Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:51 PM GMT

PDB ID	:	1HLU
Title	:	STRUCTURE OF BOVINE BETA-ACTIN-PROFILIN COMPLEX WITH ACTIN BOUND ATP PHOSPHATES SOLVENT ACCESSIBLE
Authors	:	Chik, J.K.; Lindberg, U.; Schutt, C.E.
Deposited on	:	1997-05-30
Resolution	:	2.65 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:

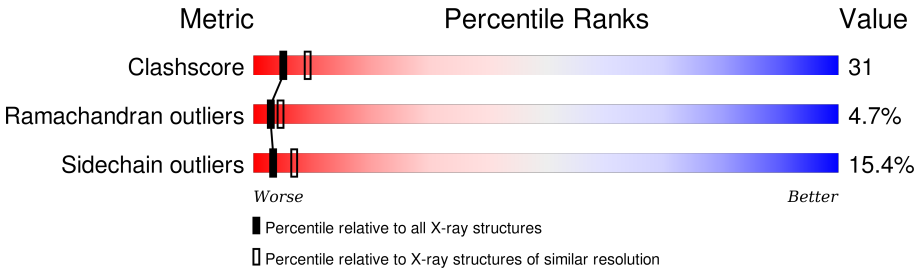
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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))
Clashscore	102246	3524 (2.70-2.62)
Ramachandran outliers	100387	3469 (2.70-2.62)
Sidechain outliers	100360	3469 (2.70-2.62)

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 <=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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	375	41% 48% 10%
2	P	140	56% 35% 9%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4001 atoms, of which 1 is hydrogen 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 BETA-ACTIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	375	Total	C	N	O	S	0	0	0
			2921	1848	490	561	22			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	73	HIC	HIS	MODIFIED RESIDUE	UNP P60712

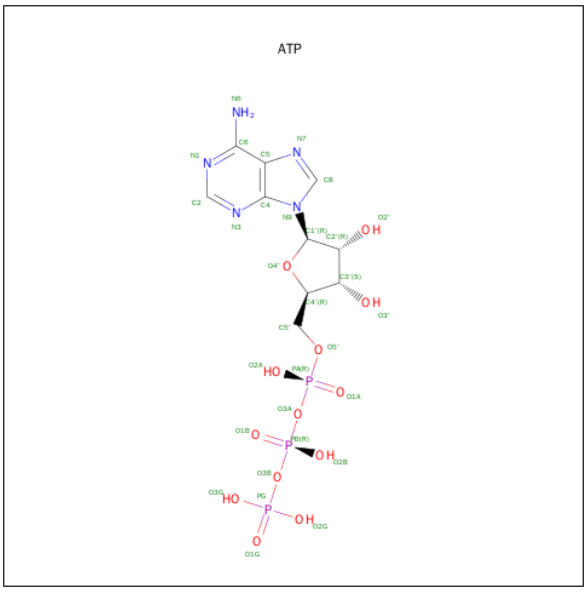
- Molecule 2 is a protein called PROFILIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	140	Total	C	N	O	S	0	0	0
			1047	659	179	200	9			

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O	P		
4	A	1	32	10	1	5	13	3	0	0

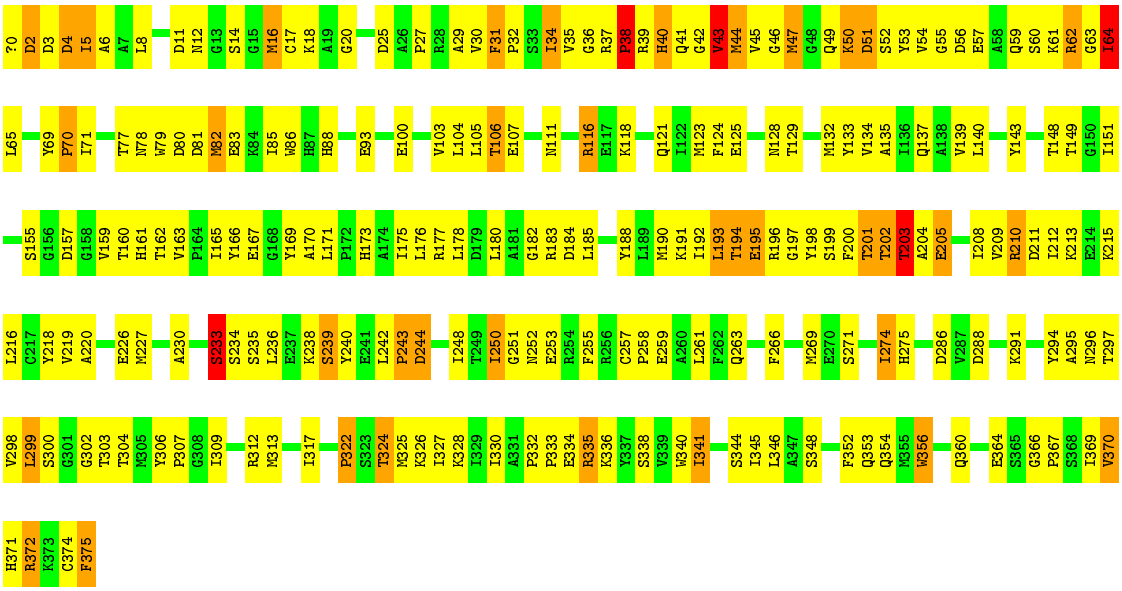
3 Residue-property plots (i)

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.

Note EDS was not executed.

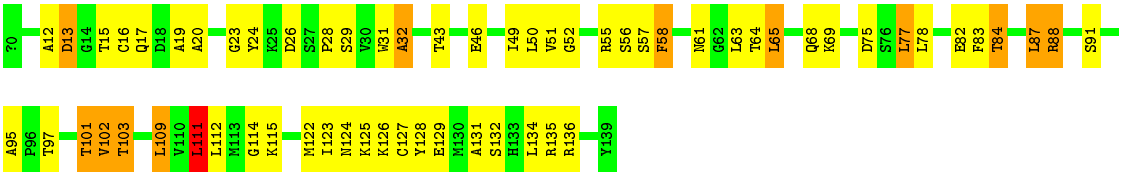
- Molecule 1: BETA-ACTIN

Chain A:



- Molecule 2: PROFILIN

Chain P:



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	38.14Å 72.24Å 185.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.65	Depositor
% Data completeness (in resolution range)	87.0 (8.00-2.65)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.201 , 0.330	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4001	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, HIC, ATP, ACE

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	0.61	0/2969	0.95	7/4019 (0.2%)
2	P	0.66	0/1063	0.88	1/1435 (0.1%)
All	All	0.63	0/4032	0.93	8/5454 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	234	SER	N-CA-C	9.30	136.10	111.00
1	A	233	SER	N-CA-C	8.76	134.66	111.00
2	P	111	LEU	CA-CB-CG	6.54	130.33	115.30
1	A	44	MET	N-CA-C	6.50	128.55	111.00
1	A	375	PHE	N-CA-C	6.50	128.54	111.00
1	A	43	VAL	CA-C-N	-5.38	105.36	117.20
1	A	43	VAL	C-N-CA	5.18	134.65	121.70
1	A	116	ARG	NE-CZ-NH1	-5.03	117.78	120.30

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	2921	0	2884	201	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	P	1047	0	1050	46	0
3	A	1	0	0	0	0
4	A	31	1	12	3	0
All	All	4000	1	3946	245	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 31.

All (245) 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:216:LEU:HD23	1:A:250:ILE:HD11	1.48	0.96
1:A:257:CYS:SG	1:A:258:PRO:HD3	2.06	0.95
2:P:16:CYS:SG	2:P:112:LEU:HD11	2.07	0.94
2:P:101:THR:HG22	2:P:123:ILE:HG22	1.50	0.93
1:A:302:GLY:HA3	4:A:1:ATP:H5'2	1.51	0.91
1:A:16:MET:HE2	1:A:32:PRO:HA	1.57	0.85
1:A:5:ILE:H	1:A:5:ILE:HD12	1.40	0.85
1:A:70:PRO:HG2	1:A:85:ILE:HD11	1.60	0.83
1:A:192:ILE:O	1:A:195:GLU:HG2	1.78	0.82
1:A:185:LEU:HB2	1:A:213:LYS:NZ	1.96	0.80
1:A:185:LEU:HD21	1:A:261:LEU:HD21	1.64	0.78
2:P:102:VAL:HB	2:P:111:LEU:HD13	1.63	0.78
2:P:91:SER:HB2	2:P:95:ALA:HB3	1.66	0.77
1:A:36:GLY:HA3	1:A:65:LEU:HD23	1.65	0.77
1:A:216:LEU:CD2	1:A:250:ILE:HD11	2.14	0.77
2:P:101:THR:CG2	2:P:123:ILE:HG22	2.14	0.77
1:A:180:LEU:HD11	1:A:269:MET:SD	2.25	0.77
2:P:64:THR:HA	2:P:68:GLN:O	1.88	0.74
1:A:35:VAL:HG21	1:A:81:ASP:HB2	1.69	0.74
1:A:160:THR:HG23	1:A:180:LEU:O	1.89	0.72
1:A:31:PHE:HE2	1:A:88:HIS:CD2	2.08	0.72
1:A:53:TYR:HB2	1:A:65:LEU:HD21	1.70	0.72
2:P:114:GLY:HA3	2:P:123:ILE:HD11	1.72	0.71
1:A:157:ASP:OD1	1:A:183:ARG:HG3	1.90	0.71
1:A:170:ALA:H	1:A:375:PHE:HE2	1.40	0.69
1:A:16:MET:HG2	1:A:30:VAL:HG12	1.75	0.69
1:A:71:ILE:HD11	1:A:82:MET:CE	2.24	0.68
1:A:107:GLU:O	1:A:137:GLN:HG3	1.94	0.67
1:A:70:PRO:O	1:A:77:THR:HB	1.95	0.67
1:A:159:VAL:HG23	1:A:178:LEU:O	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:353:GLN:NE2	1:A:356:TRP:HZ2	1.93	0.66
1:A:334:GLU:O	1:A:334:GLU:HG3	1.96	0.66
1:A:257:CYS:SG	1:A:258:PRO:CD	2.83	0.65
1:A:341:ILE:O	1:A:345:ILE:HG13	1.97	0.64
1:A:353:GLN:NE2	1:A:356:TRP:CZ2	2.65	0.64
1:A:374:CYS:O	1:A:375:PHE:HB2	1.98	0.63
1:A:372:ARG:HH21	2:P:84:THR:HG21	1.62	0.63
1:A:360:GLN:O	1:A:364:GLU:HG2	1.98	0.63
1:A:328:LYS:HG2	1:A:328:LYS:O	1.98	0.63
1:A:295:ALA:HB2	1:A:326:LYS:HG3	1.80	0.63
2:P:91:SER:HB3	2:P:97:THR:HG23	1.80	0.63
1:A:49:GLN:HG3	1:A:50:LYS:HE2	1.80	0.63
1:A:70:PRO:HG2	1:A:85:ILE:CD1	2.28	0.62
1:A:31:PHE:CE2	1:A:88:HIS:CD2	2.86	0.62
1:A:31:PHE:HE2	1:A:88:HIS:HD2	1.47	0.62
1:A:38:PRO:HB3	1:A:47:MET:SD	2.39	0.62
1:A:242:LEU:CD1	1:A:243:PRO:HD2	2.30	0.61
1:A:166:TYR:CD1	1:A:167:GLU:HG2	2.35	0.61
1:A:31:PHE:CE2	1:A:88:HIS:HD2	2.19	0.61
1:A:31:PHE:HD1	1:A:31:PHE:H	1.48	0.61
2:P:77:LEU:HD21	2:P:102:VAL:HG22	1.82	0.61
1:A:240:TYR:HB3	1:A:248:ILE:HG12	1.82	0.61
2:P:63:LEU:HD11	2:P:87:LEU:CD2	2.30	0.61
1:A:78:ASN:O	1:A:82:MET:HB2	2.01	0.61
1:A:291:LYS:HG3	1:A:325:MET:HB3	1.82	0.61
2:P:82:GLU:O	2:P:83:PHE:HB2	2.00	0.60
1:A:139:VAL:HA	1:A:165:ILE:CD1	2.32	0.60
1:A:163:VAL:HG13	1:A:175:ILE:CD1	2.31	0.60
1:A:71:ILE:HD11	1:A:82:MET:HE3	1.83	0.60
1:A:208:ILE:O	1:A:212:ILE:HG13	2.02	0.59
1:A:16:MET:CG	1:A:30:VAL:HG12	2.32	0.59
1:A:236:LEU:O	1:A:236:LEU:HD23	2.02	0.59
2:P:125:LYS:O	2:P:129:GLU:HG2	2.04	0.58
1:A:275:HIS:HA	1:A:313:MET:HE1	1.86	0.58
1:A:193:LEU:HD11	1:A:250:ILE:CG2	2.33	0.58
1:A:185:LEU:HB2	1:A:213:LYS:HZ1	1.67	0.58
1:A:198:TYR:CE2	1:A:248:ILE:HB	2.38	0.58
2:P:101:THR:HG22	2:P:123:ILE:CG2	2.30	0.58
1:A:64:ILE:HG23	1:A:65:LEU:HD13	1.85	0.58
1:A:180:LEU:CD1	1:A:269:MET:SD	2.91	0.58
1:A:139:VAL:HG12	1:A:143:TYR:CE2	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:PRO:O	1:A:325:MET:SD	2.62	0.57
1:A:166:TYR:CE1	1:A:167:GLU:HG2	2.39	0.57
1:A:193:LEU:HD11	1:A:250:ILE:HG22	1.86	0.57
1:A:161:HIS:NE2	1:A:177:ARG:HG3	2.19	0.57
1:A:140:LEU:HA	1:A:143:TYR:HD2	1.69	0.57
1:A:242:LEU:HD12	1:A:243:PRO:HD2	1.86	0.57
1:A:367:PRO:O	1:A:370:VAL:HG23	2.05	0.57
2:P:88:ARG:NH1	2:P:97:THR:OG1	2.37	0.56
1:A:41:GLN:CD	1:A:42:GLY:H	2.08	0.56
1:A:196:ARG:O	1:A:196:ARG:HG3	2.06	0.56
1:A:193:LEU:HD12	1:A:253:GLU:HG2	1.88	0.56
2:P:103:THR:CG2	2:P:127:CYS:SG	2.94	0.56
1:A:198:TYR:HB3	1:A:200:PHE:CE1	2.42	0.55
1:A:139:VAL:HA	1:A:165:ILE:HD13	1.89	0.55
1:A:61:LYS:CB	1:A:65:LEU:HD22	2.37	0.55
1:A:44:MET:HG3	1:A:46:GLY:H	1.72	0.55
1:A:53:TYR:HD1	1:A:57:GLU:OE1	1.89	0.55
1:A:61:LYS:O	1:A:63:GLY:N	2.40	0.55
1:A:54:VAL:HG22	1:A:55:GLY:N	2.22	0.54
2:P:23:GLY:O	2:P:28:PRO:HA	2.07	0.54
1:A:159:VAL:HG21	1:A:177:ARG:NH1	2.23	0.54
1:A:34:ILE:O	1:A:54:VAL:HG23	2.07	0.54
1:A:218:TYR:CE1	1:A:255:PHE:HB3	2.42	0.54
1:A:191:LYS:O	1:A:194:THR:HG22	2.07	0.54
1:A:296:ASN:HA	1:A:330:ILE:CD1	2.37	0.54
1:A:263:GLN:O	1:A:266:PHE:HB2	2.07	0.53
1:A:185:LEU:CB	1:A:213:LYS:NZ	2.71	0.53
1:A:352:PHE:CE2	1:A:356:TRP:CZ3	2.97	0.53
1:A:71:ILE:HD11	1:A:82:MET:HE2	1.89	0.53
1:A:178:LEU:HD11	1:A:271:SER:OG	2.08	0.53
1:A:275:HIS:HA	1:A:313:MET:CE	2.38	0.53
1:A:169:TYR:HA	1:A:375:PHE:CZ	2.44	0.53
1:A:70:PRO:CG	1:A:85:ILE:HD11	2.36	0.52
1:A:218:TYR:HA	1:A:307:PRO:HD2	1.91	0.52
1:A:220:ALA:HB1	1:A:226:GLU:HG3	1.90	0.52
1:A:212:ILE:HG12	1:A:240:TYR:CE2	2.45	0.52
1:A:190:MET:HB2	1:A:209:VAL:HG21	1.92	0.52
1:A:16:MET:CE	1:A:32:PRO:HA	2.33	0.52
1:A:121:GLN:O	1:A:125:GLU:HG3	2.10	0.52
1:A:201:THR:HG23	1:A:202:THR:H	1.75	0.52
1:A:219:VAL:HG23	1:A:306:TYR:HB3	1.90	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:63:LEU:HD12	2:P:63:LEU:O	2.10	0.51
1:A:151:ILE:HG23	1:A:297:THR:HA	1.93	0.51
1:A:133:TYR:HE2	1:A:135:ALA:HB2	1.74	0.51
2:P:132:SER:O	2:P:136:ARG:HG3	2.10	0.51
1:A:239:SER:HA	1:A:248:ILE:O	2.10	0.51
1:A:0:ACE:H3	1:A:6:ALA:CB	2.41	0.51
2:P:51:VAL:HG12	2:P:51:VAL:O	2.11	0.51
1:A:243:PRO:HG2	1:A:244:ASP:H	1.74	0.51
1:A:132:MET:O	1:A:356:TRP:HB2	2.11	0.50
1:A:0:ACE:H3	1:A:6:ALA:HB2	1.92	0.50
1:A:132:MET:CE	1:A:134:VAL:HG23	2.42	0.50
1:A:171:LEU:HB3	1:A:173:HIS:CE1	2.47	0.50
1:A:161:HIS:CD2	1:A:177:ARG:HG3	2.46	0.50
1:A:163:VAL:HG13	1:A:175:ILE:HD11	1.93	0.50
1:A:159:VAL:HG21	1:A:177:ARG:HH11	1.77	0.50
1:A:309:ILE:O	1:A:313:MET:HB2	2.12	0.50
1:A:116:ARG:HD3	1:A:371:HIS:CE1	2.47	0.49
1:A:159:VAL:HG22	1:A:160:THR:N	2.27	0.49
1:A:216:LEU:HD23	1:A:250:ILE:CD1	2.34	0.49
1:A:356:TRP:O	1:A:356:TRP:HD1	1.95	0.49
1:A:295:ALA:O	1:A:328:LYS:HE2	2.13	0.49
1:A:79:TRP:O	1:A:83:GLU:HB2	2.11	0.49
1:A:61:LYS:HB2	1:A:65:LEU:HD22	1.93	0.49
1:A:159:VAL:HG21	1:A:177:ARG:HG2	1.93	0.49
2:P:103:THR:HG22	2:P:127:CYS:SG	2.52	0.49
1:A:162:THR:OG1	1:A:176:LEU:HB2	2.13	0.49
2:P:91:SER:HB2	2:P:95:ALA:CB	2.39	0.49
1:A:39:ARG:HG2	1:A:64:ILE:O	2.13	0.49
1:A:325:MET:SD	1:A:325:MET:N	2.86	0.49
1:A:210:ARG:O	1:A:213:LYS:HB3	2.13	0.48
1:A:3:ASP:O	1:A:4:ASP:HB2	2.13	0.48
2:P:24:TYR:CD1	2:P:109:LEU:HD13	2.49	0.48
1:A:31:PHE:N	1:A:31:PHE:CD1	2.81	0.48
1:A:211:ASP:O	1:A:215:LYS:HB2	2.13	0.48
1:A:352:PHE:CE2	1:A:356:TRP:CE3	3.01	0.48
1:A:16:MET:HG2	1:A:30:VAL:CG1	2.43	0.48
1:A:31:PHE:HD1	1:A:31:PHE:N	2.11	0.48
1:A:159:VAL:HG21	1:A:177:ARG:CG	2.43	0.47
1:A:88:HIS:CD2	1:A:93:GLU:HG2	2.49	0.47
1:A:341:ILE:HA	1:A:341:ILE:HD13	1.70	0.47
2:P:82:GLU:O	2:P:83:PHE:CB	2.63	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:ASN:HA	1:A:330:ILE:HD11	1.97	0.47
1:A:188:TYR:HD2	1:A:257:CYS:HA	1.79	0.47
1:A:61:LYS:O	1:A:64:ILE:HG22	2.15	0.47
1:A:212:ILE:HG12	1:A:240:TYR:CD2	2.50	0.47
2:P:31:TRP:O	2:P:32:ALA:HB2	2.15	0.47
1:A:41:GLN:CG	1:A:42:GLY:H	2.27	0.47
1:A:274:ILE:HG22	1:A:275:HIS:N	2.30	0.47
1:A:69:TYR:HA	1:A:70:PRO:HD3	1.58	0.46
2:P:63:LEU:HD11	2:P:87:LEU:HD22	1.97	0.46
1:A:12:ASN:H	1:A:106:THR:HG22	1.80	0.46
1:A:60:SER:O	1:A:61:LYS:C	2.53	0.46
1:A:300:SER:HA	1:A:335:ARG:HB2	1.97	0.46
1:A:124:PHE:O	1:A:128:ASN:HA	2.16	0.46
1:A:159:VAL:CG2	1:A:160:THR:N	2.79	0.46
2:P:131:ALA:O	2:P:135:ARG:HB2	2.16	0.46
1:A:41:GLN:CG	1:A:42:GLY:N	2.78	0.46
1:A:356:TRP:CD1	1:A:356:TRP:C	2.88	0.46
1:A:372:ARG:NH2	2:P:84:THR:HG21	2.31	0.46
1:A:37:ARG:HD2	1:A:51:ASP:O	2.17	0.46
1:A:302:GLY:CA	4:A:1:ATP:H5'2	2.36	0.45
1:A:200:PHE:HA	1:A:205:GLU:HG2	1.98	0.45
2:P:101:THR:HG21	2:P:124:ASN:HA	1.97	0.45
1:A:42:GLY:O	1:A:43:VAL:HG23	2.16	0.45
1:A:294:TYR:O	1:A:327:ILE:HA	2.16	0.45
1:A:93:GLU:OE1	1:A:93:GLU:HA	2.17	0.45
1:A:159:VAL:CG2	1:A:177:ARG:HG2	2.47	0.45
1:A:116:ARG:HG2	1:A:116:ARG:NH1	2.31	0.45
1:A:35:VAL:CG2	1:A:81:ASP:HB2	2.44	0.44
2:P:77:LEU:HD21	2:P:102:VAL:CG2	2.46	0.44
1:A:334:GLU:O	1:A:334:GLU:CG	2.64	0.44
2:P:17:GLN:HB3	2:P:115:LYS:HD3	1.98	0.44
1:A:366:GLY:O	1:A:369:ILE:HG22	2.17	0.44
1:A:62:ARG:HB3	1:A:62:ARG:CZ	2.48	0.44
1:A:353:GLN:O	1:A:353:GLN:OE1	2.36	0.44
1:A:313:MET:HE2	1:A:317:ILE:HD11	2.00	0.44
1:A:298:VAL:HG12	1:A:299:LEU:N	2.32	0.44
2:P:50:LEU:HD23	2:P:63:LEU:HD22	1.99	0.44
2:P:82:GLU:O	2:P:128:TYR:HE1	2.01	0.44
1:A:133:TYR:CE2	1:A:135:ALA:HB2	2.51	0.44
1:A:259:GLU:CD	1:A:312:ARG:HH22	2.21	0.44
1:A:303:THR:O	1:A:303:THR:HG22	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:65:LEU:HD12	2:P:65:LEU:HA	1.73	0.44
1:A:100:GLU:HG2	1:A:100:GLU:O	2.18	0.44
1:A:180:LEU:HA	1:A:184:ASP:OD2	2.17	0.43
1:A:274:ILE:HA	1:A:274:ILE:HD12	1.77	0.43
1:A:160:THR:OG1	1:A:178:LEU:HB3	2.17	0.43
1:A:352:PHE:CE2	1:A:356:TRP:HZ3	2.36	0.43
1:A:203:THR:O	1:A:204:ALA:HB3	2.17	0.43
1:A:155:SER:OG	1:A:160:THR:HG22	2.17	0.43
1:A:227:MET:O	1:A:230:ALA:HB3	2.18	0.43
1:A:105:LEU:HD23	1:A:105:LEU:HA	1.72	0.43
1:A:4:ASP:OD1	1:A:5:ILE:HD12	2.19	0.43
1:A:50:LYS:C	1:A:52:SER:H	2.22	0.43
1:A:79:TRP:HZ3	1:A:86:TRP:CZ3	2.36	0.43
1:A:16:MET:HE2	1:A:16:MET:HA	2.01	0.43
1:A:334:GLU:O	1:A:336:LYS:N	2.51	0.43
1:A:8:LEU:HB2	1:A:103:VAL:HG23	2.00	0.43
1:A:304:THR:O	1:A:309:ILE:HG21	2.19	0.43
1:A:116:ARG:HD3	1:A:371:HIS:HE1	1.84	0.43
2:P:55:ARG:HD3	2:P:78:LEU:HD11	2.01	0.43
1:A:14:SER:HB2	4:A:1:ATP:PG	2.59	0.42
1:A:332:PRO:HA	1:A:333:PRO:HD2	1.79	0.42
2:P:55:ARG:HD3	2:P:78:LEU:CD1	2.49	0.42
1:A:18:LYS:HG2	1:A:27:PRO:HG3	2.00	0.42
2:P:91:SER:CB	2:P:95:ALA:HB3	2.41	0.42
1:A:17:CYS:SG	1:A:31:PHE:CE1	3.12	0.42
1:A:118:LYS:HA	1:A:118:LYS:HD2	1.86	0.42
2:P:19:ALA:O	2:P:20:ALA:HB2	2.19	0.42
1:A:53:TYR:HD1	1:A:57:GLU:CD	2.23	0.41
1:A:11:ASP:HA	1:A:106:THR:HB	2.01	0.41
2:P:69:LYS:HE2	2:P:69:LYS:HB2	1.88	0.41
1:A:240:TYR:HB3	1:A:248:ILE:CG1	2.50	0.41
1:A:105:LEU:HD11	1:A:123:MET:CG	2.50	0.41
1:A:20:GLY:HA3	1:A:340:TRP:HZ2	1.84	0.41
1:A:54:VAL:CG2	1:A:55:GLY:N	2.84	0.41
1:A:62:ARG:CB	1:A:62:ARG:NH1	2.84	0.41
1:A:39:ARG:HG3	1:A:40:HIS:N	2.35	0.41
1:A:61:LYS:HB3	1:A:65:LEU:HD13	2.03	0.41
1:A:198:TYR:HB3	1:A:200:PHE:HE1	1.84	0.41
1:A:340:TRP:CZ3	1:A:344:SER:HB2	2.55	0.41
2:P:114:GLY:HA3	2:P:123:ILE:CD1	2.48	0.41
2:P:43:THR:OG1	2:P:46:GLU:HG3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:PHE:CD2	1:A:356:TRP:CZ3	3.09	0.40
1:A:20:GLY:HA3	1:A:340:TRP:CZ2	2.56	0.40
2:P:13:ASP:HB3	2:P:15:THR:OG1	2.20	0.40
1:A:193:LEU:HD12	1:A:193:LEU:HA	1.87	0.40
2:P:112:LEU:HD21	2:P:126:LYS:HE3	2.04	0.40
2:P:50:LEU:CD2	2:P:63:LEU:HD22	2.51	0.40
2:P:84:THR:HG21	2:P:124:ASN:HD21	1.85	0.40
1:A:195:GLU:C	1:A:197:GLY:H	2.24	0.40
2:P:58:PHE:CZ	2:P:75:ASP:OD2	2.74	0.40
1:A:286:ASP:OD1	1:A:288:ASP:HB2	2.22	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	372/375 (99%)	311 (84%)	41 (11%)	20 (5%)	2 3
2	P	138/140 (99%)	120 (87%)	14 (10%)	4 (3%)	6 12
All	All	510/515 (99%)	431 (84%)	55 (11%)	24 (5%)	3 4

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4	ASP
1	A	29	ALA
1	A	62	ARG
1	A	64	ILE
1	A	203	THR
1	A	335	ARG
1	A	2	ASP
1	A	5	ILE

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Mol	Chain	Res	Type
1	A	38	PRO
1	A	40	HIS
1	A	50	LYS
1	A	182	GLY
1	A	199	SER
1	A	201	THR
1	A	233	SER
2	P	52	GLY
1	A	43	VAL
1	A	251	GLY
1	A	324	THR
2	P	26	ASP
2	P	32	ALA
1	A	202	THR
1	A	243	PRO
2	P	12	ALA

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	316/316 (100%)	269 (85%)	47 (15%)	4 8
2	P	112/112 (100%)	93 (83%)	19 (17%)	2 5
All	All	428/428 (100%)	362 (85%)	66 (15%)	3 7

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ASP
1	A	16	MET
1	A	25	ASP
1	A	31	PHE
1	A	34	ILE
1	A	38	PRO
1	A	43	VAL

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Mol	Chain	Res	Type
1	A	45	VAL
1	A	47	MET
1	A	51	ASP
1	A	56	ASP
1	A	59	GLN
1	A	64	ILE
1	A	70	PRO
1	A	80	ASP
1	A	82	MET
1	A	104	LEU
1	A	106	THR
1	A	111	ASN
1	A	129	THR
1	A	148	THR
1	A	149	THR
1	A	193	LEU
1	A	194	THR
1	A	195	GLU
1	A	203	THR
1	A	205	GLU
1	A	210	ARG
1	A	233	SER
1	A	235	SER
1	A	238	LYS
1	A	239	SER
1	A	244	ASP
1	A	250	ILE
1	A	252	ASN
1	A	274	ILE
1	A	299	LEU
1	A	322	PRO
1	A	324	THR
1	A	338	SER
1	A	341	ILE
1	A	346	LEU
1	A	348	SER
1	A	354	GLN
1	A	356	TRP
1	A	370	VAL
1	A	372	ARG
2	P	13	ASP
2	P	29	SER

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Mol	Chain	Res	Type
2	P	49	ILE
2	P	56	SER
2	P	57	SER
2	P	58	PHE
2	P	61	ASN
2	P	65	LEU
2	P	77	LEU
2	P	84	THR
2	P	87	LEU
2	P	88	ARG
2	P	101	THR
2	P	102	VAL
2	P	103	THR
2	P	109	LEU
2	P	111	LEU
2	P	122	MET
2	P	134	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	92	ASN
1	A	111	ASN
1	A	173	HIS
1	A	225	GLN
1	A	280	ASN
1	A	360	GLN
2	P	41	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
1	HIC	A	73	1	8,11,12	1.03	1 (12%)	5,14,16	0.83	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
1	HIC	A	73	1	-	0/4/6/8	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	73	HIC	CZ-NE2	-2.16	1.42	1.48

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no carbohydrates 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	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
4	ATP	A	1	3	24,33,33	0.82	1 (4%)	31,52,52	1.53	5 (16%)

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
4	ATP	A	1	3	-	0/18/38/38	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1	ATP	PG-O3G	-2.21	1.46	1.54

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	1	ATP	C2'-C1'-N9	-2.22	110.90	114.29
4	A	1	ATP	C4-C5-N7	2.12	111.43	109.48
4	A	1	ATP	O3G-PG-O1G	3.01	120.28	110.58
4	A	1	ATP	PB-O3B-PG	3.85	145.56	132.67
4	A	1	ATP	PA-O3A-PB	4.92	146.54	132.73

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
4	A	1	ATP	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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.