499:LEU:HD12	1.77	0.65
1:A:109:ASP:HA	1:A:119:ILE:CG2	2.27	0.65
1:A:341:VAL:O	1:A:345:THR:HB	1.97	0.64
1:A:121:MET:CE	1:A:126:TRP:HA	2.27	0.64
1:A:109:ASP:CA	1:A:119:ILE:HG22	2.28	0.63
1:A:169:VAL:HB	1:A:197:GLN:NE2	2.13	0.63
1:A:115:THR:O	1:A:116:GLN:HB3	2.00	0.61
1:A:360:SER:HB3	4:A:772:HOH:O	1.98	0.61
1:A:144:ILE:HA	1:A:201:LEU:HD12	1.81	0.61
1:A:103:GLY:O	1:A:106:ARG:HB3	2.01	0.60
1:A:87:LYS:HG2	1:A:158:GLN:CB	2.31	0.60
1:A:113:VAL:HG11	1:A:207:CYS:CB	2.32	0.60
1:A:110:VAL:H	1:A:119:ILE:HA	1.66	0.60
1:A:142:ASN:ND2	1:A:203:SER:OG	2.35	0.59
1:A:317:ARG:HA	4:A:757:HOH:O	2.02	0.59
1:A:280:SER:HB2	4:A:707:HOH:O	2.01	0.59
1:A:228:GLY:H	1:A:292:THR:HG22	1.67	0.59
1:A:203:SER:HB3	1:A:207:CYS:SG	2.42	0.59
1:A:494:ILE:O	1:A:498:GLU:HG3	2.03	0.58
1:A:37:THR:HB	1:A:255:ASP:HA	1.85	0.58
1:A:227:GLY:CA	1:A:292:THR:HG22	2.26	0.58
1:A:351:THR:O	1:A:355:GLN:HG3	2.04	0.57
1:A:124:ALA:HB3	1:A:125:GLN:OE1	2.05	0.57
1:A:95:VAL:O	1:A:98:VAL:HB	2.04	0.57
1:A:148:PHE:CD1	1:A:153:LEU:HB3	2.40	0.56
1:A:301:HIS:CD2	1:A:303:PHE:H	2.23	0.56
1:A:129:TYR:CE2	1:A:141:TYR:HB2	2.40	0.56
1:A:112:ASP:C	1:A:114:ASN:H	2.09	0.56
1:A:221:TRP:HZ3	4:A:734:HOH:O	1.88	0.55
1:A:485:GLU:CD	1:A:485:GLU:H	2.10	0.55
1:A:101:CYS:O	1:A:151:THR:HB	2.07	0.55
1:A:159:ARG:HB3	1:A:221:TRP:CH2	2.41	0.55
1:A:212:HIS:CE1	3:A:701:AKG:O2	2.59	0.55
1:A:36:ARG:HH11	1:A:36:ARG:HB2	1.71	0.55
1:A:213:VAL:HG11	1:A:311:ILE:HD13	1.86	0.55
1:A:318:THR:O	1:A:318:THR:HG22	2.06	0.54
1:A:351:THR:HA	1:A:483:GLU:HG2	1.89	0.54
1:A:226:GLN:HG2	1:A:293:LEU:HB3	1.90	0.54



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:323:LYS:HB2	1:A:324:PHE:CD1	2.42	0.54
1:A:172:MET:HG3	1:A:336:VAL:HA	1.89	0.53
1:A:81:SER:CB	1:A:86:ILE:HD11	2.38	0.53
1:A:121:MET:CE	1:A:126:TRP:CA	2.86	0.53
1:A:154:GLU:C	1:A:156:MET:H	2.11	0.53
1:A:301:HIS:HD2	1:A:303:PHE:H	1.57	0.53
1:A:478:VAL:HG21	1:A:485:GLU:HG2	1.91	0.53
1:A:168:TRP:CB	1:A:332:MET:HE3	2.26	0.52
1:A:160:PRO:HG2	1:A:163:VAL:CG2	2.38	0.52
1:A:103:GLY:O	1:A:106:ARG:CB	2.58	0.52
1:A:159:ARG:HB2	1:A:160:PRO:HD2	1.91	0.52
1:A:160:PRO:HB2	4:A:745:HOH:O	2.08	0.52
1:A:331:GLU:N	1:A:331:GLU:OE1	2.43	0.52
1:A:87:LYS:HG2	1:A:158:GLN:HB3	1.90	0.52
1:A:93:PHE:CE1	1:A:156:MET:CE	2.92	0.52
1:A:111:MET:HG3	1:A:117:LYS:N	2.23	0.52
1:A:115:THR:O	1:A:116:GLN:CB	2.58	0.52
1:A:115:THR:HG23	1:A:116:GLN:N	2.20	0.51
1:A:224:ILE:HG22	1:A:272:GLN:HA	1.93	0.51
1:A:323:LYS:HD2	1:A:324:PHE:HE1	1.75	0.51
1:A:110:VAL:HG22	1:A:143:VAL:HG22	1.91	0.51
1:A:92:ASP:OD2	1:A:92:ASP:N	2.43	0.51
1:A:111:MET:HG2	1:A:112:ASP:N	2.26	0.51
1:A:213:VAL:CG1	1:A:311:ILE:HD13	2.40	0.51
1:A:111:MET:CE	1:A:113:VAL:HA	2.41	0.50
1:A:305:ILE:N	1:A:306:PRO:HD2	2.26	0.50
1:A:109:ASP:HA	1:A:119:ILE:HB	1.93	0.50
1:A:324:PHE:CD1	1:A:324:PHE:N	2.80	0.50
1:A:109:ASP:HA	1:A:119:ILE:CB	2.42	0.50
1:A:147:GLU:OE2	1:A:149:SER:OG	2.26	0.50
1:A:36:ARG:O	1:A:36:ARG:HG3	2.12	0.50
1:A:69:ARG:NH1	1:A:508:LEU:HD13	2.27	0.50
1:A:132:THR:O	1:A:133:PRO:C	2.48	0.49
1:A:133:PRO:HG2	1:A:136:GLU:HB2	1.94	0.49
1:A:144:ILE:CB	1:A:201:LEU:HD12	2.43	0.49
1:A:354:PHE:HB3	4:A:747:HOH:O	2.13	0.49
1:A:473:LEU:CD2	1:A:473:LEU:H	2.25	0.49
1:A:93:PHE:CE1	1:A:156:MET:HE3	2.47	0.49
1:A:144:ILE:HB	1:A:201:LEU:HD12	1.94	0.49
1:A:100:MET:CG	1:A:101:CYS:N	2.69	0.49
1:A:168:TRP:C	1:A:332:MET:HE3	2.34	0.49



	A L C	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:102:VAL:HG23	1:A:103:GLY:N	2.29	0.48
1:A:462:ARG:HD3	4:A:764:HOH:O	2.13	0.48
1:A:113:VAL:CG2	1:A:140:LEU:HD13	2.42	0.48
1:A:111:MET:CG	1:A:116:GLN:HA	2.43	0.48
1:A:175:ARG:HD2	1:A:517:PRO:C	2.33	0.48
1:A:515:GLN:HG3	1:A:516:TRP:CE3	2.49	0.48
1:A:515:GLN:HG3	1:A:516:TRP:CZ3	2.48	0.48
1:A:172:MET:SD	1:A:514:VAL:HG11	2.54	0.48
1:A:314:ILE:HA	1:A:317:ARG:NH1	2.28	0.48
1:A:147:GLU:HG2	1:A:198:LYS:O	2.13	0.47
1:A:211:PHE:HD2	1:A:283:ILE:HG22	1.79	0.47
1:A:334:TRP:HB2	1:A:479:PRO:HD3	1.96	0.47
1:A:213:VAL:HG11	1:A:311:ILE:CD1	2.44	0.47
1:A:231:PHE:HB2	1:A:268:ILE:CG2	2.38	0.47
1:A:93:PHE:HE1	1:A:156:MET:CE	2.28	0.47
1:A:170:ASP:O	1:A:517:PRO:HB3	2.14	0.47
1:A:63:ASN:O	1:A:66:TYR:HB3	2.14	0.47
1:A:201:LEU:H	1:A:201:LEU:HD22	1.79	0.47
1:A:512:PRO:HG2	1:A:515:GLN:HE22	1.75	0.47
1:A:478:VAL:CG2	1:A:485:GLU:HG2	2.45	0.47
1:A:495:LEU:CD2	1:A:499:LEU:HD12	2.44	0.47
1:A:235:PRO:HB3	1:A:282:TRP:CH2	2.50	0.47
1:A:234:ILE:HB	1:A:283:ILE:HB	1.96	0.47
1:A:89:PRO:O	1:A:90:ASP:C	2.53	0.47
1:A:160:PRO:HG2	1:A:163:VAL:HG23	1.96	0.46
1:A:226:GLN:O	1:A:226:GLN:HG3	2.14	0.46
1:A:514:VAL:O	1:A:515:GLN:NE2	2.48	0.46
1:A:126:TRP:HZ3	1:A:141:TYR:HB3	1.80	0.46
1:A:42:GLU:HG2	1:A:46:THR:CG2	2.45	0.46
1:A:201:LEU:HD22	1:A:201:LEU:N	2.30	0.46
1:A:306:PRO:HD3	1:A:460:GLY:HA2	1.97	0.46
1:A:42:GLU:O	1:A:46:THR:HG23	2.16	0.45
1:A:346:ASN:O	1:A:346:ASN:ND2	2.50	0.45
1:A:148:PHE:O	1:A:150:HIS:N	2.50	0.45
1:A:231:PHE:O	1:A:267:ARG:HA	2.16	0.45
1:A:95:VAL:HG21	1:A:130:TYR:CD2	2.51	0.45
1:A:345:THR:O	1:A:345:THR:CG2	2.64	0.45
1:A:360:SER:O	1:A:364:GLU:HG3	2.17	0.45
1:A:488:LEU:O	1:A:492:VAL:HG23	2.17	0.45
1:A:172:MET:CG	1:A:336:VAL:HA	2.47	0.45
1:A:39:ASP:OD1	1:A:41:GLU:HB2	2.16	0.45



	1	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:50:ASN:ND2	1:A:50:ASN:H	2.15	0.45
1:A:350:LEU:HA	1:A:481:GLY:O	2.17	0.45
1:A:134:GLU:HG3	1:A:205:ARG:CZ	2.46	0.44
1:A:334:TRP:CZ3	1:A:485:GLU:HB3	2.52	0.44
1:A:361:MET:HA	1:A:364:GLU:OE1	2.17	0.44
1:A:99:LYS:HA	1:A:123:MET:HE2	1.99	0.44
1:A:113:VAL:HG22	1:A:140:LEU:HB3	1.99	0.44
1:A:175:ARG:HD3	4:A:741:HOH:O	2.17	0.44
1:A:110:VAL:O	1:A:117:LYS:O	2.36	0.44
1:A:228:GLY:N	1:A:292:THR:CG2	2.68	0.44
1:A:106:ARG:HG2	1:A:106:ARG:HH11	1.83	0.44
1:A:106:ARG:O	1:A:107:MET:C	2.56	0.44
1:A:112:ASP:C	1:A:114:ASN:N	2.71	0.44
1:A:117:LYS:O	1:A:119:ILE:N	2.51	0.44
1:A:361:MET:O	1:A:364:GLU:HB2	2.18	0.43
1:A:44:LEU:HD21	1:A:287:TYR:CD2	2.54	0.43
1:A:121:MET:CE	1:A:126:TRP:N	2.81	0.43
1:A:328:PHE:HA	1:A:331:GLU:OE2	2.18	0.43
1:A:252:SER:O	1:A:254:GLY:N	2.51	0.43
1:A:105:ARG:O	1:A:105:ARG:HG2	2.19	0.43
1:A:516:TRP:CE3	1:A:516:TRP:N	2.86	0.43
1:A:77:ILE:HD11	1:A:268:ILE:HG12	1.99	0.43
1:A:229:LYS:NZ	1:A:288:THR:OG1	2.49	0.43
1:A:220:VAL:HG21	1:A:222:TYR:CZ	2.54	0.43
1:A:504:PRO:O	1:A:508:LEU:HG	2.18	0.43
1:A:221:TRP:HA	1:A:276:PHE:O	2.18	0.43
1:A:270:LEU:HD11	1:A:276:PHE:CD1	2.54	0.43
1:A:237:THR:CG2	1:A:239:HIS:HB3	2.50	0.42
1:A:67:ILE:HD12	1:A:67:ILE:N	2.34	0.42
1:A:122:THR:H	1:A:125:GLN:NE2	2.17	0.42
1:A:309:LEU:HD12	1:A:463:CYS:HB3	2.01	0.42
1:A:516:TRP:O	1:A:517:PRO:C	2.56	0.42
1:A:62:PHE:O	1:A:162:THR:OG1	2.28	0.42
1:A:191:MET:O	1:A:191:MET:HG3	2.19	0.42
1:A:249:LEU:C	1:A:251:GLY:H	2.22	0.42
1:A:354:PHE:CZ	1:A:480:THR:HG23	2.54	0.42
1:A:354:PHE:HZ	1:A:480:THR:HG23	1.84	0.42
1:A:211:PHE:HA	1:A:284:HIS:O	2.20	0.42
1:A:87:LYS:HZ3	1:A:158:GLN:HB2	1.85	0.42
1:A:99:LYS:HA	1:A:123:MET:CE	2.50	0.42
1:A:112:ASP:OD2	1:A:114:ASN:HB2	2.20	0.42



Atom 1	Atom 2	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:305:ILE:HG21	1:A:464:LEU:HG	2.02	0.41
1:A:333:CYS:HB3	1:A:464:LEU:HD22	2.02	0.41
1:A:91:PRO:HA	1:A:225:HIS:NE2	2.36	0.41
1:A:110:VAL:HG21	1:A:121:MET:HG3	2.02	0.41
1:A:501:ASN:H	1:A:501:ASN:ND2	2.19	0.41
1:A:130:TYR:O	1:A:137:ARG:NH1	2.54	0.41
1:A:232:TRP:O	1:A:284:HIS:HA	2.21	0.41
1:A:467:LYS:HA	1:A:467:LYS:HD2	1.87	0.41
1:A:202:MET:HG2	1:A:293:LEU:CD2	2.51	0.41
1:A:208:TYR:CZ	1:A:210:ASP:HA	2.56	0.41
1:A:212:HIS:NE2	3:A:701:AKG:O1	2.54	0.41
1:A:93:PHE:HE1	1:A:156:MET:HE1	1.84	0.41
1:A:316:ASP:OD1	1:A:325:ARG:NH2	2.53	0.41
1:A:36:ARG:HB2	1:A:36:ARG:NH1	2.35	0.41
1:A:150:HIS:N	1:A:150:HIS:ND1	2.69	0.41
1:A:212:HIS:NE2	3:A:701:AKG:C1	2.84	0.41
1:A:245:GLU:O	1:A:249:LEU:HG	2.21	0.41
1:A:473:LEU:H	1:A:473:LEU:HD22	1.86	0.41
1:A:100:MET:HE2	1:A:101:CYS:N	2.36	0.41
1:A:129:TYR:CE1	1:A:137:ARG:HG3	2.56	0.41
1:A:172:MET:HG2	1:A:336:VAL:CG2	2.38	0.40
1:A:224:ILE:HD12	1:A:224:ILE:N	2.36	0.40
1:A:473:LEU:HD22	1:A:473:LEU:N	2.36	0.40
1:A:270:LEU:HD11	1:A:276:PHE:CG	2.55	0.40
1:A:104:SER:C	1:A:106:ARG:N	2.74	0.40
1:A:134:GLU:HB2	1:A:205:ARG:NH2	2.36	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	А	380/451~(84%)	341 (90%)	27 (7%)	12 (3%)	4 9

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	113	VAL
1	А	253	GLN
1	А	483	GLU
1	А	115	THR
1	А	116	GLN
1	А	118	GLY
1	А	149	SER
1	А	117	LYS
1	А	155	ASN
1	А	322	ASN
1	А	508	LEU
1	А	479	PRO

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	Perce	entiles
1	А	351/413~(85%)	324~(92%)	27~(8%)	13	30

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	36	ARG
1	А	40	LEU
1	А	50	ASN
1	А	72	LEU
1	А	90	ASP
1	А	92	ASP
1	А	99	LYS
1	А	140	LEU
1	А	145	SER



Mol	Chain	Res	Type
1	А	150	HIS
1	А	153	LEU
1	А	156	MET
1	А	161	SER
1	А	164	ASP
1	А	172	MET
1	А	193	TYR
1	А	201	LEU
1	А	215	PHE
1	А	221	TRP
1	А	237	THR
1	А	252	SER
1	А	258	LEU
1	А	357	GLU
1	А	488	LEU
1	А	491	ASP
1	А	501	ASN
1	А	516	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type	
1	А	50	ASN	
1	А	52	ASN	
1	A 142 AS		ASN	
1	А	158	GLN	
1	А	197	GLN	
1	А	298	ASN	
1	А	301	HIS	
1	А	455	HIS	
1	А	501	ASN	
1	А	515	GLN	

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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



5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

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	Dec	Tink	Bond lengths			Bond angles			
	Type	Cham	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2
3	AKG	А	701	2	3,9,9	1.07	0	4,11,11	0.71	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	AKG	А	701	2	-	0/3/9/9	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	А	701	AKG	3	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)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

